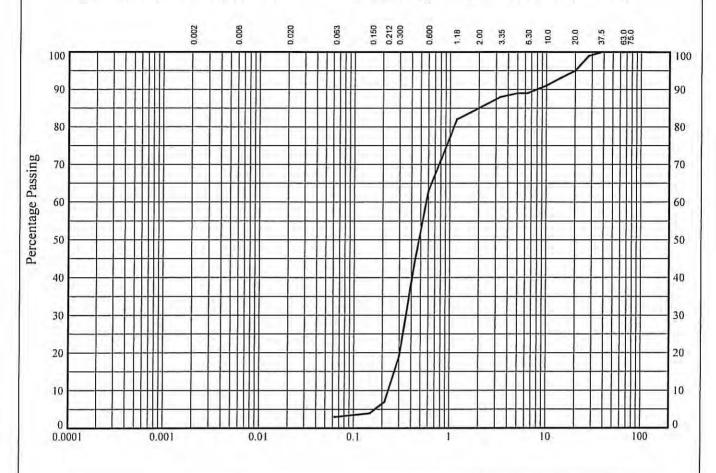
### PARTICLE SIZE DISTRIBUTION TEST

In accordance with clauses 9.2,9.5 of BS1377:Part 2:1990

Borehole: BH1 Sample Ref: Sample Type: D Depth (m): 3.50



CLAV	fine	medium	coarse	fine	medium	coarse	fine	medium	coarse	COBBLE
CLAY		SILT		SAND			GRAVEL			COBBLE

BS Test	Percentage
Sieve (mm)	Passing
125	100
90	100
75	100
63	100
50	100
37.5	100
28	99
20	95
14	93
10	91
6.3	89
5	89
3.35	88
2	85
1.18	82
0.6	63
0.425	43
0.3	19
0.212	7
0.15	4
0.063	3

Particle Diameter	Percentage Passing	Soil Fraction	Sieve Percentage	
		GRAVEL	15	
		SAND	82	
		SILT/CLAY	3	

Soil Description:

Approved Signatories: P. KENT S. CAIRNS



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PAUL KENT

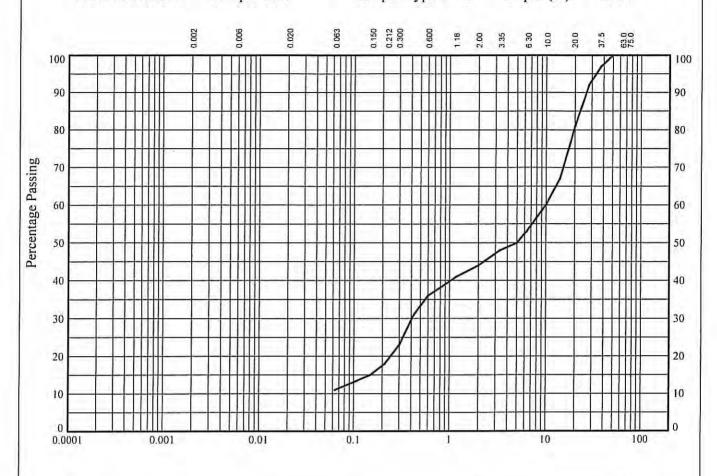
Contract
Salisbury Square, Hatfield

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### PARTICLE SIZE DISTRIBUTION TEST

In accordance with clauses 9.2,9.5 of BS1377:Part 2:1990

Borehole: BH2 Sample Ref: Sample Type: B Depth (m): 13.50



CLAV	fine	medium	coarse	fine	medium	coarse	fine	medium	coarse	COBBLES
CLAT		SILT			SAND			GRAVEI	,	CODBELL

BS Test	Percentage
Sieve (mm)	Passing
125	100
90	100
75	100
63	100
50	100
37.5	97
28	92
20	81
14	67
10	60
6.3	53
5	50
3.35	48
2	44
1.18	41
0.6	36
0.425	31
0.3	23
0.212	18
0.15	15
0.063	11

Particle	Percentage
Diameter	Passing
	1

Soil Fraction	Sieve Percentage
GRAVEL	56
SAND	33
SILT/CLAY	11

Soil Description:

Approved Signatories: P. KENT S. CAIRNS

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of 7



STRUCTURAL SOILS 18 Frogmore Road Hemel Hempstead Hertfordshire HP3 9RT

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	Date		
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Page

Salisbury Square, Hattield

## **UNCONSOLIDATED QUICK UNDRAINED (SINGLE STAGE)** TRIAXIAL COMPRESSION TEST In accordance with BS1377:Part 7:1990, Clause 8

Borehole: BH2

Sample Ref:

Mode of Failure

Sample Type:

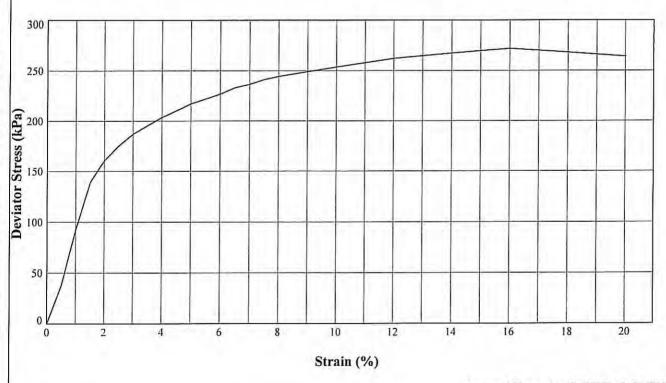
U

Depth (m): 5.70

fragments

Description: Very dark grey CLAY with occasional fine to medium chalk

STAGE NUMBER			1	2	3
SAMPLE DETAILS	Sample Condition		Undisturbed		
	Orientation of sample		Vertical		
	Diameter	(mm)	101.66		
	Height	(mm)	209.80		
	Moisture Content	(%)	19		
	Bulk Density	(Mg/m³)	2.01		
	Dry Density	(Mg/m³)	1.69		
TEST DETAILS	Membrane Thickness	(mm)	0.24		
	Rate of Axial Displacement	(%/min)	2.00		
	Cell Pressure	(kPa)	114		
	Membrane Correction	(kPa)	0.76		
	Corrected Deviator Stress	(kPa)	272		
	Undrained Shear Strength (kP		136		
	Strain at Failure	(%)	16.0		
	PACE OF VALVE OF THE STATE OF T				



Approved Signatories: P. KENT S. CAIRNS

STRUCTURAL SOILS 18 Frogmore Road Hemel Hempstead Hertfordshire HP3 9RT

Compiled By

Compound

PAUL KENT

Date 22/02/11

Contract

Salisbury Square, Hatfield

Contract Ref:

581511

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## UNCONSOLIDATED QUICK UNDRAINED (SINGLE STAGE) TRIAXIAL COMPRESSION TEST

In accordance with BS1377:Part 7:1990, Clause 8

Borehole: BH2

Sample Ref:

Sample Type:

Depth (m):

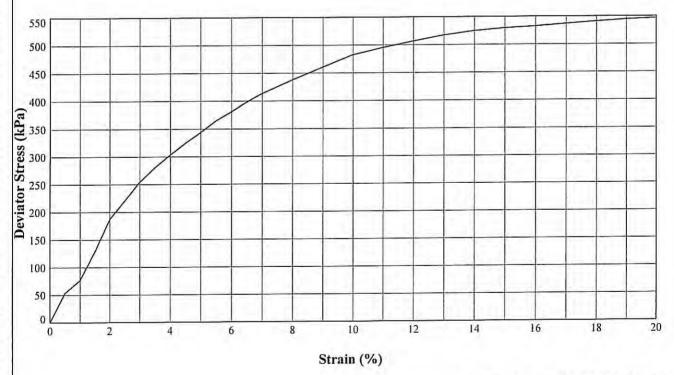
U

8.50

Description: Very dark grey CLAY with occasional fine to medium chalk

fragments

STAGE NUMBER			1	2	3
SAMPLE DETAILS	Sample Condition	Undisturbed			
	Orientation of sample		Vertical		
	Diameter	(mm)	103.13		
	Height	(mm)	210.55		
	Moisture Content	(%)	16		
	Bulk Density	(Mg/m³)	2.20		
	Dry Density	(Mg/m³)	1.91		
TEST DETAILS	Membrane Thickness	(mm)	0.24		
	Rate of Axial Displacement	(%/min)	2.00		
	Cell Pressure	(kPa)	170		
	Membrane Correction	(kPa)	0.89		
	Corrected Deviator Stress	(kPa)	547		
	Undrained Shear Strength (		273		
	Strain at Failure (%)		20.0		
	Mode of Failure		Compound		



Approved Signatories: P. KENT S. CAIRNS

STRUCTURAL SOILS 18 Frogmore Road Hemel Hempstead Hertfordshire HP3 9RT

Salisbury Square, Hatfield

Compiled By

PAUL KENT Contract Ref:

of 7

581511 Page F

AGS

Date

22/02/11

### **APPENDIX D**

**Chemical Laboratory Test Records** 





### FINAL ANALYTICAL TEST REPORT

Envirolab Job Number: 11/00569

Issue Number: 2 Date: 22 March, 2011

Client: RSK STATS Hemel Hempstead

18 Frogmore Road Hemel Hempstead

Hertfordshire

UK

HP3 9RT

Project Manager: Ben Coulston

**Project Name:** Salisbury Square, Hatfield

Project Ref: 241882

Order No: Not specified Date Samples Received: 10/02/11 Date Instructions Received: 17/03/11 Date Analysis Completed: 22/03/11

Prepared by: Approved by:

Gill Scott John Gustafson Laboratory Manager Director

Notes - Soil analysis

All results are reported as dry weight (<40℃).

Stones >10mm are removed from the sample prior to analysis and results corrected where appropriate.

Notes - General

For soil samples subscript A indicates analysis performed on the sample as received, D indicates analysis performed on dried & crushed sample.

Superscript M indicates method accredited to MCERTS.

Predominant Matrix Codes - 1 = SAND, 2 = LOAM, 3 = CLAY, 4 = LOAM/SAND, 5 = SAND/CLAY, 6 = CLAY/LOAM, 7 = OTHER. Samples with Matrix Code 7 are not predominantly a SAND/LOAM/CLAY mix and are not covered by our MCERTS accreditation. Secondary Matrix Codes - A = contains stones, B = contains construction rubble, C = contains visible hydrocarbons, D = contains glass/metal, E = contains roots/twigs.

IS indicates Insufficient sample for analysis. NDP indicates No Determination Possible. NFI indicates No Fibres Identified. Superscript # indicates method accredited to ISO 17025.

Accreditation for TPH (C6-C40) applies to the range C6-C36 only.

Analytical results reflect the quality of the sample at the time of analysis only.

Opinions and interpretations expressed are outside the scope of our accreditation.





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Lab Sample ID	11/00569/1	11/00569/2	11/00569/3	11/00569/4	11/00569/5	11/00569/6	11/00569/7	11/00569/8		
Client Sample No										
Client Sample ID	BH1	BH1	BH1	BH1	BH2	BH2	BH2	BH2		
Depth to Top	0.20	0.70	1.50	2.30	0.25	0.50	0.90	1.40		
Depth To Bottom			1.70	2.50						
Date Sampled	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11		əĘ
Sample Type	Soil - ES	Soil - ES	Soil - ES	Soil - ES	ú	Method ref				
Sample Matrix Code	7	5A	5A	1A	7	5A	7	5A	Units	Meth
ACM Screen <sub>A</sub>	-	NFI	NFI	-	NFI	NFI	-	-		Visual
pH <sub>D</sub> <sup>M#</sup>	8.1	8.8	9.4	9.0	9.0	8.6	11.6	9.0	рН	A-T-031s
Sulphate (water sol 2:1) <sub>D</sub> <sup>M#</sup>	0.02	0.05	-	0.02	-	0.03	-	-	g/l	A-T-026s
Phenois - Total by HPLC <sub>A</sub>	<0.2	-	<0.2	-	-	<0.2	-	-	mg/kg	A-T-050s
Total Organic Carbon <sub>D</sub> #	-	2.07	-	-	-	-	-	0.83	% w/w	A-T-032s
Arsenic <sub>D</sub> <sup>M#</sup>	26	23	12	7	23	-	23	11	mg/kg	A-T-024
Boron (water soluble) <sub>D</sub> <sup>M#</sup>	<1.0	<1.0	<1.0	<1.0	<1.0	-	<1.0	<1.0	mg/kg	A-T-027s
Cadmium <sub>D</sub> <sup>M#</sup>	0.6	0.6	<0.5	<0.5	0.5	-	<0.5	<0.5	mg/kg	A-T-024
Copper <sub>D</sub> <sup>M#</sup>	17	50	14	3	17	-	16	33	mg/kg	A-T-024
Chromium <sub>D</sub> <sup>M#</sup>	29	29	29	15	23	-	30	30	mg/kg	A-T-024
Lead <sub>D</sub> <sup>M#</sup>	14	278	21	5	14	-	43	46	mg/kg	A-T-024
Mercury <sub>D</sub>	<0.17	<0.17	<0.17	<0.17	<0.17	-	<0.17	<0.17	mg/kg	A-T-024
Nickel <sub>D</sub> <sup>M#</sup>	43	32	28	9	35	-	35	21	mg/kg	A-T-024
Selenium <sub>D</sub> <sup>M#</sup>	1	2	<1	<1	<1	-	<1	<1	mg/kg	A-T-024
Zinc <sub>D</sub> <sup>M#</sup>	105	177	46	17	87	-	97	62	mg/kg	A-T-024
TPH total (C6-C40) <sub>A</sub>	-	-	-	-	-	-	-	168	mg/kg	A-T-007s



Lab Sample ID 11/00569/1 1  Client Sample No  Client Sample ID BH1  Depth to Top 0.20  Depth To Bottom	11/00569/2 BH1	11/00569/3	11/00569/4	11/00569/5	11/00569/6	11/00569/7	11/00569/8		
Client Sample ID BH1  Depth to Top 0.20	BH1								
Depth to Top 0.20	BH1								
		BH1	BH1	BH2	BH2	BH2	BH2		
Denth To Bottom	0.70	1.50	2.30	0.25	0.50	0.90	1.40		
Deptil To Bottom		1.70	2.50						
Date Sampled 03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11		ef
Sample Type Soil - ES	Soil - ES	Soil - ES	Soil - ES	Soil - ES	Soil - ES	Soil - ES	Soil - ES	"	Method ref
Sample Matrix Code 7	5A	5A	1A	7	5A	7	5A	Units	Meth
TPH CWG									
Ali >C5-C6 <sub>A</sub> <0.01	-	-	-	-	-	-	-	mg/kg	A-T-022s
Ali >C6-C8 <sub>A</sub> <0.01	-	-	-	-	-	-	-	mg/kg	A-T-022s
Ali >C8-C10 <sub>A</sub> <0.01	-	-	-	-	-	-	-	mg/kg	A-T-022s
Ali >C10-C12 <sub>A</sub> # <0.1	-	-	-	-	-	-	-	mg/kg	A-T-023s
Ali >C12-C16 <sub>A</sub> # <0.1	-	-	-	-	-	-	-	mg/kg	A-T-023s
Ali >C16-C21 <sub>A</sub> # <0.1	-	-	-	-	-	-	-	mg/kg	A-T-023s
Ali >C21-C35 <sub>A</sub> # <0.1	-	-	-	-	-	-	-	mg/kg	A-T-023s
Total Aliphatics <sub>A</sub> # <0.1	-	-	-	-	-	-	-	mg/kg	A-T-022+23s
Aro >C5-C7 <sub>A</sub> <0.01	-	-	-	-	-	-	-	mg/kg	A-T-022s
Aro >C7-C8 <sub>A</sub> <0.01	-	-	-	-	-	-	-	mg/kg	A-T-022s
Aro >C8-C9 <sub>A</sub> <0.01	-	-	-	-	-	-	-	mg/kg	A-T-022s
Aro >C9-C10 <sub>A</sub> <0.01	-	-	-	-	-	-	-	mg/kg	A-T-022s
Aro >C10-C12 <sub>A</sub> # <0.1	-	-	-	-	-	-	-	mg/kg	A-T-023s
Aro >C12-C16 <sub>A</sub> # <0.1	-	-	-	-	-	-	-	mg/kg	A-T-023s
Aro >C16-C21 <sub>A</sub> # <0.1	-	-	-	-	-	-	-	mg/kg	A-T-023s
Aro >C21-C35 <sub>A</sub> # <0.1	-	-	-	-	-	-	-	mg/kg	A-T-023s
Total Aromatics <sub>A</sub> # <0.1	-	-	-	-	-	-	-	mg/kg	A-T-022+23s
TPH (Ali & Aro) <sub>A</sub> # <0.1	-	-	-	-	-	-	-	mg/kg	A-T-022+23s
MTBE <sub>A</sub> # <0.01	-	-	-	-	-	-	-	mg/kg	A-T-022s
втех									
BTEX - Benzene <sub>A</sub> # <0.01	-	-		•	-	-	-	mg/kg	A-T-022s
BTEX - Toluene <sub>A</sub> # <0.01	-	-	-	-	-	-	-	mg/kg	A-T-022s
BTEX - Ethyl Benzene <sub>A</sub> <sup>#</sup> <0.01	-	-	-	-	-	-	-	mg/kg	A-T-022s
BTEX - m & p Xylene <sub>A</sub> # <0.01	-	-		•	-	-	-	mg/kg	A-T-022s
BTEX - o Xylene <sub>A</sub> # <0.01	-	-	-	-	-	-	-	mg/kg	A-T-022s



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Lab Sample ID	11/00569/1	11/00569/2	11/00569/3	11/00569/4	11/00569/5	11/00569/6	11/00569/7	11/00569/8		
Client Sample No										
Client Sample ID	BH1	BH1	BH1	BH1	BH2	BH2	BH2	BH2		
Depth to Top	0.20	0.70	1.50	2.30	0.25	0.50	0.90	1.40		
Depth To Bottom			1.70	2.50						
Date Sampled	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11		)Ę
Sample Type	Soil - ES	Soil - ES	Soil - ES	"	Method ref					
Sample Matrix Code	7	5A	5A	1A	7	5A	7	5A	Units	Meth
PAH 16										
Acenapthene <sub>A</sub> <sup>M#</sup>	<0.01	0.01	0.02	<0.01	<0.01	-	0.04	-	mg/kg	A-T-019s
Acenapthylene <sub>A</sub> <sup>M#</sup>	<0.01	0.05	<0.01	<0.01	<0.01	-	0.14	-	mg/kg	A-T-019s
Anthracene <sub>A</sub> <sup>M#</sup>	<0.01	0.07	<0.01	<0.01	<0.01	-	0.19	-	mg/kg	A-T-019s
Benzo(a)anthracene <sub>A</sub> #	<0.01	0.36	0.01	<0.01	<0.01	-	0.67	-	mg/kg	A-T-019s
Benzo(a)pyrene <sub>A</sub> <sup>M#</sup>	0.02	0.47	<0.01	<0.01	<0.01	-	0.94	-	mg/kg	A-T-019s
Benzo(b)fluoranthene <sub>A</sub> <sup>M#</sup>	0.01	0.33	<0.01	<0.01	0.01	-	0.70	-	mg/kg	A-T-019s
Benzo(ghi)perylene <sub>A</sub> <sup>M#</sup>	<0.01	0.71	0.02	<0.01	0.01	-	1.01	-	mg/kg	A-T-019s
Benzo(k)fluoranthene <sub>A</sub>	0.02	0.48	0.02	<0.01	<0.01	-	0.76	-	mg/kg	A-T-019s
Chrysene <sub>A</sub> <sup>M#</sup>	<0.01	0.70	0.03	<0.01	0.02	-	1.48	-	mg/kg	A-T-019s
Dibenzo(ah)anthracene <sub>A</sub> #	<0.01	0.10	<0.01	<0.01	<0.01	-	0.14	-	mg/kg	A-T-019s
Fluoranthene <sub>A</sub> <sup>M#</sup>	0.01	0.78	0.07	<0.01	0.03	-	1.90	-	mg/kg	A-T-019s
Fluorene <sub>A</sub> <sup>M#</sup>	<0.01	<0.01	0.01	<0.01	<0.01	-	0.03	-	mg/kg	A-T-019s
Indeno(123-cd)pyrene <sub>A</sub> #	<0.01	0.27	<0.01	<0.01	<0.01	-	0.58	-	mg/kg	A-T-019s
Napthalene <sub>A</sub> <sup>M#</sup>	<0.01	0.03	0.11	0.02	<0.01	-	0.04	-	mg/kg	A-T-019s
Phenanthrene <sub>A</sub> <sup>M#</sup>	0.02	0.17	0.06	<0.01	0.02	-	0.65	-	mg/kg	A-T-019s
Pyrene <sub>A</sub> <sup>M#</sup>	0.01	0.75	0.06	0.02	0.03	-	1.71	-	mg/kg	A-T-019s
Total PAH <sub>A</sub> #	0.10	5.28	0.41	0.03	0.13	-	11	-	mg/kg	A-T-019s



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Lab Sample ID	11/00569/9	11/00569/10	11/00569/11	11/00569/12	11/00569/13	11/00569/14	11/00569/15	11/00569/16		
Client Sample No										
Client Sample ID	BH2	BH2	WS1	WS1	WS2	WS2	WS3	WS3		
Depth to Top	3.00	4.90	0.20	0.50	0.20	0.50	0.20	0.50		
Depth To Bottom			0.30	0.60	0.30	0.60	0.30	0.60		
Date Sampled	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11		ef
Sample Type	Soil - ES	Soil - ES	Soil - ES	Soil - ES	Soil - ES	Soil - ES	Soil - ES	Soil - ES	S	Method ref
Sample Matrix Code	5A	1A	5AE	5AE	5AE	7	5A		Units	Meth
ACM Screen <sub>A</sub>	-	-	NFI	-	NFI	NFI	NFI	NFI		Visual
pH <sub>D</sub> <sup>M#</sup>	8.7	8.9	-	8.7	-	7.6	-	-	рН	A-T-031s
Sulphate (water sol 2:1) <sub>D</sub> <sup>M#</sup>	-	0.02	-	<0.01	-	-	-	-	g/l	A-T-026s
Phenols - Total by HPLC <sub>A</sub>	<0.2	-	<0.2	-	<0.2	-	-	-	mg/kg	A-T-050s
Total Organic Carbon <sub>D</sub> #	-	-	-	-	-	0.10	-	-	% w/w	A-T-032s
Arsenic <sub>D</sub> <sup>M#</sup>	11	14	-	12	-	23	-	-	mg/kg	A-T-024
Boron (water soluble) <sub>D</sub> <sup>M#</sup>	<1.0	<1.0	-	<1.0	-	<1.0	-	-	mg/kg	A-T-027s
Cadmium <sub>D</sub> <sup>M#</sup>	<0.5	<0.5	-	<0.5	-	<0.5	-	-	mg/kg	A-T-024
Copper <sub>D</sub> <sup>M#</sup>	18	9	-	26	-	17	-	-	mg/kg	A-T-024
Chromium <sub>D</sub> <sup>M#</sup>	30	18	-	30	-	29	-	-	mg/kg	A-T-024
Lead <sub>D</sub> <sup>M#</sup>	35	16	-	68	-	17	-	-	mg/kg	A-T-024
Mercury <sub>D</sub>	<0.17	<0.17	-	<0.17	-	<0.17	-	-	mg/kg	A-T-024
Nickel <sub>D</sub> <sup>M#</sup>	21	15	-	21	-	42	-	-	mg/kg	A-T-024
Selenium <sub>D</sub> <sup>M#</sup>	<1	<1	-	<1	-	1	-	-	mg/kg	A-T-024
Zinc <sub>D</sub> <sup>M#</sup>	55	40	-	80	-	95	-	-	mg/kg	A-T-024
TPH total (C6-C40) <sub>A</sub>	-	-	-	<10	<10	ı	<10	-	mg/kg	A-T-007s



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Lab Sample ID	11/00569/9	11/00569/10	11/00569/11	11/00569/12	11/00569/13	11/00569/14	11/00569/15	11/00569/16		
Client Sample No										
Client Sample ID	BH2	BH2	WS1	WS1	WS2	WS2	WS3	WS3		
Depth to Top	3.00	4.90	0.20	0.50	0.20	0.50	0.20	0.50		
Depth To Bottom			0.30	0.60	0.30	0.60	0.30	0.60		
Date Sampled	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11		e F
Sample Type	Soil - ES	Soil - ES	Soil - ES	Soil - ES	Soil - ES	Soil - ES	Soil - ES	Soil - ES	60	Method ref
Sample Matrix Code	5A	1A	5AE	5AE	5AE	7	5A		Units	Meth
PAH 16										
Acenapthene <sub>A</sub> <sup>M#</sup>	-	0.02	-	0.02	-	<0.01	-	-	mg/kg	A-T-019s
Acenapthylene <sub>A</sub> <sup>M#</sup>	-	<0.01	-	<0.01	-	<0.01	-	-	mg/kg	A-T-019s
Anthracene <sub>A</sub> <sup>M#</sup>	-	<0.01	-	0.02	-	0.01	-	-	mg/kg	A-T-019s
Benzo(a)anthracene <sub>A</sub> #	-	<0.01	-	0.04	-	<0.01	-	-	mg/kg	A-T-019s
Benzo(a)pyrene <sub>A</sub> <sup>M#</sup>	-	<0.01	-	0.04	-	<0.01	-	-	mg/kg	A-T-019s
Benzo(b)fluoranthene <sub>A</sub> <sup>M#</sup>	-	<0.01	-	0.03	-	<0.01	-	-	mg/kg	A-T-019s
Benzo(ghi)perylene <sub>A</sub> <sup>M#</sup>	-	<0.01	-	0.09	-	<0.01	-	-	mg/kg	A-T-019s
Benzo(k)fluoranthene <sub>A</sub>	-	<0.01	-	0.03	-	<0.01	-	-	mg/kg	A-T-019s
Chrysene <sub>A</sub> <sup>M#</sup>	-	0.02	-	0.13	-	0.01	-	-	mg/kg	A-T-019s
Dibenzo(ah)anthracene <sub>A</sub> #	-	<0.01	-	<0.01	-	<0.01	-	-	mg/kg	A-T-019s
Fluoranthene <sub>A</sub> <sup>M#</sup>	-	0.03	-	0.14	-	0.03	-	-	mg/kg	A-T-019s
Fluorene <sub>A</sub> <sup>M#</sup>	-	<0.01	-	<0.01	-	<0.01	-	-	mg/kg	A-T-019s
Indeno(123-cd)pyrene <sub>A</sub> #	-	<0.01	-	0.03	-	<0.01	-	-	mg/kg	A-T-019s
Napthalene <sub>A</sub> <sup>M#</sup>	-	0.02	-	0.02	-	<0.01	-	-	mg/kg	A-T-019s
Phenanthrene <sub>A</sub> <sup>M#</sup>	-	0.02	-	0.05	-	0.02	-	-	mg/kg	A-T-019s
Pyrene <sub>A</sub> <sup>M#</sup>	-	0.03	-	0.13	-	0.02	-	-	mg/kg	A-T-019s
Total PAH <sub>A</sub> #	-	0.14	-	0.78	-	0.09	-	-	mg/kg	A-T-019s



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Lab Sample ID	11/00569/17	11/00569/18	11/00569/19	11/00569/20	11/00569/23	11/00569/25			
Client Sample No									
Client Sample ID	WS4	WS4	TP1	TP1	TP2	TP3			
Depth to Top	0.20	0.50	0.10	0.40	0.50	0.10			
Depth To Bottom	0.30	0.60							
Date Sampled	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11			əĘ
Sample Type	Soil - ES		s	Method ref					
Sample Matrix Code	7	7	4AE		4AE	4AE		Units	Meth
ACM Screen <sub>A</sub>	NFI	-	-	NFI	-	-			Visual
pH <sub>D</sub> <sup>M#</sup>	8.4	8.4	8.3	-	8.3	8.2		рН	A-T-031s
Sulphate (water sol 2:1) <sub>D</sub> <sup>M#</sup>	-	0.01	-	-	-	-		g/l	A-T-026s
Total Organic Carbon <sub>D</sub> #	0.07	-	-	-	1.08	-		% w/w	A-T-032s
Arsenic <sub>D</sub> <sup>M#</sup>	22	18	10	-	12	22		mg/kg	A-T-024
Boron (water soluble) <sub>D</sub> <sup>M#</sup>	<1.0	<1.0	<1.0	-	<1.0	<1.0		mg/kg	A-T-027s
Cadmium <sub>D</sub> <sup>M#</sup>	<0.5	<0.5	<0.5	-	<0.5	0.9		mg/kg	A-T-024
Copper <sub>D</sub> <sup>M#</sup>	14	11	22	-	37	174		mg/kg	A-T-024
Chromium <sub>D</sub> <sup>M#</sup>	21	20	18	-	20	29		mg/kg	A-T-024
Lead <sub>D</sub> <sup>M#</sup>	39	10	66	-	84	345		mg/kg	A-T-024
Mercury <sub>D</sub>	<0.17	<0.17	0.17	-	<0.17	1.03		mg/kg	A-T-024
Nickel <sub>D</sub> <sup>M#</sup>	30	33	14	-	18	33		mg/kg	A-T-024
Selenium <sub>D</sub> <sup>M#</sup>	<1	1	<1	-	1	2		mg/kg	A-T-024
Zinc <sub>D</sub> <sup>M#</sup>	73	70	80	-	112	306		mg/kg	A-T-024



Lab Sample ID	11/00569/17	11/00569/18	11/00569/19	11/00569/20	11/00569/23	11/00569/25			
Client Sample No									
Client Sample ID	WS4	WS4	TP1	TP1	TP2	TP3			
Depth to Top	0.20	0.50	0.10	0.40	0.50	0.10			
Depth To Bottom	0.30	0.60							
Date Sampled	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11			<b>5</b>
Sample Type	Soil - ES		l	92 po					
Sample Matrix Code	7	7	4AE		4AE	4AE		Units	Method ref
TPH CWG									
Ali >C5-C6 <sub>A</sub>	-	<0.01	-	-	-	-		mg/kg	A-T-022s
Ali >C6-C8 <sub>A</sub>	-	<0.01	-	-	-	-		mg/kg	A-T-022s
Ali >C8-C10 <sub>A</sub>	-	<0.01	-	-	-	-		mg/kg	A-T-022s
Ali >C10-C12 <sub>A</sub> #	-	<0.1	-	-	-	-		mg/kg	A-T-023s
Ali >C12-C16 <sub>A</sub> #	-	<0.1	-	-	-	-		mg/kg	A-T-023s
Ali >C16-C21 <sub>A</sub> #	-	<0.1	-	-	-	-		mg/kg	A-T-023s
Ali >C21-C35 <sub>A</sub> #	-	<0.1	-	-	-	-		mg/kg	A-T-023s
Total Aliphatics <sub>A</sub> #	-	<0.1	-	-	-	-		mg/kg	A-T-022+23s
Aro >C5-C7 <sub>A</sub>	-	<0.01	-	-	-	-		mg/kg	A-T-022s
Aro >C7-C8 <sub>A</sub>	-	<0.01	-	-	-	-		mg/kg	A-T-022s
Aro >C8-C9 <sub>A</sub>	-	<0.01	-	-	-	-		mg/kg	A-T-022s
Aro >C9-C10 <sub>A</sub>	-	<0.01	-	-	-	-		mg/kg	A-T-022s
Aro >C10-C12 <sub>A</sub> #	-	<0.1	-	-	-	-		mg/kg	A-T-023s
Aro >C12-C16 <sub>A</sub> #	-	<0.1	-	-	-	-		mg/kg	A-T-023s
Aro >C16-C21 <sub>A</sub> #	-	<0.1	-	-	-	-		mg/kg	A-T-023s
Aro >C21-C35 <sub>A</sub> #	-	<0.1	-	-	-	-		mg/kg	A-T-023s
Total Aromatics <sub>A</sub> #	-	<0.1	-	-	-	-		mg/kg	A-T-022+23s
TPH (Ali & Aro) <sub>A</sub> #	-	<0.1	-	-	-	-		mg/kg	A-T-022+23s
MTBE <sub>A</sub> #	-	<0.01	-	-	-	-		mg/kg	A-T-022s
ВТЕХ									
BTEX - Benzene <sub>A</sub> #	-	<0.01	-	-	-	-		mg/kg	A-T-022s
BTEX - Toluene <sub>A</sub> #	-	<0.01	-	-	-	-		mg/kg	A-T-022s
BTEX - Ethyl Benzene <sub>A</sub> #	-	<0.01	-	-	-	-		mg/kg	A-T-022s
BTEX - m & p Xylene <sub>A</sub> #	-	<0.01	-	-	-	-		mg/kg	A-T-022s
BTEX - o Xylene <sub>A</sub> #	-	<0.01	-	-	-	-		mg/kg	A-T-022s



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Lab Sample ID	11/00569/17	11/00569/18	11/00569/19	11/00569/20	11/00569/23	11/00569/25			
Client Sample No									
Client Sample ID	WS4	WS4	TP1	TP1	TP2	TP3			
Depth to Top	0.20	0.50	0.10	0.40	0.50	0.10			
Depth To Bottom	0.30	0.60							
Date Sampled	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11			e.
Sample Type	Soil - ES		"0	Method ref					
Sample Matrix Code	7	7	4AE		4AE	4AE		Units	Meth
PAH 16									
Acenapthene <sub>A</sub> <sup>M#</sup>	<0.01	<0.01	0.02	-	0.13	-		mg/kg	A-T-019s
Acenapthylene <sub>A</sub> <sup>M#</sup>	<0.01	<0.01	0.05	-	0.01	-		mg/kg	A-T-019s
Anthracene <sub>A</sub> <sup>M#</sup>	<0.01	0.01	0.09	-	1.60	-		mg/kg	A-T-019s
Benzo(a)anthracene <sub>A</sub> #	<0.01	<0.01	0.33	-	3.24	-		mg/kg	A-T-019s
Benzo(a)pyrene <sub>A</sub> <sup>M#</sup>	<0.01	<0.01	0.47	-	2.33	-		mg/kg	A-T-019s
Benzo(b)fluoranthene <sub>A</sub> M#	<0.01	<0.01	0.44	-	2.29	-		mg/kg	A-T-019s
Benzo(ghi)perylene <sub>A</sub> <sup>M#</sup>	<0.01	0.01	0.68	-	1.90	-		mg/kg	A-T-019s
Benzo(k)fluoranthene <sub>A</sub>	<0.01	<0.01	0.36	-	2.64	-		mg/kg	A-T-019s
Chrysene <sub>A</sub> <sup>M#</sup>	<0.01	0.02	0.71	-	5.27	-		mg/kg	A-T-019s
Dibenzo(ah)anthracene <sub>A</sub> #	<0.01	<0.01	0.05	-	0.42	-		mg/kg	A-T-019s
Fluoranthene <sub>A</sub> <sup>M#</sup>	<0.01	0.03	0.84	-	8.88	-		mg/kg	A-T-019s
Fluorene <sub>A</sub> <sup>M#</sup>	<0.01	<0.01	<0.01	-	0.17	-		mg/kg	A-T-019s
Indeno(123-cd)pyrene <sub>A</sub> #	<0.01	<0.01	0.26	-	1.42	-		mg/kg	A-T-019s
Napthalene <sub>A</sub> <sup>M#</sup>	<0.01	0.01	<0.01	-	0.02	-		mg/kg	A-T-019s
Phenanthrene <sub>A</sub> <sup>M#</sup>	<0.01	0.03	0.22	-	3.49	-		mg/kg	A-T-019s
Pyrene <sub>A</sub> <sup>M#</sup>	<0.01	0.02	0.79	-	6.11	-		mg/kg	A-T-019s
Total PAH <sub>A</sub> #	<0.01	0.14	5.30	-	39.9	-		mg/kg	A-T-019s



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Lab Sample ID	11/00569/17	11/00569/18	11/00569/19	11/00569/20	11/00569/23	11/00569/25			
Client Sample No									
Client Sample ID	WS4	WS4	TP1	TP1	TP2	TP3			
Depth to Top	0.20	0.50	0.10	0.40	0.50	0.10			
Depth To Bottom	0.30	0.60							
Date Sampled	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11	03-Feb-11			eĘ.
Sample Type	Soil - ES		ω.	Method ref					
Sample Matrix Code	7	7	4AE		4AE	4AE		Units	Meth
Spec PCB-WHO12									
PCB BZ 81 <sub>D</sub>	-	<0.005	-	-	-	-		mg/kg	A-T-004/5s
PCB BZ 105 <sub>D</sub>	-	<0.005	-	-	-	-		mg/kg	A-T-004/5s
PCB BZ 114 <sub>D</sub>	-	<0.005	-	-	-	-		mg/kg	A-T-004/5s
PCB BZ 118/123 <sub>D</sub>	-	<0.005	-	-	-	-		mg/kg	A-T-004/5s
PCB BZ 126 <sub>D</sub>	-	<0.005	-	-	-	-		mg/kg	A-T-004/5s
PCB BZ 156 <sub>D</sub>	-	<0.005	-	-	-	-		mg/kg	A-T-004/5s
PCB BZ 157 <sub>D</sub>	-	<0.005	-	-	-	-		mg/kg	A-T-004/5s
PCB BZ 167 <sub>D</sub>	-	<0.005	-	-	-	-		mg/kg	A-T-004/5s
PCB BZ 169 <sub>D</sub>	-	<0.005	-	-	-	-		mg/kg	A-T-004/5s
PCB BZ 189 <sub>D</sub>	-	<0.005	-	-	-	-		mg/kg	A-T-004/5s
PCB BZ 77 <sub>D</sub>	-	<0.005	-	-	-	-		mg/kg	A-T-004/5s



### FINAL ANALYTICAL TEST REPORT

**Envirolab Job Number:** 11/01190

Issue Number: 1 Date: 22 March, 2011

Client: RSK STATS Hemel Hempstead

18 Frogmore Road Hemel Hempstead

Hertfordshire

UK

HP3 9RT

Project Manager: Ben Coulston

**Project Name:** Salisbury Square, Hatfield

Project Ref: 241882

Order No: Not specified Date Samples Received: 18/03/11 Date Instructions Received: 18/03/11

Date Analysis Completed: 22/03/11

Prepared by: Approved by:

John Gustafson Gill Scott

Director Laboratory Manager

Notes - Soil analysis

All results are reported as dry weight (<40 ℃).

Stones >10mm are removed from the sample prior to analysis and results corrected where appropriate.

Notes - General

For soil samples subscript A indicates analysis performed on the sample as received, D indicates analysis performed on dried & crushed sample.

Superscript M indicates method accredited to MCERTS.

Predominant Matrix Codes - 1 = SAND, 2 = LOAM, 3 = CLAY, 4 = LOAM/SAND, 5 = SAND/CLAY, 6 = CLAY/LOAM, 7 = OTHER. Samples with Matrix Code 7 are not predominantly a SAND/LOAM/CLAY mix and are not covered by our MCERTS accreditation. Secondary Matrix Codes - A = contains stones, B = contains construction rubble, C = contains visible hydrocarbons, D = contains glass/metal, E = contains roots/twigs.

IS indicates Insufficient sample for analysis. NDP indicates No Determination Possible. NFI indicates No Fibres Identified. Superscript # indicates method accredited to ISO 17025.

Accreditation for TPH (C6-C40) applies to the range C6-C36 only.

Analytical results reflect the quality of the sample at the time of analysis only.

Opinions and interpretations expressed are outside the scope of our accreditation.







Lab Sample ID	11/01190/1	11/01190/2					
Client Sample No							
Client Sample ID	BH2	BH1					
Depth to Top	12.20	6.80					
Depth To Bottom							
Date Sampled	03-Feb-11	03-Feb-11					eŧ
Sample Type	Soil - D	Soil - D				,	Method ref
Sample Matrix Code	5A	5A				Units	Meth
pH <sub>D</sub> <sup>M#</sup>	8.2	8.0				pН	A-T-031s
Sulphate (water sol 2:1) <sub>D</sub> <sup>M#</sup>	0.15	0.28				g/I	A-T-026s

### **APPENDIX E**

**CLEA Software Output Reports** 





# Generic Assessment Criteria for Human Health Residential Scenario – Private Gardens

The human health generic assessment criteria (GAC) have been developed during a period of regulatory review and updating of the Contaminated Land Exposure Assessment (CLEA) project. Hence, the Environment Agency (EA) is in the process of publishing updated reports relating to the CLEA project and the GAC presented in this document may change to reflect these updates. This issue was prepared following the publication of soil guideline value reports and associated publications<sup>(1)</sup> for mercury, selenium, benzene, toluene, ethylbenzene and xylene in March 2009, arsenic and nickel in May 2009, cadmium and phenol in June 2009, dioxins, furans and dioxin-like PCBs in September 2009. It was also produced following publication of GAC by LQM<sup>(6)</sup>. Where available, the published soil guideline values (SGV)<sup>(1)</sup> have been used as GAC.

### 1. Model Selection

Soil assessment criteria (SAC) were calculated using CLEA v1.06 and the supporting UK guidance <sup>(1-6)</sup>. Groundwater assessment criteria (GrAC) protective of human health via the inhalation pathway were derived using the RBCA 1.3b model. RSK has updated the inputs within RBCA to reflect the UK guidance <sup>(1-5)</sup>. The SAC and GrAC collectively are termed GAC.

#### 2. Conceptual Model

In accordance with EA Science Report SC050221/SR3<sup>(3)</sup>, the residential with private garden scenario considers risks to a female child between the ages of 0 and 6 years old. In accordance with Box 3.1, SR3<sup>(3)</sup>, the pathways considered for production of the SAC in the residential with gardens scenario are:

- Direct soil and dust ingestion;
- Consumption of homegrown produce;
- Consumption of soil attached to homegrown produce;
- Dermal contact with soil and indoor dust, and
- Inhalation of indoor and outdoor dust and vapours.

Figure 1 is a conceptual model illustrating these linkages.

The pathway considered in production of the GrAC is the volatilisation of compounds from groundwater and subsequent vapour inhalation by residents whilst indoors. Figure 2 illustrates this linkage. Although the outdoor air inhalation pathway is also valid, this contributes little to the overall risks owing to the dilution in outdoor air. Within RBCA, the solubility limit of the determinant restricts the extent of volatilisation, which in turn drives the indoor air inhalation pathway. Whilst the same restriction is not built into the CLEA model, the CLEA model output cells are flagged red where the soil saturation limit has been exceeded.

An assumption used in the CLEA model is that of simple linear partitioning of a chemical in the soil between the sorbed, dissolved and vapour phase<sup>(4)</sup>. The upper boundaries of this partitioning are represented by the aqueous solubility and pure saturated vapour concentration of the chemical. The CLEA software uses a traffic light system to identify when individual and/or combined assessment criteria exceed the lower of either the aqueous or vapour based saturation limits. Where model output cells are flagged red the soil or vapour saturation limit has been exceeded and further consideration of the SAC to be used within the assessment is required. One approach that could be adopted is to use the 'modelled' solubility saturation limit or vapour saturation limit of the compound as the SAC. However, as stated within the CLEA Handbook<sup>(4)</sup> this is likely to not be practical in many cases because of the very low limits



and is in any case highly conservative. Unless free-phase product is present, concentrations of the chemical are unlikely to be present at sufficient concentration to result in an exceedance of the health criteria value (HCV).

RSK has adopted an approach for petroleum hydrocarbons in accordance with LQM/CIEH<sup>(6)</sup> whereby the concentration modelled for each petroleum hydrocarbon fraction has been tabulated as the SAC with the corresponding solubility or vapour saturation limit given in brackets. Therefore, when using the SAC's to screen laboratory analysis the assessor should take note if a given SAC has a corresponding solubility or vapour saturation limit (in brackets), and subsequently incorporate this piece of information within the screening analytical discussion. If further assessment is required following this process then an additional approach can be utilised as detailed within Section 4.12 of the CLEA model manual<sup>(4)</sup> which explains how to calculate an effective assessment criterion manually.

### 3. Input Selection

Chemical data was obtained from EA Report SC050021/SR7<sup>(5)</sup> and the health criteria values (HCV) from the UK TOX<sup>(1)</sup> reports where available. For SAC for total petroleum hydrocarbons (TPH) and polycyclic aromatic hydrocarbons (PAH) toxicological and chemical specific parameters were obtained from the LQM/CIEH report <sup>(6)</sup>. Similarly, toxicological and specific chemical parameters for the volatile organic compound 1,2,4-trimethylbenzene were obtained from EIC/AGS/CL:AIRE<sup>(7)</sup>.

For total petroleum hydrocarbons (TPH), aromatic hydrocarbons  $C_5$ - $C_8$  were not modelled since benzene and toluene are being modelled separately. The aromatic  $C_8$ - $C_9$  hydrocarbon fraction comprises ethylbenzene, xylene and styrene. Since ethylbenzene and xylene are being modelled separately, the physical, chemical and toxicological data for this band has been taken from styrene.

Owing to the lack of UK-specific data, default information in the RBCA model was used to evaluate methyl tertiary butyl ether (MTBE). No published UK data was available for 1,3,5-trimethylbenzene, so information was obtained from the US EPA as in the RBCA model. RBCA uses toxicity data for the inhalation pathway in different units to the CLEA model and cannot consider separately the mean daily intake (MDI), occupancy periods or breathing rates. Therefore, the HCV in RBCA was amended to take account of:

- Amendments to the MDI using Table 3.4 of SR2<sup>(2)</sup>;
- A child weighing 13.3kg (average of 0-6 year old female in accordance with Table 4.6 of SR3<sup>(3)</sup>) and breathing 11.85m<sup>3</sup> (average daily inhalation rate for a 0-6 yr old female in accordance with Table 4.14 of SR3<sup>(3)</sup>; and
- The 50% rule (for petroleum hydrocarbons, 1,3,5-trimethylbenzene and MTBE)<sup>(2)</sup> where MDI data is not available but background exposure is considered important in the overall exposure.

### 4. Physical Parameters

For the residential with private gardens scenario, the CLEA default building is a small two-storey terrace house with concrete ground bearing slab. The house is assumed to have a 100m<sup>2</sup> private garden consisting of lawn, flowerbeds and incorporating a 20m<sup>2</sup> plot for growing fruit and vegetables consumed by the residents. SR3<sup>(3)</sup> notes this residential building type to be the most conservative in terms of protection from vapour intrusion. The building parameters are outlined in Table 5.

The parameters for a sandy loam soil type were used in line with SR3<sup>(3)</sup>. This includes a value of 6% for the percentage soil organic matter (SOM) within the soil. In RSK's experience, this is rather high for many sites. To avoid undertaking site specific risk assessments for this parameter, RSK has produced an



additional set of SAC for an SOM of 1% and 2.5%. For the GrAC, the depth to groundwater was taken as 2.5m based on RSK's experience of assessing the volatilisation pathway from groundwater.

### 5. GAC

The SAC were produced using the input parameters in Tables 1 to 5 and the GrAC using input parameters in Table 6. The final selected GAC are presented by pathway in Table 7 and the combined GAC in Table 8.



Figure 1
Conceptual Model for CLEA Residential Scenario – Private Gardens

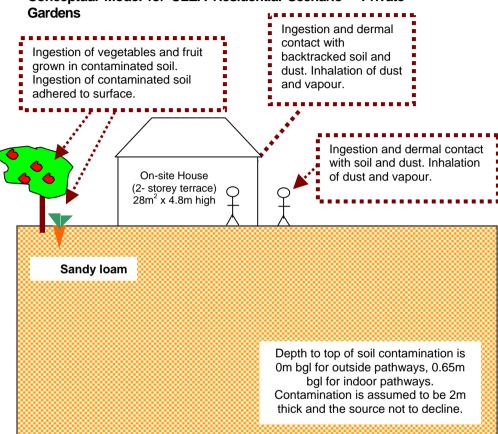


Table 1
Exposure Assessment Parameters for Residential Scenario - Private Gardens – Inputs for CLEA Model

Parameter	Value	Justification
Land use	Residential with homegrown produce	Chosen land use
Receptor	Female child age 1 to 6	Key generic assumption given in Box 3.1, report SC050021/SR3 <sup>(3)</sup>
Building	Small terraced house	Key generic assumption given in Box 3.1, report SC050021/SR3. Two storey small terraced house chosen as it is the most conservative residential building type in terms of protection from vapor intrusion (Section 3.4.6, report SC050021/SR3) <sup>(3)</sup>
Soil type	Sandy Loam	Most common UK soil type (Section 4.3.1, From Table 3.1, report SC050021/SR3) <sup>(3)</sup>
Start AC (age class)	1	Range of age classes corresponding to key generic assumption that the
End AC (age class)	6	critical receptor is a young female child aged zero to six. From Box 3.1, report SC050021/SR3 <sup>(3)</sup>
SOM (%)	6	Representative of sandy loamy soil according to EA Guidance note dated January 2009 entitled 'Changes We Have Made to the CLEA Framework Documents' (8)
	2.5	To provide SAC for sites where SOM <6% as often observed by RSK
рН	7	Model default



# Table 2 Residential with Private Gardens –Homegrown Produce Data for CLEA Model

	Con	sumpt day	ion Ra <sup>-1</sup> ) by <i>i</i>	te (g F Age Cl	W kg <sup>-1</sup> ass	BW	Dry Weight Conversion Factor	Homegrown Fraction (average)	Homegrown Fraction (high end)	Soil loading factor	Preparation correction factor
Name	1	2	3	4	5	6	g DW g <sup>-1</sup> FW	-	-	g g <sup>-1</sup> DW	-
Green vegetables	7.12	6.85	6.85	6.85	3.74	3.74	0.096	0.05	0.33	1.00E-03	2.00E-01
Root vegetables	10.69	3.30	3.30	3.30	1.77	1.77	0.103	0.06	0.4	1.00E-03	1.00E+00
Tuber vegetables	16.03	5.46	5.46	5.46	3.38	3.38	0.21	0.02	0.13	1.00E-03	1.00E+00
Herbaceous fruit	1.83	3.96	3.96	3.96	1.85	1.85	0.058	0.06	0.4	1.00E-03	6.00E-01
Shrub fruit	2.23	0.54	0.54	0.54	0.16	0.16	0.166	0.09	0.6	1.00E-03	6.00E-01
Tree fruit	3.82	11.96	11.96	11.96	4.26	4.26	0.157	0.04	0.27	1.00E-03	6.00E-01
Justification		Та	able 4.1	7, SR3	3 <sup>(3)</sup>		Table 6.3, SR3 <sup>(3)</sup>	Table 4.1	19, SR3 <sup>(3)</sup>	Table 6	5.3, SR3 <sup>(3)</sup>



Table 3
Residential with Private Gardens – Land Use Data for CLEA Model

Deserrator			OSC Data IOI		e Class					
Parameter	Unit	1	2	3	4	5	6			
EF (soil and dust ingestion)	day yr <sup>-1</sup>	180	365	365	365	365	365			
EF (consumption of homegrown produce)	day yr <sup>-1</sup>	180	365	365	365	365	365			
EF (skin contact, indoor)	day yr <sup>-1</sup>	180	365	365	365	365	365			
EF (skin contact, outdoor)	day yr <sup>-1</sup>	180	365	365	365	365	365			
EF (inhalation of dust and vapour, indoor)	day yr <sup>-1</sup>	365	365	365	365	365	365			
EF (inhalation of dust and vapour, outdoor)	day yr <sup>-1</sup>	365	365	365	365	365	365			
Justification	•			Table	3.1, SR3 <sup>(3)</sup>		•			
Occupancy period (indoor)	hr day <sup>-1</sup>	23	23	23	23	19	19			
Occupancy period (outdoor)	hr day <sup>-1</sup>	1	1	1	1	1	1			
Justification				Table	3.2, SR3 <sup>(3)</sup>					
Soil to skin adherence factor (indoor)	mg cm <sup>-2</sup> day <sup>-1</sup>	6.00E-02	6.00E-02	6.00E-02	6.00E-02	6.00E-02	6.00E-02			
Soil to skin adherence factor (outdoor)	mg cm <sup>-2</sup> day <sup>-1</sup>	1.00E+00	1.00E+00	1.00E+00	1.00E+00	1.00E+00	1.00E+00			
Justification		-		Table	8.1, SR3 <sup>(3)</sup>		-			
Soil and dust ingestion rate	g day <sup>-1</sup>	1.00E-01	1.00E-01	1.00E-01	1.00E-01	1.00E-01	1.00E-01			
Justification		Table 6.2, SR3 <sup>(3)</sup>								

Of note, for **cadmium**, the exposure assessment for a residential land use is based on estimates representative of lifetime exposure AC1-18. This is because the  $TDI_{oral}$  and  $TDI_{inh}$  – are based on considerations of the kidney burden accumulated over 50 years. It is therefore reasonable to consider exposure not only in childhood but averaged over a longer time period. See the Environment Agency Science report: SC05002 / TOX 3  $^{(1)}$  and Science Report SC050021 / Cadmium SGV  $^{(1)}$  for more information.

Table 4
Residential with Private Gardens – Receptor Data for CLEA Model

1. Coldonilla intant intanto candonio 1. Cocopto: Data 10: Clear inicato:											
Parameter	Unit			Age (	Class			Justification			
raiailletei	Offic	1	2	3	4	5	6	Justilication			
Body weight	kg	5.6	9.8	12.7	15.1	16.9	19.7	Table 4.6, SR3 <sup>(3)</sup>			
Body height	m	0.7	0.8	0.9	0.9	1	1.1				
Inhalation rate	m <sup>3</sup> day <sup>-1</sup>	8.5	13.3	12.7	12.2	12.2	12.2	Table 4.14, SR3 <sup>(3)</sup>			
Max exposed skin fraction (indoor)	m <sup>2</sup> m <sup>-2</sup>	0.32	0.33	0.32	0.35	0.35	0.33	Table 4.8, SR3 <sup>(3)</sup>			
Max exposed skin fraction (outdoor)	m <sup>2</sup> m <sup>-2</sup>	0.26	0.26	0.25	0.28	0.28	0.26	1 abic 4.0, 3N3			

See cadmium note as per Table 3 above.



Table 5
Residential with Private Gardens – Soil and Building Inputs for CLEA Model

Residential with Private Gardens			
Parameter	Unit	Value	Justification
		ROPERTIES for	sandy loam
Porosity, total	cm <sup>3</sup> cm <sup>-3</sup>	0.53	
Porosity, air filled	cm <sup>3</sup> cm <sup>-3</sup>	0.20	
Porosity, water filled	cm <sup>3</sup> cm <sup>-3</sup>	0.33	Default soil type is sandy loam, section 4.3.1,
Residual soil water content	cm <sup>3</sup> cm <sup>-3</sup>	0.12	SR3 <sup>(3)</sup>
Saturated hydraulic conductivity	cm s <sup>-1</sup>	3.56E-03	Parameters for sandy loam from Table 4.4, SR3 <sup>(3)</sup>
van Genuchten shape parameter (m)	-	3.20E-01	Since Since
Bulk density	g cm <sup>-3</sup>	1.21	
Threshold value of wind speed at 10m	m s <sup>-1</sup>	7.20	Default value taken from Section 9.2.2, SR3 <sup>(3)</sup>
Empirical function (F <sub>x</sub> ) for dust model	-	1.22	Value taken from Section 9.2.2, SR3 <sup>(3)</sup>
Ambient soil temperature	К	283	Annual average soil temperature representative of UK surface soils. Section 4.3.1, SR3 <sup>(3)</sup>
		DISPERSION MO	
Mean annual wind speed (10 m)	m s <sup>-1</sup>	5.00	Default value taken from Section 9.2.2, SR3 <sup>(3)</sup>
Air dispersion factor at height of 0.8 m	g m <sup>-2</sup> s <sup>-1</sup> per kg m <sup>-3</sup> g m <sup>-2</sup> s <sup>-1</sup>	2400	Values for a 0.01 ha site, appropriate to a residential land use in Newcastle (most
Air dispersion factor at height of 1.6 m	g m <sup>-2</sup> s <sup>-1</sup> per kg m <sup>-3</sup>	0	representative city for UK). (from Table 9.1, SR3) <sup>(3)</sup> Assumed child of 6 is not tall enough to reach 1.6m
Fraction of site with hard or vegetative cover	$\mathrm{m}^2\mathrm{m}^{-2}$	0.75	Section 3.2.6, SR3 (3) based on residential land use
BUILDING PROPE	RTIES for sr	mall terrace house	with ground-bearing floor slab
Building footprint	$m^2$	28	
Living space air exchange rate	hr <sup>-1</sup>	0.50	From Table 3.3 and 4.21, SR3 <sup>(3)</sup>
Living space height (above ground)	m	4.8	
Living space height (below ground)	m	0.0	Assumed no basement
Pressure difference (soil to enclosed space)	Pa	3.1	T 11 0 0 0Ds(3)
Foundation thickness	m	0.15	From Table 3.3, SR3 <sup>(3)</sup>
Floor crack area	cm <sup>2</sup>	423	
Dust loading factor	μg m <sup>-3</sup>	50	Default value for a residential site taken from Section 9.3, SR3 <sup>(3)</sup>
		VAPOUR MOD	EL
Default soil gas ingress rate	cm <sup>3</sup> s <sup>-1</sup>	25	Generic flow rate, Section 10.3, SR3 <sup>(3)</sup>
Depth to top of source (beneath building)	cm	50	Section 3.2.6, SR3 <sup>(3)</sup> states source is 50cm below building or 65cm below ground surface
Depth to top of source (no building)	cm	0	Section 10.2, SR3 <sup>(3)</sup> assumes impact from 0-1m for outdoor inhalation pathway
Thickness of contaminant layer	cm	200	Model default for indoor air, Section 4.9, SR4 <sup>(4)</sup>
Time average period for surface emissions	years	6	Time period of a 0 to 6 year old, Box 3.5, SR3 <sup>(3)</sup>
User-defined effective air permeability	cm <sup>2</sup>	3.05E-08	Calculated for sandy loam using equations in Appendix 1, SR3 <sup>(3)</sup>
-			• • • • • • • • • • • • • • • • • • • •



Figure 2 GrAC Conceptual Model for RBCA Residential with Private Gardens Scenario

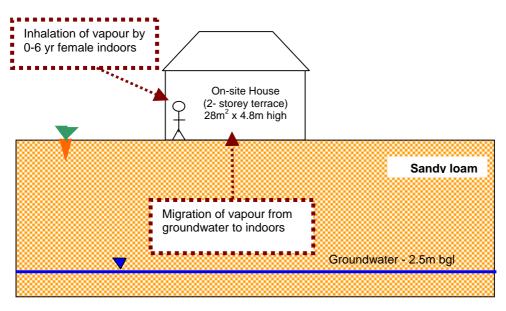


Table 6
Residential with Private Gardens RBCA Inputs

Parameter	Unit	Value	Justification
			RECEPTOR
Averaging time	Years	6	From Box 3.1, SR3 <sup>(3)</sup>
Receptor weight	kg	13.3	Average of CLEA 0-6 year old female data, Table 4.6, SR3 <sup>(3)</sup>
Exposure duration	Years	6	From Box 3.1, report, SR3 <sup>(3)</sup>
Exposure frequency	Days/yr	350	Weighted using occupancy period of 23 hours per day for 365 days of the year
		SOIL TY	PE – SANDY LOAM
Total porosity	-	0.53	
Volumetric water content	-	0.33	CLEA value for sandy loam. Parameters for sandy loam from
Volumetric air content	-	0.20	Table 4.4, SR3 <sup>(3)</sup>
Dry bulk density	g cm <sup>-3</sup>	1.21	
Vertical hydraulic conductivity	cm s <sup>-1</sup>	3.56E-3	CLEA value for saturated conductivity of sandy loam, Table 4.4, SR3 <sup>(3)</sup>
Vapour permeability	m <sup>2</sup>	3.05E-12	Calculated for sandy loam using equations in Appendix 1, SR3 <sup>(3)</sup>
Capillary zone thickness	m	0.1	Professional judgement
Fraction organic carbon	%	(i) 0.0348	Representative of sandy loam according to EA Guidance note dated January 2009 entitled Changes We Have Made to the CLEA Framework Documents (ref)
		(ii) 0.0058	To provide SAC for site's where SOM < 6% as often observed by RSK
			BUILDING
Building volume/area ratio	m	4.8	Table 3.3, SR3 <sup>(3)</sup>
Foundation area	m <sup>2</sup>	28	·
Foundation perimeter	m	22	Calculated assuming building measures 7m x 4m to give 28m <sup>2</sup> foundation area
Building air exchange rate	d <sup>-1</sup>	12	
Depth to bottom of foundation slab	m	0.15	Table 3.3, SR3 <sup>(3)</sup>
Foundation thickness	m	0.15	
Foundation crack fraction	-	0.0151	Calculated from floor crack area of 423 cm <sup>2</sup> and building footprint of 28m <sup>2</sup> in Table 4.21, SR3 <sup>(3)</sup>
Volumetric water content of cracks	-	0.33	Assumed equal to underlying soil type in assumption that cracks
Volumetric air content of cracks	-	0.2	become filled with soil over time. Parameters for sandy loam from Table 4.4, SR3 <sup>(3)</sup>
Indoor/outdoor differential pressure	Pa	3.1	From Table 3.3, SR3 <sup>(3)</sup>

Table 7

Human Health Generic Assessment Criteria by Pathway for Residential Scenario - Private Gardens



	Ž	GrAC	SAC Appropri	ate to Pathway S	OM 1% (ma/ka)		SAC Appropri	iate to Pathway SO	M 2 5% (ma/ka)		SAC Annrons	iate to Pathway S	OM 6% (ma/ka)	
Compound	Notes	(mg/l)	Oral	Inhalation	Combined	Soil Saturation Limit (mg/kg)	Oral	Inhalation	Combined	Soil Saturation	Oral	Inhalation	Combined	Soil Saturation
Compound	u	(mg/i)	Orai	innalation	Combined	Limit (mg/kg)	Orai	innalation	Combined	Limit (mg/kg)	Orai	innalation	Combined	Limit (mg/kg)
Metals														
Arsenic	(b)(c)	-	3.24E+01	8.50E+01	-	NR	3.24E+01	8.50E+01	-	NR	3.24E+01	8.50E+01	-	NR
Cadmium	(b)	-	1.12E+01	1.85E+02	1.10E+01	NR	1.12E+01	1.85E+02	1.10E+01	NR	1.12E+01	1.85E+02	1,10E+01	NR
Chromium (III) - oxide		-	1.84E+04	3.55E+03	2.98E+03	NR	1.84E+04	3.55E+03	2.98E+03	NR	1.84E+04	3.55E+03	2.98E+03	NR
Chromium (VI) - hexavalent		-	1.02E+01	4.25E+00	3.21E+00	NR	1.02E+01	4.25E+00	3.21E+00	NR	1.02E+01	4.25E+00	3,21E+00	NR
Copper		_	2.66E+03	1.04E+04	2.33E+03	NR	2.66E+03	1.04E+04	2.33E+03	NR	2.66E+03	1.04E+04	2.33E+03	NR NR
Lead	(a)	-	4.50E+02	-	-	NR	4.50E+02	-	-	NR	4.50E+02	-	-	NR
Elemental Mercury (Hq <sup>0</sup> )	(b)(d)	9.40E-03	-	1.70E-01	_	4.31E+00	-	4.24E-01	_	1.07E+01	-	1.02E+00	_	2.58E+01
Inorganic Mercury (Hg <sup>2+</sup> )	(b)	-	1.81E+02	2.55E+03	1.69E+02	NR	1.81E+02	2.55E+03	1.69E+02	NR	1.81E+02	2.55E+03	1.69E+02	NR
Methyl Mercury (Hq <sup>4+</sup> )	(b)	2.00E+01	1.39E+01	1.59E+01	7.40E+00	7.33E+01	1.39E+01	3.08E+01	9.55E+00	1.42E+02	1.39E+01	6.53E+01	1.14E+01	3.04E+02
Nickel	(b)(d)	-	5.31E+02	1.27E+02	-	NR	5.31E+02	1.27E+02	-	NR	5.31E+02	1.27E+02	-	NR
Selenium	(b)(c)	_	3.50E+02	1.272102	_	NR	3.50E+02	NR	_	NR	3.50E+02	-	-	NR
Zinc	(c)	-	3.75E+03	2.55E+07	_	NR	3.75E+03	2.55E+07	_	NR	3.75E+03	2.55E+07	_	NR
Cyanide	(0)		2.66E+01	3.97E+00	3.68E+00	NR	2.66E+01	3.97E+00	3.68E+00	NR	2.66E+01	3.97E+00	3.68E+00	NR
Cyariide			2.002+01	3.97 = +00	J.00L+00	INIX	2.00L+01	3.97 L+00	3.00L+00		2.00L+01	3.97 E+00	3.00L+00	IVIX
Volatile Organic Compounds														
Benzene	(b)	7.20E+00	1.12E-01	2.69E-01	7.92E-02	1.22E+03	2.28E-01	4.99E-01	1.57E-01	2.26E+03	4.89E-01	1.04E+00	3.32E-01	4.71E+03
Toluene	(b)	1.90E+03	1.47E+02	6.26E+02	1.19E+02	8.69E+02	3.35E+02	1.38E+03	2.70E+02	1.92E+03	7.59E+02	3.14E+03	6.11E+02	4.36E+03
Ethylbenzene	(b)	2.60E+02	1.06E+02	1.70E+02	6.52E+01	5.18E+02	2.51E+02	3.98E+02	1.54E+02	1.22E+03	5.70E+02	9.32E+02	3.54E+02	2.84E+03
Xylene - m	(b)	8.40E+01	2.02E+02	5.56E+01	4.36E+01	6.25E+02	4.80E+02	1.31E+02	1.03E+02	1.47E+03	1.09E+03	3.07E+02	2.40E+02	3.46E+03
	(b)	1.00E+02	1.85E+02	5.98E+01	4.52E+01	4.78E+02	4.80E+02 4.38E+02	1.40E+02	1.06E+02	1.12E+03	9.96E+02	3.07E+02 3.27E+02	2.46E+02	
Xylene - o	(6)	8.70E+01	1.91E+02	5.34E+01	4.17E+01		4.38E+02 4.51E+02	1.40E+02 1.26E+02	9.82E+01		9.96E+02 1.02E+03	3.27E+02 2.94E+02	2.28E+02	2.62E+03
Xylene - p		8.40E+01	2.02E+02	5.56E+01	4.17E+01 4.36E+01	5.76E+02		<del>                                     </del>	<del>                                     </del>	1.35E+03		<del>                                     </del>		3.17E+03
Total xylene		2.20E+03		1	1	6.25E+02	4.80E+02	1.31E+02	1.03E+02	1.47E+03	1.09E+03	3.07E+02	2.40E+02	3.46E+03
Methyl t-Butyl ether		1.80E+00	1.75E+00	1.84E+02	1.75E+00	1.66E+04	3.68E+00	2.40E+02	3.67E+00	2.16E+04	7.41E+00	3.70E+02	7.37E+00	3.34E+04
Trichloroethene		3.60E+00	2.83E+00	1.10E-01	1.06E-01	1.54E+03	6.25E+00	2.30E-01	2.22E-01	3.22E+03	1.40E+01	5.11E-01	4.93E-01	7.14E+03
Tetrachloroethene		2.60E+00	1.06E+01	1.03E+00	9.36E-01	4.24E+02	2.44E+01	2.30E+00	2.10E+00	9.51E+02	5.55E+01	5.28E+00	4.82E+00	2.18E+03
1,1,1-Trichloroethane			3.20E+02	6.33E+00	6.21E+00	1.43E+03	6.97E+02	1.29E+01	1.27E+01	2.92E+03	1.55E+03	2.84E+01	2.79E+01	6.39E+03
1,1,1,2 Tetrachloroethane		1.40E+01	5.19E+00	1.08E+00	8.93E-01	2.60E+03	1.22E+01	2.50E+00	2.08E+00	6.02E+03	2.78E+01	5.83E+00	4.82E+00	1.40E+04
1,1,2,2-Tetrachloroethane		1.40E+01	2.70E+00	2.76E+00	1.37E+00	2.67E+03	5.85E+00	5.65E+00	2.87E+00	5.46E+03	1.30E+01	1.24E+01	6.34E+00	1.20E+04
Carbon Tetrachloride		5.50E-02	1.05E+00	1.81E-02	1.79E-02	1.52E+03	2.41E+00	3.97E-02	3.93E-02	3.32E+03	5.44E+00	8.99E-02	8.92E-02	7.54E+03
1,2-Dichloroethane		3.00E-01	3.06E-02	6.46E-03	5.34E-03	3.41E+03	5.53E-02	9.32E-03	7.98E-03	4.91E+03	1.05E-01	1.60E-02	1.39E-02	8.43E+03
Vinyl Chloride		1.90E-02	3.69E-03	5.43E-04	4.73E-04	1.36E+03	6.64E-03	7.02E-04	6.35E-04	1.76E+03	1.21E-02	1.07E-03	9.86E-04	2.69E+03
1,2,4-Trimethylbenzene		7.50E-02	-	3.51E-01	-	5.57E+02	-	8.55E-01	-	1.36E+03	-	2.10E+00	-	3.25E+03
1,3,5-Trimethylbenzene		4.70E-02	1.45E+01	4.60E-01	4.56E-01	9.47E+01	3.47E+01	1.10E+00	1.09E+00	2.26E+02	7.94E+01	2.59E+00	2.56E+00	5.33E+02
Semi-Volatile Organic Compounds						•				1			T .	1
Acenaphthene		3.20E+00	2.18E+02	3.46E+03	2.05E+02	5.70E+01	5.08E+02	8.54E+03	4.79E+02	1.41E+02	1.06E+03	2.03E+04	1.01E+03	3.36E+02
Acenaphthylene		4.20E+00	1.78E+02	3.27E+03	1.68E+02	8.61E+01	4.17E+02	8.03E+03	3.97E+02	2.12E+02	8.90E+02	1.91E+04	8.51E+02	5.06E+02
Anthracene		2.10E-02	2.31E+03	1.08E+05	2.26E+03	1.17E+00	5.03E+03	2.65E+05	4.93E+03	2.91E+00	9.33E+03	6.15E+05	9.19E+03	6.96E+00
Benzo(a)anthracene		3.80E-03	7.00E+00	5.55E+00	3.10E+00	1.71E+00	8.98E+00	9.83E+00	4.69E+00	4.28E+00	1.01E+01	1.41E+01	5.88E+00	1.03E+01
Benzo(b)fluoranthene		2.00E-03	8.06E+00	1.79E+01	5.56E+00	1.22E+00	9.78E+00	1.97E+01	6.53E+00	3.04E+00	1.07E+01	2.05E+01	7.02E+00	7.29E+00
Benzo(g,h,i)perylene		2.60E-04	6.68E+01	1.27E+02	4.38E+01	1.54E-02	7.04E+01	1.32E+02	4.59E+01	3.85E-02	7.19E+01	1.34E+02	4.68E+01	9.23E-02
Benzo(k)fluoranthene		8.00E-04	1.25E+01	2.66E+01	8.51E+00	6.87E-01	1.44E+01	2.83E+01	9.56E+00	1.72E+00	1.53E+01	2.91E+01	1.00E+01	4.12E+00
Chrysene		2.00E-03	8.76E+00	1.95E+01	6.00E+00	4.40E-01	1.20E+01	2.45E+01	8.04E+00	1.10E+00	1.41E+01	2.72E+01	9.27E+00	2.64E+00
Dibenzo(a,h)anthracene		6.00E-04	1.19E+00	2.13E+00	7.62E-01	3.93E-03	1.33E+00	2.42E+00	8.58E-01	9.82E-03	1.39E+00	2.56E+00	9.03E-01	2.36E-02
Fluoranthene		2.30E-01	2.59E+02	2.69E+04	2.57E+02	1.89E+01	4.67E+02	6.23E+04	4.63E+02	4.73E+01	6.78E+02	1.28E+05	6.74E+02	1.13E+02
Fluorene		1.90E+00	1.70E+02	4.35E+03	1.63E+02	3.09E+01	3.91E+02	1.07E+04	3.77E+02	7.65E+01	8.00E+02	2.54E+04	7.76E+02	1.83E+02
Indeno(1,2,3-cd)pyrene		2.00E-04	4.58E+00	1.04E+01	3.18E+00	6.13E-02	5.74E+00	1.17E+01	3.85E+00	1.53E-01	6.37E+00	1.22E+01	4.19E+00	3.68E-01
Phenanthrene		5.30E-01	9.35E+01	5.04E+03	9.18E+01	3.60E+01	2.04E+02	1.23E+04	2.01E+02	8.96E+01	3.81E+02	2.86E+04	3.76E+02	2.14E+02
Pyrene		1.30E-01	5.69E+02	6.18E+04	5.63E+02	2.20E+00	1.05E+03	1.44E+05	1.04E+03	5.49E+00	1.56E+03	2.97E+05	1.56E+03	1.32E+01
Benzo(a)pyrene		3.80E-03	1.21E+00	2.62E+00	8.26E-01	9.11E-01	1.42E+00	2.81E+00	9.43E-01	2.28E+00	1.52E+00	2.90E+00	9.98E-01	5.46E+00
	_	1.90E+01	2.68E+01	1.64E+00	1.54E+00	7.64E+01	6.36E+01	3.93E+00	3.70E+00	1.83E+02	1.43E+02	9.27E+00	8.71E+00	4.32E+02
Naphthalene														

Table 7 RSK GAC\_2010\_03\_Rev01

#### Table 7

Human Health Generic Assessment Criteria by Pathway for Residential Scenario - Private Gardens



	ON	GrAC	SAC Appropri	ate to Pathway So	OM 1% (mg/kg)	Soil Saturation	SAC Appropri	ate to Pathway SOI	VI 2.5% (mg/kg)	Soil Saturation	SAC Appropri	iate to Pathway S	OM 6% (mg/kg)	Soil Saturation
Compound	tes	(mg/l)	Oral	Inhalation	Combined	Limit (mg/kg)	Oral	Inhalation	Combined	Limit (mg/kg)	Oral	Inhalation	Combined	Limit (mg/kg)
Total Petroleum Hydrocarbons														
Aliphatic hydrocarbons EC <sub>5</sub> -EC <sub>6</sub>		1.00E+01	4.79E+03	2.98E+01	2.97E+01	3.04E+02	1.08E+04	5.47E+01	5.46E+01	5.58E+02	2.35E+04	1.13E+02	1.13E+02	1.15E+03
Aliphatic hydrocarbons >EC <sub>6</sub> -EC <sub>8</sub>		5.40E+00	1.43E+04	7.27E+01	7.26E+01	1.44E+02	3.21E+04	1.62E+02	1.62E+02	3.22E+02	6:36E+04	3.72E+02	3.71E+02	7.36E+02
Aliphatic hydrocarbons >EC <sub>8</sub> -EC <sub>10</sub>		2.30E-01	1.46E+03	1.89E+01	1.88E+01	7.77E+01	2.44E+03	4.60E+01	4.58E+01	1.90E+02	3.30E+03	1.09E+02	1.08E+02	4.51E+02
Aliphatic hydrocarbons >EC <sub>10</sub> -EC <sub>12</sub>		3.40E-02	3.52E+03	9.34E+01	9.28E+01	4.75E+01	4.01E+03	2:32E+02	2.29E+02	1.18E+02	4.24E+03	5.57E+02	5.37E+02	2.83E+02
Aliphatic hydrocarbons >EC <sub>12</sub> -EC <sub>16</sub>		7.60E-04	4.37E+03	7.82E+02	7.44E+02	2.37E+01	4.40E+03	1.95E+03	1.69E+03	5.91E+01	4.41E+03	4.68E+03	3.03E+03	1.42E+00
Aliphatic hydrocarbons >EC <sub>16</sub> -EC <sub>35</sub>	(c)	-	4.51E+04	-	-	8.48E+00	6.38E+04	-	-	2.12E+01	7.61E+04	-	-	5.09E+01
Aliphatic hydrocarbons >EC <sub>35</sub> -EC <sub>44</sub>	(c)	-	4.51E+04	-	-	8.48E+00	6.38E+04	-	-	2.12E+01	7.61E+04	-	-	5.09E+01
Aromatic hydrocarbons >EC <sub>8</sub> -EC <sub>9</sub> (styrene)		7.40E+00	1.66E+02	2.65E+02	1.33E+02	6.20E+02	3.92E+02	6.47E+02	3.16E+02	1.52E+03	8.50E+02	1.54E+03	7.02E+02	3.61E+03
Aromatic hydrocarbons >EC <sub>9</sub> -EC <sub>10</sub>		7.40E+00	5.55E+01	3.33E+01	2.69E+01	6.13E+02	1.31E+02	8.16E+01	6.54E+01	1.50E+03	2.84E+02	1.94E+02	1.51E+02	3.58E+02
Aromatic hydrocarbons >EC <sub>10</sub> -EC <sub>12</sub>		2.50E+01	7.97E+01	1.82E+02	6.91E+01	3.64E+02	1.86E+02	4.48E+02	1.62E+02	8.99E+02	3.87E+02	1.07E+03	3.46E+02	2.15E+03
Aromatic hydrocarbons >EC <sub>12</sub> -EC <sub>16</sub>		5.80E+00	1.40E+02	2.00E+03	1.38E+02	1.69E+02	3.13E+02	4.96E+03	3.08E+02	4.19E+02	6.01E+02	1.18E+04	5.93E+02	1.00E+03
Aromatic hydrocarbons >EC <sub>16</sub> -EC <sub>21</sub>	(c)	-	2.47E+02	-	-	5.37E+01	4.82E+02	-	-	1.34E+02	7.66E+02	-	-	3.21E+02
Aromatic hydrocarbons >EC <sub>21</sub> -EC <sub>35</sub>	(c)	-	8.88E+02	-	-	4.83E+00	1.11E+03	-	-	1.21E+01	1.22E+03	-	-	2.90E+01
Aromatic hydrocarbons >EC <sub>35</sub> -EC <sub>44</sub>	(c)	-	8.88E+02	-	-	4.83E+00	1.11E+03	-	-	1.21E+01	1.22E+03	-	-	2.90E+01

#### Notes:

-' Generic assessment criteria not calculated owing to low volatility of substance and therefore no pathway, or an absence of toxicological data.

NR - the compound is not volatile and therefore a soil saturation limit not calculated within CLEA

EC - equivalent carbon. GrAC - groundwater assessment criteria. SAC - soil assessment criteria.

The CLEA model output is colour coded depending upon whether the soil saturation limit has been exceeded.



Calculated SAC exceeds soil saturation limit and may significantly effect the interpretation of any exceedances since the contribution of the indoor and outdoor vapour pathway to total exposure is

>10%. This shading has also been used for the RBCA output where the theoretical solubility limit has been exceeded. The SAC has been set as the model calculated SAC with the saturation limits shown in brackets. Calculated SAC exceeds soil saturation limit but will not effect the SSV significantly since the contribution of the indoor and outdoor vapour pathway to total exposure is <10%. Calculated SAC does not exceed the soil saturation limit.

For consistency where the theoretical solubility limit within RBCA has been exceeded in production of the GrAC, these cellls have also been hatched red.

The SAC for organic compounds are dependant upon soil organic matter (SOM) (%) content. To obtain SOM from total organic carbon (TOC) (%) divide by 0.58. 1% SOM is 0.58% TOC. DL Rowell Soil Science: Methods and Applications, Longmans, 1994. SAC for TPH fractions, polycyclic aromatic hydrocarbons, MTBE, BTEX and trimethylbenzene compounds were produced using an attenuation factor for the indoor air inhalation pathway of 10 to reduce conservatism associated with the vapour

inhalation pathway, section 10.1.1, SR3

(a) GAC taken as former Soil Guideline Value owing to uncertainty regarding toxicological approach to be adopted by the Environment Agency.

- (b) GAC taken from the Environment Agency SGV reports published 2009.
- (c) SAC for selenium, aliphatic and aromatic hydrocarbons >EC16 does not include inhalation pathway owing to absence of toxicity data. SAC for arsenic is only based on oral contribution (rather than combined) owing to the relative small contribution from inhalation in accordance with the SGV report. The same approach has been adopted for zinc.
- (d) SAC for elemental mercury, chromium VI and nickel is based on the inhalation pathway only owing to an absence of toxicity for elemental mercury, in accordance with the SGV report for nickel and LQM report for chromium VI.



Table 8 Human Health Generic Assessment Criteria for Residential Scenario - Private Gardens

Compound	GrAC for Groundwater (mg/l)	SAC for Soil SOM 1% (mg/kg)	SAC for Soil SOM 2.5% (mg/kg)	SAC for Soil SOM 6% (mg/kg)
Metals				
Arsenic	-	32	32	32
Cadmium	-	10	10	10
Chormium (III) - oxide	-	3,000	3,000	3,000
Chromium (VI) - hexavalent	-	4.3	4.3	4.3
Copper	-	2,300	2,300	2,300
Lead	-	450	450	450
Elemental Mercury (Hg <sup>0</sup> )	0.009	0.17	0.42	1.0
Inorganic Mercury (Hg <sup>2+</sup> )	-	170	170	170
Methyl Mercury (Hg4+)	20	7.4	9.6	11
Nickel	-	130	130	130
Selenium	-	350	350	350
Zinc	-	3,800	3,800	3,800
Cyanide	-	3.7	3.7	3.7
Volatile Organic Compounds				
Benzene	7	0.079	0.157	0.33
Toluene	1,900	120	270	610
Ethylbenzene	260	65	154	350
Xylene - m	100	44	103	240
Xylene - o	87	45	106	250
Xylene - p	84	42	98	230
Total xylene	84	44	103	240
Methyl tertiary butyl ether (MTBE)	2,200	1.8	3.7	7.4
Trichloroethene	1.8	0.11	0.2	0.49
Tetrachloroethene	3.6 26	0.94 6.2	2.1 12.7	4.8 28
1,1,1-Trichloroethane	14	0.89	2.1	4.8
1,1,1,2Tetrachloroethane 1,1,2,2-Tetrachloroethane	14	1.4	2.87	6.3
Carbon Tetrachloride	0.055	0.018	0.039	0.089
1,2-Dichloroethane	0.30	0.0053	0.0080	0.009
Vinyl Chloride	0.019	0.00047	0.0006	0.001
1,2,4-Trimethylbenzene	0.075	0.35	0.85	2.1
1,3,5-Trimethylbenzene	0.047	0.46	1.1	2.6
Semi-Volatile Organic Compounds				
Acenaphthene	3.2	210	480	1,000
Acenaphthylene	4.2	170	400	850
Anthracene	0.021	2,300	4,900	9,200
Benzo(a)anthracene	0.0038	3.1	4.7	5.9
Benzo(b)fluoranthene	0.0020	5.6	6.5	7.0
Benzo(g,h,i)perylene	0.00026	44	46	47
Benzo(k)fluoranthene	0.00080	8.5	9.6	10
Chrysene	0.0020	6.0	8.0	9.3
Dibenzo(a,h)anthracene	0.00060	0.76	0.86	0.90
Fluoranthene	0.23	260	460	670
Fluorene	1.9	160	380	780
Indeno(1,2,3-cd)pyrene	0.0002	3.2	3.8	4.2
Phenanthrene	0.53	92	200	380
Pyrene	0.13	560	1,000	1,600
Benzo(a)pyrene Naphthalene	0.0038	0.83	0.94	1.0
Naphthalene Phenol	19	1.5 180	3.7 290	8.7 420
	-	180	290	420
Total Petroleum Hydrocarbons  Aliphatic hydrocarbons EC <sub>5</sub> -EC <sub>6</sub>	10	30	55	110
Aliphatic hydrocarbons > EC <sub>6</sub> -EC <sub>8</sub>	5.4	73	160	370
Aliphatic hydrocarbons >EC <sub>8</sub> -EC <sub>10</sub>	0.23	19 <b>93 (48)</b>	46	110
Aliphatic hydrocarbons >EC <sub>10</sub> -EC <sub>12</sub> Aliphatic hydrocarbons >EC <sub>12</sub> -EC <sub>16</sub>	0.034 0.00076	744 (24)	230 (118) 1,700 (59)	540 (283) 3,000 (142)
Aliphatic hydrocarbons >EC <sub>16</sub> -EC <sub>35</sub>	-	45,100 (8.48)	64,000 (21)	76,000
Aliphatic hydrocarbons >EC <sub>35</sub> -EC <sub>44</sub>	-	45.100 (8.48)	64,000 (21)	76,000
Aromatic hydrocarbons >EC <sub>8</sub> -EC <sub>9</sub> (styrene)	7.4	130	316	700
Aromatic hydrocarbons >EC <sub>9</sub> -EC <sub>10</sub>	7.4	27	65	150
Aromatic hydrocarbons >EC <sub>10</sub> -EC <sub>12</sub>	25	69	160	346
Aromatic hydrocarbons >EC <sub>12</sub> -EC <sub>16</sub>	5.8	140	310	593
Aromatic hydrocarbons >EC <sub>16</sub> -EC <sub>21</sub>	-	250	480	770
Aromatic hydrocarbons >EC <sub>21</sub> -EC <sub>35</sub>	_	890	1,100	1,230
Aromatic hydrocarbons >EC <sub>35</sub> -EC <sub>44</sub>	-	890	1,100	1,230
	·	030	1,100	1,230

#### Notes

The SAC has been set as the model calculated SAC with the saturation limit shown in brackets.

For consistency where the GrAC exceeds the solubility limit, GrAC has been set at the solubility limit. The GrAC conservative since concentrations of the chemical are very unlikely to be at sufficient concentration to result in an exceedance of the health criteria value at the point of exposure (i.e. indoor air) provided free-phase product is absent.

Table 8 RSK GAC 2010 03 Rev01

<sup>-</sup> Generic assessment criteria not calculated owing to low volatility of substance and therefore no pathway, or an absence of toxicological data.

EC - equivalent carbon. GrAC - groundwater assessment criteria. SAC - soil assessment criteria.

The SAC for organic compounds are dependent on Soil Organic Matter (SOM) (%) content. To obtain SOM from total organic carbon (TOC) (%) divide by 0.58. 1% SOM is 0.58% TOC. DL Rowell Soil Science: Methods and Applications, Longmans, 1994.

SAC for TPH fractions, polycyclic aromatic hydrocarbons, MTBE, BTEX and trimethylbenzene compounds were produced using an attenuation factor for the indoor air inhalation pathway of 10 to reduce conservatism associated with the vapour inhalation pathway, section 10.1.1, SR3.



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- 7) The Soil Generic Assessment Criteria for Human Health Risk Assessment, Report ref. ISBN 978-1-905046-20-1, December 2009, published by CL:AIRE
- 8) Changes made to the CLEA framework documents after the 3 month evaluation period in 2008, released January 2009 by the Environment Agency.



# Generic Assessment Criteria (GAC) for Human Health Residential Scenario – Without Gardens

The human health generic assessment criteria (GAC) have been developed during a period of regulatory review and updating of the Contaminated Land Exposure Assessment (CLEA) project. Hence, the Environment Agency (EA) is in the process of publishing updated reports relating to the CLEA project and the GAC presented in this document may change to reflect these updates. This issue was prepared following the publication of soil guideline value reports and associated publications<sup>(1)</sup> for mercury, selenium, benzene, toluene, ethylbenzene and xylene in March 2009 plus arsenic and nickel in May 2009, cadmium and phenol in June 2009, dioxins, furans and dioxin-like PCB's in September 2009. It was also produced following publication of GAC by LQM<sup>(6)</sup>. Where available, the published soil guideline values (SGV)<sup>(1)</sup> were used as the GAC.

#### 1. Model Selection

Soil assessment criteria (SAC) were calculated using CLEA v1.06 and the supporting UK guidance<sup>(1-6)</sup>. Groundwater assessment criteria (GrAC) protective of human health via the inhalation pathway were derived using the RBCA 1.3b model. RSK has updated the inputs within RBCA to reflect the UK guidance<sup>(1-5)</sup>. The SAC and GrAC collectively are termed GAC.

#### 2. Conceptual Model

In accordance with EA Science Report SC050021/SR3<sup>(3)</sup>, the residential without gardens scenario considers risks to a female child between the ages of 0 and 6 years old. In accordance with Box 3.1, SR3, the pathways considered for production of the SAC in the residential without gardens scenario are:

- Direct soil and dust ingestion;
- Dermal contact with soil and indoor dust, and
- Inhalation of indoor and outdoor dust and vapours.

Figure 1 is a conceptual model illustrating these linkages.

The pathway considered in production of the GrAC is the volatilisation of compounds from groundwater and subsequent vapour inhalation by residents whilst indoors. Figure 2 illustrates this linkage. Although the outdoor air inhalation pathway is also valid, this contributes little to the overall risks owing to the dilution in outdoor air. Within RBCA, the solubility limit of the determinant restricts the extent of volatilisation, which in turn drives the indoor air inhalation pathway. Whilst the same restriction is not built into the CLEA model, the CLEA model output cells are flagged red where the soil saturation limit has been exceeded.

An assumption used in the CLEA model is that of simple linear partitioning of a chemical in the soil, between the sorbed, dissolved and vapour phase<sup>(4)</sup>. The upper boundaries of this partitioning are represented by the aqueous solubility and pure saturated vapour concentration of the chemical. The CLEA software uses a traffic light system to identify when individual and/or combined assessment criteria exceed the lower of either the aqueous or vapour based saturation limits. Where model output cells are flagged red the soil or vapour saturation has been exceeded and further consideration of the SAC to be used within the assessment is required. One approach that could be adopted is to use the 'modelled' solubility saturation limit or vapour saturation limit of the compound as the SAC. However, as stated within the CLEA Handbook<sup>(4)</sup> this is likely not to be practical in many cases because of the subsequent very low solubility/vapour saturation limits and is in any case highly conservative and unless free-phase product is present, concentrations of the chemical are unlikely to be present at sufficient concentration to result in an exceedance of the health criteria value (HCV).



RSK has adopted an approach for petroleum hydrocarbons in accordance with LQM/CIEH<sup>(6)</sup> whereby the concentration modelled for each petroleum hydrocarbon fraction has been tabulated as the SAC with the corresponding solubility or vapour saturation limits given in brackets. Therefore, when using the SAC to screen laboratory analysis the assessor should take note if a given SAC has a corresponding solubility or vapour saturation limit (in brackets) and subsequently incorporate this information within the screening analytical discussion. If further assessment is required following this process then an additional approach can be utilised as detailed within Section 4.12 of the CLEA model Handbook<sup>(4)</sup> which explains how to calculate an effective assessment criterion manually.

#### 3. Input Selection

Chemical data was obtained from EA Report SC050021/SR7<sup>(5)</sup> and the health criteria values (HCV) from the UK TOX<sup>(1)</sup> reports where available. For total petroleum hydrocarbons (TPH) and polycyclic aromatic hydrocarbons (PAH) toxicological and chemical specific parameters were obtained from the LQM/CIEH report<sup>(6)</sup>. Similarly, toxicological and specific chemical parameters for the volatile organic compound 1,2,4-trimethylbenzene were obtained from EIC/AGS/CL:AIRE<sup>(7)</sup>.

For total petroleum hydrocarbons (TPH), aromatic hydrocarbons  $C_5$ - $C_8$  were not modelled since benzene and toluene are being modelled separately. The aromatic  $C_8$ - $C_9$  hydrocarbon fraction comprises ethylbenzene, xylene and styrene. Since ethylbenzene and xylene are being modelled separately, the physical, chemical and toxicological data for this band has been taken from styrene.

Owing to the lack of UK-specific data, default information in the RBCA model was used to evaluate methyl tertiary butyl ether (MTBE). No published UK data was available for 1,3,5-trimethylbenzene, so information was obtained from the RBCA model. RBCA uses toxicity data for the inhalation pathway in different units to the CLEA model and cannot consider separately the mean daily intake (MDI), occupancy periods or breathing rates. Therefore, the HCV in RBCA was amended to take account of:

- Amendments to the MDI using Table 3.4 of SR2<sup>(2)</sup>;
- A child weighing 13.3kg (average of 0-6 year old female in accordance with Table 4.6 of SR3<sup>(3)</sup>) and breathing 11.85m<sup>3</sup> (average daily inhalation rate for a 0-6yr old female in accordance with Table 4.14 of SR3<sup>(3)</sup>; and
- The 50% rule (for petroleum hydrocarbons, trimethylbenzenes and MTBE)<sup>(2)</sup> where MDI data is not available but background exposure is considered important in the overall exposure.

#### 4. Physical Parameters

For the residential without gardens scenario, the CLEA default building is a small two-storey terrace house with concrete ground bearing slab. SR3<sup>(3)</sup> notes this residential building type to be the most conservative in terms of protection from vapour intrusion. The building parameters are outlined in Table 3.

The parameters for a sandy loam soil type were used in line with SR3<sup>(3)</sup>. This includes a value of 6% for the percentage soil organic matter (SOM) within the soil. In RSK's experience, this is rather high for many sites. To avoid undertaking site-specific risk assessments for this parameter, RSK has produced an additional set of SAC for an SOM of 1% and 2.5%.

For the GrAC, the depth to groundwater was taken as 2.5m based on RSK's experience of assessing the volatilisation pathway from groundwater.



### <u>5. GAC</u>

The SAC were produced using the input parameters in Tables 1 to 3 and the GrAC using the input parameters in Table 4. The GAC by pathway are presented in Table 5 with the combined GAC presented in Table 6.



Figure 1 Conceptual Model for CLEA Residential Scenario – without Gardens

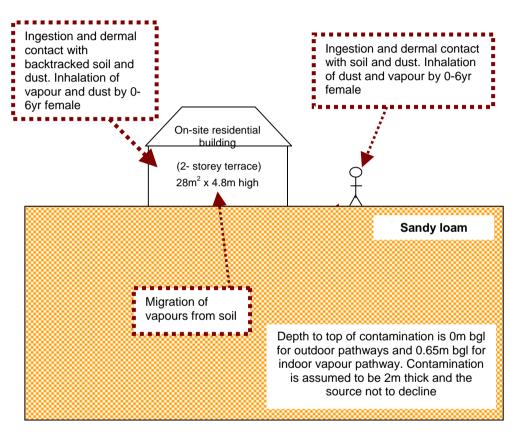


Table 1
Exposure Assessment Parameters for Residential Scenario – without Gardens – Inputs for CLEA model.

Parameter	Value	Justification
Land use	Residential without homegrown produce	Chosen land use
Receptor	Female Child	Taken as female child exposed over 6 years from 0-6 years, Box 3.1, SR3 <sup>(3)</sup>
Building	Small terraced house	Key generic assumption given in Box 3.1, SR3 <sup>(3)</sup> . Two storey small terraced house chosen as it is the most conservative residential building type in terms of protection from vapour intrusion (Section 3.2.6, report SC050021/SR3 <sup>(3)</sup> ). Table 3 presents building specific input data
Soil type	Sandy loam	Most common UK soil type (Section 4.3.1, Table 4.4, SR3 <sup>(3)</sup> ). Table 4 presents soil-specific input data
Start age class (AC)	1	Range of AC corresponding to key generic assumption that the critical receptor is a young female child aged
End AC	6	zero to six. From Box 3.1, SR3 <sup>(3)</sup> .  Data specific to the receptor is presented in Table 2
SOM (%)	6	Representative of sandy loam according to EA Guidance note dated January 2009 entitled 'Changes We Have Made to the CLEA Framework Documents' <sup>(8)</sup>
	1	To provide SAC for site's where SOM
pH	2.5	< 6% as often observed by RSK  Model default
ווק	ı	เขเบนต์ นต์เลนเเ



Table 2
Residential without Gardens – Land use and Receptor Data for CLEA Model

Parameter	Unit			Age	Class		
		1	2	3	4	5	6
Exposure frequency (EF) (soil and dust ingestion)	day yr <sup>-1</sup>	180	365	365	365	365	365
EF (skin contact, indoor)	day yr <sup>-1</sup>	180	365	365	365	365	365
EF (skin contact, outdoor)	day yr <sup>-1</sup>	180	365	365	365	365	365
EF (inhalation of dust and vapour, indoor)	day yr <sup>-1</sup>	365	365	365	365	365	365
EF (inhalation of dust and vapour, outdoor)	day yr <sup>-1</sup>	365	365	365	365	365	365
Justification				Table 3	.1, SR3 <sup>(3)</sup>		
Occupancy period (indoor)	hr day <sup>-1</sup>	23	23	23	23	19	19
Occupancy period (outdoor)	hr day <sup>-1</sup>	1	1	1	1	1	1
Justification				Table 3	.2, SR3 <sup>(3)</sup>		
Soil ingestion rate	g/day	0.1	0.1	0.1	0.1	0.1	0.1
Justification				Table 6	.2, SR3 <sup>(3)</sup>		
Soil to skin adherence factor – (indoor)	mg soil/cm <sup>2</sup> skin	0.06	0.06	0.06	0.06	0.06	0.06
Soil to skin adherence factor – (outdoor)	mg soil/cm <sup>2</sup> skin	1	1	1	1	1	1
Justification				Table 8	.1, SR3 <sup>(3)</sup>		
Body Weight	kg	5.6	9.8	12.7	15.1	16.9	19.7
Body height	m	0.7	0.8	0.9	0.9	1	1.1
Justification				Table 4	.6, SR3 <sup>(3)</sup>		
Inhalation Rate	m <sup>3</sup> day <sup>-1</sup>	8.5	13.3	12.7	12.2	12.2	12.2
Justification				Table 4.	14, SR3 <sup>(3)</sup>		
Max exposed skin fraction (indoor)	m <sup>2</sup> m <sup>-2</sup>	0.32	0.33	0.32	0.35	0.35	0.33
Max exposed skin fraction (outdoor)	m <sup>2</sup> m <sup>-2</sup>	0.26	0.26	0.25	0.28	0.28	0.26
Justification				Table 4	.8, SR3 <sup>(3)</sup>		

Note: for **cadmium**, the exposure assessment for a residential land use is based on estimates representative of lifetime exposure AC1-18. This is because the  $TDI_{oral}$  and  $TDI_{inh}$  are based on considerations of the kidney burden accumulated over 50 years. It is therefore reasonable to consider exposure not only in childhood but averaged over a longer time period. See the Environment Agency Science report:  $SC050021 / TOX 3^{(1)}$  and Science Report  $SC050021 / Cadmium SGV^{(1)}$  for the full AC1-18 Land use Data suite.



Table 3
Residential without Gardens – Soil, Air and Building Specific Inputs for CLEA Model

Parameter	Unit	Value	Justification
	SOIL P	ROPERTIES for	sandy loam
Porosity, total	cm <sup>3</sup> <sub>3</sub> cm <sup>-</sup>	0.53	
Porosity, air filled	cm <sup>3</sup> cm <sup>-</sup>	0.20	
Porosity, water filled	cm <sup>3</sup> cm <sup>-</sup>	0.33	Default soil type is sandy loam, section 4.3.1, SR3
Residual soil water content	cm <sup>3</sup> <sub>3</sub> cm <sup>-</sup>	0.12	Parameters for sandy loam from Table 4.4, SR3 <sup>(3)</sup>
Saturated hydraulic conductivity	cm s <sup>-1</sup>	0.00356	
Van Genuchten shape parameter (m)	-	0.3201	
Bulk density	g cm <sup>-3</sup>	1.21	
Threshold value of wind speed at 10m	m s <sup>-1</sup>	7.2	Default value taken from Section 9.2.2, SR3 <sup>(3)</sup>
Empirical function (F <sub>x</sub> ) for dust model	-	1.22	Value taken from Section 9.2.2, SR3 <sup>(3)</sup>
Ambient soil temperature	К	283	Annual average soil temperature of UK surface soils. Section 4.3.1, SR3 <sup>(3)</sup>
	Al	R DISPERSION	
Mean annual wind speed (10 m)	m s <sup>-1</sup>	5.0	Default value taken from Section 9.2.2, SR3 <sup>(3)</sup>
Air dispersion factor at height of 0.8 m	g m <sup>-2</sup> s <sup>-1</sup> per kg m <sup>-3</sup>	2400	From Table 9.1, SR3 <sup>(3)</sup> . Values for a 0.01 ha site, appropriate to a residential land use in Newcastle (representative city for UK, section 9.2.1, SR3 <sup>(3)</sup> )
Fraction of site with hard or vegetative cover	$m^2 m^{-2}$	0.75	Section 3.2.6, SR3 <sup>(3)</sup> for residential land use
BUILDING PR		S for house wit	h ground-bearing floor slab
Building footprint	$m^2$	28	
Living space air exchange rate	hr <sup>-1</sup>	0.50	From Table 3.3 and 4.21, SR3 <sup>(3)</sup>
Living space height (above ground)	m	4.8	
Living space height (below ground)	m	0.0	Assumed no basement
Pressure difference (soil to enclosed space)	Pa	3.1	5 T. H. O.O. H.4.04 ODO(3)
Foundation thickness	m	0.15	From Table 3.3 and 4.21, SR3 <sup>(3)</sup>
Floor crack area	cm <sup>2</sup>	423	
Dust loading factor	μg m <sup>-3</sup>	50	Default value for a residential site taken from Section 9.3, SR3 <sup>(3)</sup>
		VAPOUR MOD	
Default soil gas ingress rate	cm <sup>3</sup> s <sup>-1</sup>	25	Generic flow rate, Section 10.3, SR3 <sup>(3)</sup>
Depth to top of source (beneath building for indoor exposure)	cm	50	Section 3.2.6, SR3 <sup>(3)</sup> states source is 50cm below building or 65cm below ground surface
Depth to top of source (outdoors)	cm	0	Section 10.2, SR3 <sup>(3)</sup> assumes impact from 0-1m for outdoor inhalation pathway
Thickness of contaminant layer	cm	200	Model default for indoor air, Section 4.9, SR4 <sup>(4)</sup>
Time average period for surface emissions	years	6	Time period of a 0–6 year old, Box 3.5, SR3 <sup>(3)</sup>
User-defined effective air permeability	cm <sup>2</sup>	3.05E-08	Calculated for sandy loam using equations in Appendix 1, SR3 <sup>(3)</sup>



Figure 2
GrAC Conceptual Model for RBCA Residential without Gardens Scenario

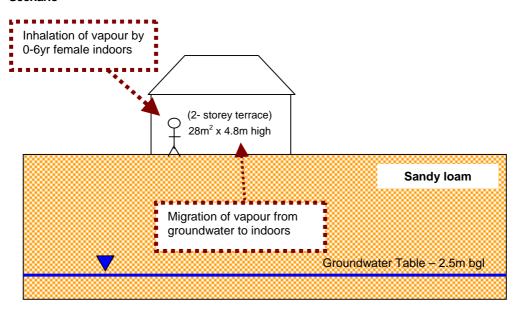


Table 4
Residential without Gardens RBCA Inputs

Residential without Garde			
Parameter	Unit	Value	Justification
			RECEPTOR
Averaging time	Years	6	From Box 3.1, SR3 <sup>(3)</sup>
Receptor weight	kg	13.3	Average of CLEA 0-6 year old female data, Table 4.6, SR3 <sup>(3)</sup>
Exposure duration	Years	6	From Box 3.1, report, SR3 <sup>(3)</sup>
Exposure frequency	Days/yr	350	Weighted using occupancy period of 23 hours per day for 365 days of the year
		SOIL TY	PE - SANDY LOAM
Total porosity	-	0.53	
Volumetric water content	-	0.33	CLEA value for sandy loam. Parameters for sandy loam from
Volumetric air content	-	0.20	Table 4.4, SR3 <sup>(3)</sup>
Dry bulk density	g cm <sup>-3</sup>	1.21	
Vertical hydraulic conductivity	cm s <sup>-1</sup>	3.56E-3	CLEA value for saturated conductivity of sandy loam, Table 4.4, SR3 <sup>(3)</sup>
Vapour permeability	m <sup>2</sup>	3.05E-12	Calculated for sandy loam using equations in Appendix 1, SR3 <sub>(3)</sub>
Capillary zone thickness	m	0.1	Professional judgement
			BUILDING
Building volume/area ratio	m	4.8	Table 3.3, SR3 <sup>(3)</sup>
Foundation area	m <sup>2</sup>	28	,
Foundation perimeter	m	22	Calculated assuming building measures 7m x 4m to give 28m <sup>2</sup> foundation area
Building air exchange rate	d <sup>-1</sup>	12	
Depth to bottom of foundation slab	m	0.15	Table 3.3, SR3 <sup>(3)</sup>
Foundation thickness	m	0.15	
Foundation crack fraction	-	0.0151	Calculated from floor crack area of 423 cm <sup>2</sup> and building footprint of 28m <sup>2</sup> in Table 4.21, SR3 <sup>(3)</sup>
Volumetric water content of cracks	-	0.33	Assumed equal to underlying soil type in assumption that cracks become filled with soil over time. Parameters for sandy loam
Volumetric air content of cracks	-	0.2	from Table 4.4, SR3 <sup>(3)</sup>
Indoor/outdoor differential pressure	Pa	3.1	From Table 3.3, SR3 <sup>(3)</sup>

### Table 5





	N <sub>O</sub>	GrAC	SAC Appropri	ate to Pathway St	OM 1% (mg/kg)	Soil Saturation	SAC Appropri	ate to Pathway SO	M 2.5% (mg/kg)	Soil Saturation	SAC Approp	riate to Pathway S	OM 6% (mg/kg)	Soil Saturation
Compound	es	(mg/l)	Oral	Inhalation	Combined	Limit (mg/kg)	Oral	Inhalation	Combined	Limit (mg/kg)	Oral	Inhalation	Combined	Limit (mg/kg)
Metals														
Arsenic	(c)	-	3.50E+01	8.50E+01	-	NR	3.50E+01	8.50E+01	-	NR	3.50E+01	8.50E+01	-	NR
Cadmium		-	1.21E+02	1.85E+02	8.49E+01	NR	1.21E+02	1.85E+02	8.49E+01	NR	1.21E+02	1.85E+02	8.49E+01	NR
Chromium (III) -oxide		-	1.98E+04	3.55E+03	3.01E+03	NR	1.98E+04	3.55E+03	3.01E+03	NR	1.98E+04	3.55E+03	3.01E+03	NR
Chromium (VI) - hexavalent		-	8.40E+01	4.25E+00	4.12E+00	NR	8.40E+01	4.25E+00	4.12E+00	NR	8.40E+01	4.25E+00	4.12E+00	NR
Copper		-	1.08E+04	1.04E+04	6.20E+03	NR	1.08E+04	1.04E+04	6.20E+03	NR	1.08E+04	1.04E+04	6.20E+03	NR
Lead	(a)	-	4.50E+02	-	-	NR	4.50E+02	-	-		4.50E+02	-	-	NR
Elemental Mercury (Hg <sup>0</sup> )	(d)	9.40E-03	-	1.70E-01	-	4.31E+00	-	4.24E-01	-	1.07E+01	-	1.02E+00	-	2.58E+01
Inorganic Mercury (Hg <sup>2+</sup> )		-	2.62E+02	2.55E+03	2.38E+02	NR	2.62E+02	2.55E+03	2.38E+02	NR	2.62E+02	2.55E+03	2.38E+02	NR
Methyl Mercury (Hg <sup>4+</sup> )		2.00E+01	1.80E+01	1.59E+01	8.43E+00	7.33E+01	1.80E+01	1.59E+01	1.13E+01	1.42E+02	1.80E+01	6.53E+01	1.41E+01	3.04E+02
Nickel	(d)	-	7.86E+02	1.27E+02	-	NR	7.86E+02	1.27E+02	-	NR	7.86E+02	1.27E+02	-	NR
Selenium	(c)		5.95E+02	-	-	NR	5.95E+02	-	-	NR	5.95E+02	-	-	NR
Zinc	(c)	-	4.05E+04	2.55E+07	-	NR	4.05E+04	2.55E+07	-	NR	4.05E+04	2.55E+07	-	NR
Cyanide		-	7.69E+02	1.15E+02	1.06E+02	NR	7.69E+02	1.15E+02	1.06E+02	NR	7.69E+02	1.15E+02	1.06E+02	NR
Volatile Organic Compounds														
Benzene		7.00E+00	2.58E+01	2.69E-01	2.66E-01	1.22E+03	2.58E+01	4.99E-01	4.90E-01	2.26E+03	2.58E+01	1.04E+00	9.98E-01	4.71E+03
Toluene		1.90E+03	1.98E+04	6.26E+02	6.07E+02	8.69E+02	1.98E+04	1.38E+03	1.29E+03	1.92E+03	1.98E+04	3.14E+03	2.71E+03	4.36E+03
Ethylbenzene		2.60E+02	8.88E+03	1.70E+02	1.67E+02	5.18E+02	8.88E+03	3.98E+02	3.81E+02	1.22E+03	8.88E+03	9.32E+02	8.43E+02	2.84E+03
Xylene - m		8.40E+01	1.60E+04	5.56E+01	5.54E+01	6.25E+02	1.60E+04	1.31E+02	1.30E+02	1.47E+03	1.60E+04	3.07E+02	3.02E+02	3.46E+03
Xylene - o		1.00E+02	1.60E+04	5.98E+01	5.95E+01	4.78E+02	1.60E+04	1.40E+02	1.39E+02	1.12E+03	1.60E+04	3.27E+02	3.21E+02	2.62E+03
Xylene - p		8.70E+01	1.60E+04	5.34E+01	5.33E+01	5.76E+02	1.60E+04	1.26E+02	1.25E+02	1.35E+03	1.60E+04	2.94E+02	2.88E+02	3.17E+03
Total xylene		8.40E+01	1.60E+04	5.56E+01	5.54E+01	6.25E+02	1.60E+04	1.31E+02	1.30E+02	1.47E+03	1.60E+04	3.07E+02	3.02E+02	3.46E+03
Methyl tertiary butyl ether (MTBE)		2.20E+03	4.45E+02	1.84E+02	1.61E+02	1.66E+04	4.45E+02	2.40E+02	2.00E+02	2.16E+04	4.45E+02	3.70E+02	2.68E+02	3.34E+04
Trichloroethene		1.80E+00	4.63E+02	1.10E-01	1.10E-01	1.54E+03	4.63E+02	2.30E-01	2.30E-01	3.22E+03	4.63E+02	5.11E-01	5.11E-01	7.14E+03
Tetrachloroethene		3.60E+00	1.20E+03	1.03E+00	1.03E+00	4.24E+02	1.20E+03	2.30E+00	2.30E+00	9.51E+02	1.20E+03	5.28E+00	5.26E+00	2.18E+03
1,1,1-Trichloroethane		2.60E+01	5.34E+04	6.33E+00	6.33E+00	1.43E+03	5.34E+04	1.29E+01	1.29E+01	2.92E+03	5.34E+04	2.84E+01	2.84E+01	6.39E+03
1,1,1,2-Tetrachloroethane		1.40E+01	5.07E+02	1.08E+00	1.08E+00	2.60E+03	5.07E+02	2.50E+00	2.49E+00	6.02E+03	5.07E+02	5.83E+00	5.76E+00	1.40E+04
1,1,2,2-Tetrachloroethane		1.40E+01	5.07E+02	2.76E+00	2.74E+00	2.67E+03	5.07E+02	5.65E+00	5.58E+00	5.46E+03	5.07E+02	1.24E+01	1.21E+01	1.20E+04
Carbon tetrachloride		5.50E-02	1.25E+02	1.81E-02	1.81E-02	1.52E+03	1.25E+02	3.97E-02	3.96E-02	3.32E+03	1.25E+02	8.99E-02	8.99E-02	7.54E+03
1,2-Dichloroethane		3.00E-01	1.07E+01	6.46E-03	6.46E-03	3.41E+03	1.07E+01	9.32E-03	9.31E-03	4.91E+03	1.07E+01	1.60E-02	1.60E-02	8.43E+03
Vinyl chloride		1.90E-02	1.25E+00	5.43E-04	5.43E-04	1.36E+03	1.25E+00	7.02E-04	7.02E-04	1.76E+03	1.25E+00	1.07E-03	1.07E-03	2.69E+03
1,2,4-Trimethylbenzene		7.50E-02	-	4.08E-01	-	5.57E+02	-	9.91E-01	-	1.36E+03	ı	2.33E+00		3.25E+03
1,3,5-Trimethylbenzene		4.70E-02	1.28E+03	4.60E-01	4.60E-01	9.47E+01	1.28E+03	1.10E+00	1.10E+00	2.26E+02	1.28E+03	2.59E+00	2.58E+00	5.33E+02
Semi-Volatile Organic Compounds														
Acenaphthene		3.20E+00	4.85E+03	3.46E+03	2.02E+03	5.70E+01	4.85E+03	8.54E+03	3.09E+03	1.41E+02	4.85E+03	2.30E+04	3.91E+03	3.36E+02
Acenaphthylene		4.20E+00	4.85E+03	3.27E+03	1.95E+03	8.61E+01	4.85E+03	8.03E+03	3.02E+03	2.12E+02	4.85E+03	1.91E+04	3.87E+03	5.06E+02
Anthracene		2.10E-02	2.43E+04	1.08E+05	1.98E+04	1.17E+00	2.43E+04	2.65E+05	2.22E+04	2.91E+00	2.43E+04	6.15E+05	2.33E+04	6.96E+00
Benzo(a)anthracene		3.80E-03	1.12E+01	5.55E+00	3.71E+00	1.71E+00	1.12E+01	9.83E+00	5.23E+00	4.28E+00	1.12E+01	1.41E+01	6.22E+00	1.03E+01
Benzo(b)fluoranthene		2.00E-03	1.15E+00	1.79E+01	6.99E+00	1.22E+00	1.15E+01	1.97E+01	7.25E+00	3.04E+00	1.15E+01	2.05E+01	7.36E+00	7.29E+00
Benzo(g,h,i)perylene		2:60E-04	7.35E+01	1.27E+02	4.66E+01	1.54E-02	7.35E+01	1.32E+02	4.72E+01	3.85E-02	7.35E+01	1.34E+02	4.75E+01	9.23E-02
Benzo(k)fluoranthene		8.00E-04	1.62E+01	2.66E+01	1.01E+01	6.87E-01	1.62E+01	2.83E+01	1.03E+01	1.72E+00	1.62E+01	2.91E+01	1.04E+01	4.12E+00
Chrysene		2.00E-03	1.62E+01	1.95E+01	8.84E+00	4.40E-01	1.62E+01	2.45E+01	9.74E+00	1.10E+00	1.62E+01	2.72E+01	1.01E+01	2.64E+00
Dibenzo(a,h)anthracene		6.00E-04	1.46E+00	2.13E+00	8.65E-01	3.93E-03	1.46E+00	2.42E+00	9.09E-01	9.82E-03	1.46E+00	2.56E+00	9.28E-01	2.36E-02
Fluoranthene		2:30E-01	1.01E+03	2.69E+04	9.72E+02	1.89E+01	1.01E+03	6.23E+04	9.93E+02	4.73E+01	1.01E+03	1.28E+05	1.00E+03	1.13E+02
Fluorene		1.90E+00	3.23E+03	4.35E+03	1.85E+03	3.09E+01	3.23E+03	1.07E+04	2.48E+03	7.65E+01	3.23E+03	2.54E+04	2.87E+03	1.83E+02
Indeno(1,2,3-cd)pyrene		2.00E-04	6.95E+00	1.04E+01	4.17E+00	6.13E-02	6.95E+00	1.17E+01	4.35E+00	1.53E-01	6.95E+00	1.22E+01	4.43E+00	3.68E-01
Phenanthrene		5.30E-01	1.00E+03	5.04E+03	8.37E+02	3.60E+01	1.00E+03	1.23E+04	9.28E+02	8.96E+01	1.00E+03	2.86E+04	9.70E+02	2.14E+02
Pyrene		1.30E-01	2.42E+03	6.18E+04	2.33E+03	2.20E+00	2.42E+03	1.44E+05	2.38E+03	5.49E+00	2.42E+03	2.97E+05	2.40E+03	1.32E+01
Benzo(a)pyrene		3.80E-03	1.62E+00	2.62E+00	1.00E+00	9.11E-01	1.62E+00	2.81E+00	1.03E+00	2.28E+00	1.62E+00	2.90E+00	1.04E+00	5.46E+00
Naphthalene		1.90E+01	1.58E+03	1.64E+00	1.64E+00	7.64E+01	1.58E+03	3.93E+00	3.92E+00	1.83E+02	1.58E+03	9.27E+00	9.22E+00	4.32E+02
Phenol		-	9.17E+04	3.11E+02	3.10E+02	4.16E+04	9.17E+04	4.20E+02	4.18E+02	8.15E+04	9.17E+04	5.21E+02	5.19E+02	1.74E+05
	-													

Table 5 RSK GAC\_2010\_03\_Rev01

#### Table 5

Human Health Generic Assessment Criteria by Pathway for Residential Scenario Without Gardens



	٥N	GrAC	SAC Appropris	ate to Pathway SC	OM 1% (mg/kg)	Soil Saturation	SAC Appropri	iate to Pathway SOI	VI 2.5% (mg/kg)	Soil Saturation	SAC Appropri	iate to Pathway S	OM 6% (mg/kg)	Soil Saturation
Compound	tes	(mg/l)	Oral	Inhalation	Combined	Limit (mg/kg)	Oral	Inhalation	Combined	Limit (mg/kg)	Oral	Inhalation	Combined	Limit (mg/kg)
Total Petroleum Hydrocarbons														
Aliphatic hydrocarbons EC <sub>5</sub> -EC <sub>6</sub>		1.00E+01	2.23E+05	2.98E+01	2.98E+01	3.04E+02	2.23E+05	5.47E+01	5.47E+01	5.58E+02	2,23E+05	1.13E+02	1.13E+02	1.15E+03
Aliphatic hydrocarbons >EC <sub>6</sub> -EC <sub>8</sub>		5.40E+00	2.23E+05	7.27E+01	7.27E+01	1.44E+02	2.23E+05	1.62E+02	1.62E+02	3.22E+02	2.23E+05	3.72E+02	3.71E+02	7.36E+02
Aliphatic hydrocarbons >EC <sub>8</sub> -EC <sub>10</sub>		2.30E-01	4.45E+03	1.89E+01	1.88E+01	7.77E+01	4.45E+03	4.60E+01	4.59E+01	1.90E+02	4.45E+03	1.09E+02	1.09E+02	4.51E+02
Aliphatic hydrocarbons >EC <sub>10</sub> -EC <sub>12</sub>		3.00E-02	4.45E+03	9.34E+01	9.29E+01	4.75E+01	4.45E+03	2.32E+02	2.29E+02	1.18E+02	4.45E+03	5.57E+02	5.38E+02	2.83E+02
Aliphatic hydrocarbons >EC <sub>12</sub> -EC <sub>16</sub>		8.00E-04	4.45E+03	7.82E+02	7.45E+02	2.37E+01	4.45E+03	1.95E+03	1.69E+03	5.91E+01	4.45E+03	4.68E+03	3.04E+03	1.42E+02
Aliphatic hydrocarbons >EC <sub>16</sub> -EC <sub>35</sub>	(c)	-	4.53E+04	-	-	8.48E+00	6.41E+04	-	-	2.12E+01	7.66E+04	-	-	5.09E+01
Aliphatic hydrocarbons >EC35-EC44	(c)	-	4.53E+04	-	-	8.48E+00	6.41E+04	-	-	2.12E+01	7.66E+04	-	-	5.09E+01
Aromatic hydrocarbons >EC <sub>5</sub> -EC <sub>7</sub>		-	1.98E+04	2.66E+02	2.63E+02	1.22E+03	1.98E+04	4.95E+02	4.83E+02	2.26E+03	1.98E+04	1.03E+03	9.78E+02	4.71E+03
Aromatic hydrocarbons >EC <sub>7</sub> -EC <sub>8</sub>		-	1.98E+04	6.26E+02	6.07E+02	8.69E+02	1.98E+04	1.38E+03	1.29E+03	1.92E+03	1.98E+04	3.14E+03	2.71E+03	4.36E+03
Aromatic hydrocarbons >EC <sub>8</sub> -EC <sub>9</sub> (styre	ene)	7.40E+00	5.34E+03	2.65E+02	2.61E+02	6.20E+02	5.34E+03	6.47E+02	6.27E+02	1.52E+03	5.34E+03	1.54E+03	1.41E+03	3.61E+03
Aromatic hydrocarbons >EC <sub>9</sub> -EC <sub>10</sub>		7.40E+00	1.78E+03	3.33E+01	3.32E+01	6.13E+02	1.78E+03	8.16E+01	8.07E+01	1.50E+03	1.78E+03	1.94E+02	1.89E+02	3.58E+03
Aromatic hydrocarbons >EC <sub>10</sub> -EC <sub>12</sub>		2.50E+01	1.78E+03	1.82E+02	1.77E+02	3.64E+02	1.78E+03	4.48E+02	4.17E+02	8.99E+02	1.78E+03	1.07E+03	8.66E+02	2.15E+03
Aromatic hydrocarbons >EC <sub>12</sub> -EC <sub>16</sub>		5.80E+00	1.78E+03	2.00E+03	1.25E+03	1.69E+02	1.78E+03	4.96E+03	1.59E+03	4.19E+02	1.78E+03	1.18E+04	1.71E+03	1.00E+03
Aromatic hydrocarbons >EC <sub>16</sub> -EC <sub>21</sub>	(c)	-	1.29E+03	-	-	5.37E+01	1.31E+03	-	-	1.34E+02	1.32E+03	-	-	3.21E+02
Aromatic hydrocarbons >EC <sub>21</sub> -EC <sub>35</sub>	(c)	-	1.33E+03	-	-	4.83E+00	1.33E+03	-	-	1.21E+01	1.33E+03	-	-	2.90E+01
Aromatic hydrocarbons >EC <sub>35</sub> -EC <sub>44</sub>	(c)	-	1.33E+03	-	-	4.83E+00	1.33E+03	-	-	1.21E+01	1.33E+03	-	-	2.90E+01

#### Notes:

-' Generic assessment criteria not calculated owing to low volatility of substance and therefore no pathway, or an absence of toxicological data.

NR - the compound is not volatile and therefore a soil saturation limit not calculated within CLEA

EC - equivalent carbon. GrAC - groundwater assessment criteria. SAC - soil assessment criteria.

The CLEA model output is colour coded depending upon whether the soil saturation limit has been exceeded.



Calculated SAC exceeds soil saturation limit and may significantly effect the interpretation of any exceedances since the contribution of the indoor and outdoor vapour pathway to total exposure is

>10%. This shading has also been used for the RBCA output where the theoretical solubility limit has been exceeded. The SAC has been set as the model calculated SAC with the saturation limits shown in brackets. Calculated SAC exceeds soil saturation limit but will not effect the SAC significantly since the contribution of the indoor and outdoor vapour pathway to total exposure is <10%. Calculated SAC does not exceed the soil saturation limit.

For consistency where the theoretical solubility limit within RBCA has been exceeded in production of the GrAC, these cellls have also been hatched red.

The SAC for organic compounds are dependant upon soil organic matter (SOM) (%) content. To obtain SOM from total organic carbon (TOC) (%) divide by 0.58. 1% SOM is 0.58% TOC. DL Rowell Soil Science: Methods and Applications, Longmans, 1994. SAC for TPH fractions, polycyclic aromatic hydrocarbons, MTBE, BTEX and trimethylbenzene compounds were produced using an attenuation factor for the indoor air inhalation pathway of 10 to reduce conservatism associated with the vapour inhalation pathway, section 10.1.1, SR3

- (a) GAC taken as former Soil Guideline Value owing to uncertainty regarding toxicological approach to be adopted by the Environment Agency.
- (b) GAC taken from the Environment Agency SGV reports published 2009.
- (c) SAC for selenium, aliphatic and aromatic hydrocarbons >EC16 do not include inhalation pathway owing to absence of toxicity data. SAC for arsenic is only based on oral contribution (rather than combined) owing to the relative small contribution from inhalation in accordance with the SGV report.
- (d) SAC for elemental mercury, chromium VI and nickel are based on the inhalation pathway only owing to an absence of toxicity for elemental mercury, in accordance with the SGV report for nickel and LQM report for chromium VI.



Table 6 Human Health Generic Assessment Criteria for Residential Without Gardens

Compound	GrAC for Groundwater (mg/l)	SAC for Soil SOM 1% (mg/kg)	SAC for Soil SOM 2.5% (mg/kg)	SAC for Soil SOM 6% (mg/kg)
Metals Arsenic		35	35	35
Cadmium	-	85	85	85
Chromium (III) - oxide		3,000	3,000	3,000
Chromium (VI) - hexavalent	-	4.3	4.3	4.3
Copper	-	6,200	6,200	6,200
Lead	-	450	450	450
Elemental Mercury (Hg0)	0.0094	0.17 240	0.42 240	1.0 240
Inorganic Mercury (Hg2+) Methyl Mercury (Hg4+)	20	8.4	11	14
Nickel	-	130	130	130
Selenium	-	600	600	600
Zinc	-	41,000	41,000	41,000
Cyanide	-	110	110	110
Valatila Organia Campayada				
Volatile Organic Compounds Benzene	7	0.27	0.49	1.0
Toluene	1,900	610	1,289	2,700
Ethylbenzene	260	170	381	840
Xylene - m	84	55	130	300
Xylene - o	100	60	139	320
Xylene - p	87	53	125	290
Total xylene	84	55	130	300
Methyl tertiary butyl ether (MTBE) Trichloroethene	2,200 1.8	160 0.11	199.55 0.2	270
Tetrachloroethene	3.6	1.0	2.3	0.51 5.3
1,1,1-Trichloroethane	26	6.3	12.9	28
1,1,1,2-Tetrachloroethane	14	1.1	2.5	5.8
1,1,2,2-Tetrachloroethane	14	2.7	5.58	12
Carbon tetrachloride	0.055	0.02	0.040	0.09
1,2-Dichloroethane	0.30	0.006	0.0093	0.02
Vinyl chloride	0.019	0.0005	0.0007	0.001
1,2,4-Trimethylbenzene	0.075	0.4	0.99	2.3
1,3,5-Trimethylbenzene	0.047	0.5	1.10	2.6
Semi-Volatile Organic Compounds				
Acenaphthene	3.2	2,000 (57)	3,100 (141)	3,900 (340)
Acenaphthylene Anthracene	4.2 0.021	2,000 (86) 20,000 (1.2)	3,000 (212) 22,000	3.900 (510) 23,000
Benzo(a)anthracene	0.004	3.7	5.2	6.2
		7.0	7.3	7.4
Benzo(b)fluoranthene	0.002			
Benzo(b)fluoranthene Benzo(g,h,i)perylene	0.002 0.0003	47	47	48
	0.0003 0.0008	47 10	10	10
Benzo(g,h,i)perylene Benzo(k)fluoranthene Chrysene	0.0003 0.0008 0.002	47 10 8.8	10 9.7	10 10
Benzo(g,h,i)perylene Benzo(k)fluoranthene Chrysene Dibenzo(a,h)anthracene	0.0003 0.0008 0.002 0.0006	47 10 8.8 0.87	10 9.7 0.91	10 10 0.93
Benzo(g,h,i)perylene Benzo(k)fluoranthene Chrysene Dibenzo(a,h)anthracene Fluoranthene	0.0003 0.0008 0.002 0.0006 0.23	47 10 8.8 0.87 970	10 9.7 0.91 993	10 10 0.93 1,000
Benzo(g,h,i)perylene Benzo(k)fluoranthene Chrysene Dibenzo(a,h)anthracene Fluoranthene Fluorene	0.0003 0.0008 0.002 0.0006 0.23 1.9	47 10 8.8 0.87 970 1,900 (31)	10 9.7 0.91 993 2,500 (77)	10 10 0.93 1,000 2,900 (180)
Benzo(g,h,i)perylene Benzo(k)fluoranthene Chrysene Dibenzo(a,h)anthracene Fluoranthene Fluorene Indeno(1,2,3-cd)pyrene	0.0003 0.0008 0.002 0.0006 0.23 1.9	47 10 8.8 0.87 970 1.900 (31) 4.2	10 9.7 0.91 993 2,500 (77) 4.4	10 10 0.93 1,000 2,900 (180) 4.4
Benzo(g,h,i)perylene Benzo(k)fluoranthene Chrysene Dibenzo(a,h)anthracene Fluoranthene Fluorene	0.0003 0.0008 0.002 0.0006 0.23 1.9 0.0002	47 10 8.8 0.87 970 1.900 (31) 4.2 840 (36)	10 9.7 0.91 993 2,500 (77)	10 10 0.93 1,000 2,900 (180)
Benzo(g,h,i)perylene Benzo(k)fluoranthene Chrysene Dibenzo(a,h)anthracene Fluoranthene Fluorene Indeno(1,2,3-cd)pyrene Phenanthrene	0.0003 0.0008 0.002 0.0006 0.23 1.9	47 10 8.8 0.87 970 1.900 (31) 4.2	10 9.7 0.91 993 2,500 (77) 4.4 930	10 10 0.93 1,000 2,900 (180) 4.4 970
Benzo(g,h,i)perylene Benzo(k)fluoranthene Chrysene Dibenzo(a,h)anthracene Fluoranthene Fluorene Indeno(1,2,3-cd)pyrene Phenanthrene Pyrene Benzo(a)pyrene Naphthalene	0.0003 0.0008 0.002 0.0006 0.23 1.9 0.0002 0.53 0.13	47 10 8.8 0.87 970 1,900 (31) 4.2 840 (36) 2,300 1.0	10 9.7 0.91 993 2,500 (77) 4.4 930 2,400 1.0 3.9	10 10 0.93 1,000 2,900 (180) 4.4 970 2,400 1.0 9.2
Benzo(g,h,i)perylene Benzo(k)fluoranthene Chrysene Dibenzo(a,h)anthracene Fluoranthene Fluorene Indeno(1,2,3-cd)pyrene Phenanthrene Pyrene Benzo(a)pyrene	0.0003 0.0008 0.002 0.0006 0.23 1.9 0.0002 0.53 0.13	47 10 8.8 0.87 970 1,900 (31) 4.2 840 (36) 2,300 1.0	10 9.7 0.91 993 2,500 (77) 4.4 930 2,400 1.0	10 10 0.93 1,000 2,900 (180) 4.4 970 2,400
Benzo(g,h,i)perylene Benzo(k)fluoranthene Chrysene Dibenzo(a,h)anthracene Fluoranthene Fluorene Indeno(1,2,3-cd)pyrene Phenanthrene Pyrene Benzo(a)pyrene Naphthalene Phenol	0.0003 0.0008 0.0002 0.0006 0.23 1.9 0.0002 0.53 0.13 0.004	47 10 8.8 0.87 970 1,900 (31) 4.2 840 (36) 2,300 1.0	10 9.7 0.91 993 2,500 (77) 4.4 930 2,400 1.0 3.9	10 10 0.93 1,000 2,900 (180) 4.4 970 2,400 1.0 9.2
Benzo(g,h,i)perylene Benzo(k)fluoranthene Chrysene Dibenzo(a,h)anthracene Fluoranthene Fluorene Indeno(1,2,3-cd)pyrene Phenanthrene Pyrene Benzo(a)pyrene Naphthalene Phenol Total Petroleum Hydrocarbons	0.0003 0.0008 0.0002 0.0006 0.23 1.9 0.0002 0.53 0.13 0.004	47 10 8.8 0.87 970 1.900 (31) 4.2 840 (36) 2,300 1.0 1.6 310	10 9.7 0.91 993 2,500 (77) 4.4 930 2,400 1.0 3.9 420	10 10 0.93 1,000 2,900 (180) 4.4 970 2,400 1.0 9.2
Benzo(g,h,i)perylene Benzo(k)fluoranthene Chrysene Dibenzo(a,h)anthracene Fluoranthene Fluorene Indeno(1,2,3-cd)pyrene Phenanthrene Pyrene Benzo(a)pyrene Naphthalene Phenol  Total Petroleum Hydrocarbons Aliphatic hydrocarbons EC <sub>5</sub> -EC <sub>6</sub>	0.0003 0.0008 0.002 0.0006 0.23 1.9 0.0002 0.53 0.13 0.004 19	47 10 8.8 0.87 970 1.906 (31) 4.2 840 (36) 2,300 1.0 1.6 310	10 9.7 0.91 993 2,500 (77) 4.4 930 2,400 1.0 3.9 420	10 10 0.93 1,000 2,900 (180) 4.4 970 2,400 1.0 9.2 520
Benzo(g,h,i)perylene Benzo(k)fluoranthene Chrysene Dibenzo(a,h)anthracene Fluoranthene Fluorene Indeno(1,2,3-cd)pyrene Phenanthrene Pyrene Benzo(a)pyrene Naphthalene Phenol  Total Petroleum Hydrocarbons Aliphatic hydrocarbons >EC <sub>6</sub> -EC <sub>6</sub> Aliphatic hydrocarbons >EC <sub>6</sub> -EC <sub>6</sub>	0.0003 0.0008 0.002 0.0006 0.23 1.9 0.0002 0.53 0.13 0.004 19	47 10 8.8 0.87 970 1,900 (31) 4.2 840 (36) 2,300 1.0 1.6 310	10 9.7 0.91 993 2500 (77) 4.4 930 2,400 1.0 3.9 420	10 10 0.93 1,000 2,900 (180) 4.4 970 2,400 1.0 9.2 520
Benzo(g,h,i)perylene Benzo(k)fluoranthene Chrysene Dibenzo(a,h)anthracene Fluoranthene Fluorene Indeno(1,2,3-cd)pyrene Phenanthrene Pyrene Benzo(a)pyrene Naphthalene Phenol  Total Petroleum Hydrocarbons Aliphatic hydrocarbons >EC <sub>5</sub> -EC <sub>6</sub> Aliphatic hydrocarbons >EC <sub>6</sub> -EC <sub>8</sub> Aliphatic hydrocarbons >EC <sub>6</sub> -EC <sub>8</sub>	0.0003 0.0008 0.002 0.0006 0.23 1.9 0.0002 0.53 0.13 0.004 19 - 10 5.4 0.23	47 10 8.8 0.87 970 1.900 (31) 4.2 840 (36) 2,300 1.0 1.6 310  30 73	10 9.7 0.91 993 2,500 (77) 4.4 930 2,400 1.0 3.9 420 55 160 46	10 10 0.93 1,000 2,900 (180) 4.4 970 2,400 1.0 9.2 520
Benzo(g,h,i)perylene Benzo(k)fluoranthene Chrysene Dibenzo(a,h)anthracene Fluoranthene Fluorene Indeno(1,2,3-cd)pyrene Phenanthrene Pyrene Benzo(a)pyrene Naphthalene Phenol  Total Petroleum Hydrocarbons Aliphatic hydrocarbons >EC <sub>6</sub> -EC <sub>6</sub> Aliphatic hydrocarbons >EC <sub>6</sub> -EC <sub>6</sub>	0.0003 0.0008 0.002 0.0006 0.23 1.9 0.0002 0.53 0.13 0.004 19	47 10 8.8 0.87 970 1,900 (31) 4.2 840 (36) 2,300 1.0 1.6 310	10 9.7 0.91 993 2500 (77) 4.4 930 2,400 1.0 3.9 420	10 10 0.93 1,000 2,900 (180) 4.4 970 2,400 1.0 9.2 520
Benzo(g,h,i)perylene Benzo(k)fluoranthene Chrysene Dibenzo(a,h)anthracene Fluoranthene Fluorene Indeno(1,2,3-cd)pyrene Phenanthrene Pyrene Benzo(a)pyrene Naphthalene Phenol  Total Petroleum Hydrocarbons Aliphatic hydrocarbons >EC <sub>5</sub> -EC <sub>6</sub> Aliphatic hydrocarbons >EC <sub>6</sub> -EC <sub>8</sub> Aliphatic hydrocarbons >EC <sub>6</sub> -EC <sub>8</sub>	0.0003 0.0008 0.002 0.0006 0.23 1.9 0.0002 0.53 0.13 0.004 19 - 10 5.4 0.23	47 10 8.8 0.87 970 1.900 (31) 4.2 840 (36) 2,300 1.0 1.6 310  30 73	10 9.7 0.91 993 2,500 (77) 4.4 930 2,400 1.0 3.9 420 55 160 46	10 10 0.93 1,000 2,900 (180) 4.4 970 2,400 1.0 9.2 520
Benzo(g,h,i)perylene Benzo(k)fluoranthene Chrysene Dibenzo(a,h)anthracene Fluoranthene Fluorene Indeno(1,2,3-cd)pyrene Phenanthrene Pyrene Benzo(a)pyrene Naphthalene Phenol  Total Petroleum Hydrocarbons Aliphatic hydrocarbons >EC <sub>5</sub> -EC <sub>6</sub> Aliphatic hydrocarbons >EC <sub>6</sub> -EC <sub>8</sub> Aliphatic hydrocarbons >EC <sub>6</sub> -EC <sub>10</sub> Aliphatic hydrocarbons >EC <sub>10</sub> -EC <sub>12</sub>	0.0003 0.0008 0.0002 0.0006 0.23 1.9 0.0002 0.53 0.13 0.004 19 - 10 5.4 0.23	47 10 8.8 0.87 970 1.900 (31) 4.2 840 (36) 2,300 1.0 1.6 310 30 73 19 93 (48)	10 9.7 0.91 993 2,500 (77) 4.4 930 2,400 1.0 3.9 420 55 160 46	10 10 0.93 1,000 2,900 (180) 4.4 970 2,400 1.0 9.2 520  110 370 110 540 (280)
Benzo(g,h,i)perylene Benzo(k)fluoranthene Chrysene Dibenzo(a,h)anthracene Fluoranthene Fluoranthene Fluorene Indeno(1,2,3-cd)pyrene Phenanthrene Pyrene Benzo(a)pyrene Naphthalene Phenol  Total Petroleum Hydrocarbons Aliphatic hydrocarbons EC <sub>8</sub> -EC <sub>6</sub> Aliphatic hydrocarbons >EC <sub>8</sub> -EC <sub>8</sub> Aliphatic hydrocarbons >EC <sub>8</sub> -EC <sub>10</sub> Aliphatic hydrocarbons >EC <sub>10</sub> -EC <sub>12</sub> Aliphatic hydrocarbons >EC <sub>10</sub> -EC <sub>12</sub> Aliphatic hydrocarbons >EC <sub>12</sub> -EC <sub>16</sub>	0.0003 0.0008 0.0002 0.0006 0.23 1.9 0.0002 0.53 0.13 0.004 19 - 10 5.4 0.23 0.03 0.03	47 10 8.8 0.87 970 1.900(31) 4.2 840(36) 2,300 1.0 1.6 310  30 73 19 33 (48) 746 (24) 45,000	10 9.7 0.91 993 2,500 (77) 4.4 930 2,400 1.0 3.9 420 55 160 46 230 (118) 1,700 (59) 84,900 (21)	10 10 0.93 1,000 2,900 (180) 4.4 970 2,400 1.0 9.2 520  110 370 110 540 (280) 3,000 (140) 77,000
Benzo(g,h,i)perylene Benzo(k)fluoranthene Chrysene Dibenzo(a,h)anthracene Fluoranthene Fluorene Indeno(1,2,3-cd)pyrene Phenanthrene Pyrene Benzo(a)pyrene Naphthalene Phenol  Total Petroleum Hydrocarbons Aliphatic hydrocarbons EC <sub>8</sub> -EC <sub>6</sub> Aliphatic hydrocarbons >EC <sub>8</sub> -EC <sub>8</sub> Aliphatic hydrocarbons >EC <sub>8</sub> -EC <sub>10</sub> Aliphatic hydrocarbons >EC <sub>10</sub> -EC <sub>12</sub> Aliphatic hydrocarbons >EC <sub>12</sub> -EC <sub>16</sub> Aliphatic hydrocarbons >EC <sub>15</sub> -EC <sub>35</sub> Aliphatic hydrocarbons >EC <sub>35</sub> -EC <sub>44</sub>	0.0003 0.0008 0.0002 0.0006 0.23 1.9 0.0002 0.53 0.13 0.004 19 - 10 5.4 0.23 0.03 0.03	47 10 8.8 0.87 970 1.900 (31) 4.2 840 (36) 2,300 1.0 1.6 310  30 73 19 93 (48) 746 (24) 45,000	10 9.7 0.91 993 2,500 (77) 4.4 930 2,400 1.0 3.9 420 55 160 46 230 (118) 1,700 (59) 64,000 (21)	10 10 0.93 1,000 2,900 (180) 4.4 970 2,400 1.0 9.2 520  110 370 110 540 (280) 3,000 (140) 77,000
Benzo(g,h,i)perylene Benzo(k)fluoranthene Chrysene Dibenzo(a,h)anthracene Fluoranthene Fluorene Indeno(1,2,3-cd)pyrene Phenanthrene Pyrene Benzo(a)pyrene Naphthalene Phenol  Total Petroleum Hydrocarbons Aliphatic hydrocarbons EC <sub>8</sub> -EC <sub>8</sub> Aliphatic hydrocarbons >EC <sub>8</sub> -EC <sub>8</sub> Aliphatic hydrocarbons >EC <sub>8</sub> -EC <sub>10</sub> Aliphatic hydrocarbons >EC <sub>10</sub> -EC <sub>12</sub> Aliphatic hydrocarbons >EC <sub>10</sub> -EC <sub>13</sub> Aliphatic hydrocarbons >EC <sub>10</sub> -EC <sub>13</sub> Aliphatic hydrocarbons >EC <sub>10</sub> -EC <sub>13</sub> Aliphatic hydrocarbons >EC <sub>10</sub> -EC <sub>35</sub> Aliphatic hydrocarbons >EC <sub>35</sub> -EC <sub>44</sub> Aromatic hydrocarbons >EC <sub>8</sub> -EC <sub>9</sub> (styrene)	0.0003 0.0008 0.0008 0.0002 0.0006 0.23 1.9 0.0002 0.53 0.13 0.004 19 10 5.4 0.23 0.03 0.03 0.008 7.4	47 10 8.8 0.87 970 1.900 (31) 4.2 840 (36) 2,300 1.0 1.6 310  30 73 19 93 (48) 746 (24) 45,000 45,000	10 9.7 0.91 993 2,500 (77) 4.4 930 2,400 1.0 3.9 420 55 160 46 230 (118) 1,700 (59) 64,000 (21) 627	10 10 0.93 1,000 2,900 (180) 4.4 970 2,400 1.0 9.2 520  110 370 110 540 (280) 3,000 (140) 77,000 77,000 1,400
Benzo(g,h,i)perylene Benzo(k)fluoranthene Chrysene Dibenzo(a,h)anthracene Fluoranthene Fluorene Indeno(1,2,3-cd)pyrene Phenanthrene Pyrene Benzo(a)pyrene Naphthalene Phenol  Total Petroleum Hydrocarbons Aliphatic hydrocarbons EC <sub>5</sub> -EC <sub>6</sub> Aliphatic hydrocarbons >EC <sub>6</sub> -EC <sub>8</sub> Aliphatic hydrocarbons >EC <sub>6</sub> -EC <sub>12</sub> Aliphatic hydrocarbons >EC <sub>10</sub> -EC <sub>13</sub> Aliphatic hydrocarbons >EC <sub>10</sub> -EC <sub>13</sub> Aliphatic hydrocarbons >EC <sub>10</sub> -EC <sub>35</sub> Aliphatic hydrocarbons >EC <sub>35</sub> -EC <sub>44</sub> Aromatic hydrocarbons >EC <sub>3</sub> -EC <sub>9</sub> (styrene) Aromatic hydrocarbons >EC <sub>8</sub> -EC <sub>9</sub> (styrene)	0.0003 0.0008 0.0008 0.0002 0.0008 0.23 1.9 0.0002 0.53 0.13 0.004 19 10 5.4 0.23 0.03 0.008 7.4 7.4	47 10 8.8 0.87 970 1.900 (31) 4.2 840 (36) 2,300 1.0 1.6 310  30 73 19 33 (48) 746 (24) 45,000 45,000 260 33	10 9.7 0.91 993 2,500 (77) 4.4 930 2,400 1.0 3.9 420 55 160 46 230 (118) 1,700 (59) 64,000 (21) 627 81	10 10 0.93 1,000 2,900 (180) 4.4 970 2,400 1.0 9.2 520  110 370 110 540 (280) 3,000 (140) 77,000 77,000 1,400 190
Benzo(g,h,i)perylene Benzo(k)fluoranthene Chrysene Dibenzo(a,h)anthracene Fluoranthene Fluorene Indeno(1,2,3-cd)pyrene Phenanthrene Pyrene Benzo(a)pyrene Naphthalene Phenol  Total Petroleum Hydrocarbons Aliphatic hydrocarbons EC <sub>8</sub> -EC <sub>8</sub> Aliphatic hydrocarbons >EC <sub>8</sub> -EC <sub>8</sub> Aliphatic hydrocarbons >EC <sub>8</sub> -EC <sub>10</sub> Aliphatic hydrocarbons >EC <sub>12</sub> -EC <sub>12</sub> Aliphatic hydrocarbons >EC <sub>12</sub> -EC <sub>12</sub> Aliphatic hydrocarbons >EC <sub>12</sub> -EC <sub>13</sub> Aliphatic hydrocarbons >EC <sub>12</sub> -EC <sub>14</sub> Aliphatic hydrocarbons >EC <sub>12</sub> -EC <sub>16</sub> Aliphatic hydrocarbons >EC <sub>12</sub> -EC <sub>16</sub> Aliphatic hydrocarbons >EC <sub>3</sub> -EC <sub>3</sub> Aliphatic hydrocarbons >EC <sub>3</sub> -EC <sub>10</sub> Aromatic hydrocarbons >EC <sub>8</sub> -EC <sub>9</sub> (styrene) Aromatic hydrocarbons >EC <sub>9</sub> -EC <sub>10</sub> Aromatic hydrocarbons >EC <sub>10</sub> -EC <sub>12</sub>	0.0003 0.0008 0.0008 0.0002 0.0006 0.23 1.9 0.0002 0.53 0.13 0.004 19 10 5.4 0.23 0.03 0.03 0.008 7.4 7.4	47 10 8.8 0.87 970 1.900 (31) 4.2 840 (36) 2,300 1.0 1.6 310  30 73 19 93 (48) 746 (24) 45,000 45,000 260 33 180	10 9.7 0.91 993 2,500 (77) 4.4 930 2,400 1.0 3.9 420  55 160 46 230 (118) 1,700 (59) 64,000 (21) 627 81 417	10 10 0.93 1,000 2,900 (180) 4.4 970 2,400 1.0 9.2 520  110 370 110 540 (280) 3,000 (140) 77,000 77,000 1,400 190 870
Benzo(g,h,i)perylene Benzo(k)fluoranthene Chrysene Dibenzo(a,h)anthracene Fluoranthene Fluorene Indeno(1,2,3-cd)pyrene Phenanthrene Pyrene Benzo(a)pyrene Naphthalene Phenol  Total Petroleum Hydrocarbons Aliphatic hydrocarbons EC <sub>5</sub> -EC <sub>6</sub> Aliphatic hydrocarbons >EC <sub>6</sub> -EC <sub>8</sub> Aliphatic hydrocarbons >EC <sub>6</sub> -EC <sub>10</sub> Aliphatic hydrocarbons >EC <sub>10</sub> -EC <sub>12</sub> Aliphatic hydrocarbons >EC <sub>10</sub> -EC <sub>12</sub> Aliphatic hydrocarbons >EC <sub>10</sub> -EC <sub>12</sub> Aliphatic hydrocarbons >EC <sub>10</sub> -EC <sub>13</sub> Aliphatic hydrocarbons >EC <sub>10</sub> -EC <sub>13</sub> Aliphatic hydrocarbons >EC <sub>10</sub> -EC <sub>35</sub> Aliphatic hydrocarbons >EC <sub>35</sub> -EC <sub>44</sub> Aromatic hydrocarbons >EC <sub>3</sub> -EC <sub>9</sub> (styrene) Aromatic hydrocarbons >EC <sub>3</sub> -EC <sub>10</sub>	0.0003 0.0008 0.0008 0.0002 0.0008 0.23 1.9 0.0002 0.53 0.13 0.004 19 10 5.4 0.23 0.03 0.008 7.4 7.4	47 10 8.8 0.87 970 1.900 (31) 4.2 840 (36) 2,300 1.0 1.6 310  30 73 19 33 (48) 746 (24) 45,000 45,000 260 33	10 9.7 0.91 993 2,500 (77) 4.4 930 2,400 1.0 3.9 420 55 160 46 230 (118) 1,700 (59) 64,000 (21) 627 81	10 10 0.93 1,000 2,900 (180) 4.4 970 2,400 1.0 9.2 520  110 370 110 540 (280) 3,000 (140) 77,000 77,000 1,400 190
Benzo(g,h,i)perylene Benzo(k)fluoranthene Chrysene Dibenzo(a,h)anthracene Fluoranthene Fluorene Indeno(1,2,3-cd)pyrene Phenanthrene Pyrene Benzo(a)pyrene Naphthalene Phenol  Total Petroleum Hydrocarbons Aliphatic hydrocarbons EC <sub>8</sub> -EC <sub>8</sub> Aliphatic hydrocarbons >EC <sub>8</sub> -EC <sub>8</sub> Aliphatic hydrocarbons >EC <sub>8</sub> -EC <sub>10</sub> Aliphatic hydrocarbons >EC <sub>12</sub> -EC <sub>12</sub> Aliphatic hydrocarbons >EC <sub>12</sub> -EC <sub>12</sub> Aliphatic hydrocarbons >EC <sub>12</sub> -EC <sub>13</sub> Aliphatic hydrocarbons >EC <sub>12</sub> -EC <sub>14</sub> Aliphatic hydrocarbons >EC <sub>12</sub> -EC <sub>16</sub> Aliphatic hydrocarbons >EC <sub>12</sub> -EC <sub>16</sub> Aliphatic hydrocarbons >EC <sub>3</sub> -EC <sub>3</sub> Aliphatic hydrocarbons >EC <sub>3</sub> -EC <sub>10</sub> Aromatic hydrocarbons >EC <sub>8</sub> -EC <sub>9</sub> (styrene) Aromatic hydrocarbons >EC <sub>9</sub> -EC <sub>10</sub> Aromatic hydrocarbons >EC <sub>10</sub> -EC <sub>12</sub>	0.0003 0.0008 0.0008 0.0002 0.0006 0.23 1.9 0.0002 0.53 0.13 0.004 19 10 5.4 0.23 0.03 0.03 0.008 7.4 7.4	47 10 8.8 0.87 970 1.900 (31) 4.2 840 (36) 2,300 1.0 1.6 310  30 73 19 93 (48) 746 (24) 45,000 45,000 260 33 180	10 9.7 0.91 993 2,500 (77) 4.4 930 2,400 1.0 3.9 420  55 160 46 230 (118) 1,700 (59) 64,000 (21) 627 81 417	10 10 0.93 1,000 2,900 (180) 4.4 970 2,400 1.0 9.2 520  110 370 110 540 (280) 3,000 (140) 77,000 77,000 1,400 190 870
Benzo(g,h,i)perylene Benzo(k)fluoranthene Chrysene Dibenzo(a,h)anthracene Fluoranthene Fluorene Indeno(1,2,3-cd)pyrene Phenanthrene Pyrene Benzo(a)pyrene Naphthalene Phenol  Total Petroleum Hydrocarbons Aliphatic hydrocarbons EC <sub>5</sub> -EC <sub>6</sub> Aliphatic hydrocarbons >EC <sub>6</sub> -EC <sub>8</sub> Aliphatic hydrocarbons >EC <sub>6</sub> -EC <sub>10</sub> Aliphatic hydrocarbons >EC <sub>10</sub> -EC <sub>12</sub> Aliphatic hydrocarbons >EC <sub>12</sub> -EC <sub>16</sub> Aliphatic hydrocarbons >EC <sub>3</sub> -EC <sub>9</sub> (styrene) Aromatic hydrocarbons >EC <sub>8</sub> -EC <sub>9</sub> (styrene) Aromatic hydrocarbons >EC <sub>9</sub> -EC <sub>12</sub> Aromatic hydrocarbons >EC <sub>12</sub> -EC <sub>12</sub>	0.0003 0.0008 0.0008 0.0002 0.0006 0.23 1.9 0.0002 0.53 0.13 0.004 19 10 5.4 0.23 0.03 0.03 0.008 7.4 7.4 7.4 25 5.8	47 10 8.8 0.87 970 1.900 (31) 4.2 840 (36) 2,300 1.0 1.6 310  30 73 19 33 (48) 746 (24) 45,000 45,000 260 33 180 1.300 (170)	10 9.7 0.91 993 2,500 (77) 4.4 930 2,400 1.0 3.9 420  55 160 46 230 (118) 1,700 (59) 64,000 (21) 627 81 417 1,600 (419)	10 10 0.93 1,000 2,900 (180) 4.4 970 2,400 1.0 9.2 520  110 370 110 540 (280) 3,000 (140) 77,000 77,000 1,400 190 870 1,700
Benzo(g,h,i)perylene Benzo(k)fluoranthene Chrysene Dibenzo(a,h)anthracene Fluoranthene Fluorene Indeno(1,2,3-cd)pyrene Phenanthrene Pyrene Benzo(a)pyrene Naphthalene Phenol  Total Petroleum Hydrocarbons Aliphatic hydrocarbons EC <sub>5</sub> -EC <sub>6</sub> Aliphatic hydrocarbons >EC <sub>6</sub> -EC <sub>8</sub> Aliphatic hydrocarbons >EC <sub>10</sub> -EC <sub>12</sub> Aromatic hydrocarbons >EC <sub>3</sub> -EC <sub>9</sub> (styrene) Aromatic hydrocarbons >EC <sub>3</sub> -EC <sub>10</sub> Aromatic hydrocarbons >EC <sub>10</sub> -EC <sub>12</sub>	0.0003 0.0008 0.0008 0.0008 0.0008 0.0008 0.23 1.9 0.0002 0.53 0.13 0.004 19 10 5.4 0.23 0.03 0.03 0.008 7.4 7.4 25 5.8	47 10 8.8 0.87 970 1.900 (31) 4.2 840 (36) 2,300 1.0 1.6 310  30 73 19 33 (48) 746 (24) 45,000 45,000 260 33 180 1,300 (170) 1,300	10 9.7 0.91 993 2,500 (77) 4.4 930 2,400 1.0 3.9 420 55 160 46 230 (118) 1,700 (59) 64,000 (21) 627 81 417 1,600 (419) 1,300	10 10 0.93 1,000 2,900 (180) 4.4 970 2,400 1.0 9.2 520  110 370 110 540 (280) 3,000 (140) 77,000 77,000 1,400 190 870 1,700 1,300

#### Notes:

The SAC has been set as the model calculated SAC with the saturation limit shown in brackets.

For consistency where the GrAC exceeds the solubility limit, GrSV has been set at the solubility limit. These are highly conservative since concentrations of the chemical are very unlikely to be at sufficient concentration to result in an exceedance of the health criteria value at the point of exposure (i.e. indoor air) provided free-phase product is absent.

Table 6 RSK GAC\_2010\_03\_Rev01

L'Generic assessment criteria not calculated owing to low volatility of substance and therefore no pathway, or an absence of toxicological data.

EC - equivalent carbon. GrAC - groundwater assessment criteria. SAC - soil assessment criteria.

The SAC for organic compounds are dependent on Soil Organic Matter (SOM) (%) content. To obtain SOM from total organic carbon (TOC) (%) divide by 0.58. 1% SOM is 0.58% TOC. DL Rowell Soil Science: Methods and Applications, Longmans, 1994.

SAC for TPH fractions, polycyclic aromatic hydrocarbons, MTBE, BTEX and trimethylbenzene compounds were produced using an attenuation factor for the indoor air inhalation pathway of 10 to reduce conservatism associated with the vapour inhalation pathway, section 10.1.1, SR3.



### **REFERENCES**

- 1) Environment Agency, 31 March 2009, May 2009 and September 2009. Science Report SC050021 / benzene SGV, toluene SGV, ethylbenzene SGV, xylene SGV, mercury SGV, selenium SGV, nickel SGV, arsenic SGV, cadmium SGV, phenol SGV, dioxins, furans and dioxinlike PCBs SGVs. Supplementary information for the derivation of SGV for: benzene, toluene, ethylbenzene, xylene, mercury, selenium, nickel, arsenic, cadmium, phenol, dioxins, furans and dioxin-like PCBs. Contaminants in soil: updated collation of toxicological data and intake values for humans: benzene, toluene, ethylbenzene, xylene, mercury, selenium, nickel, arsenic, cadmium, phenol, dioxins, furans and dioxin-like PCBs.
- 2) Environment Agency, January 2009. Science Report SC050021/SR2 Human Health Toxicological Assessment of Contaminants in Soil.
- 3) Environment Agency, January 2009. Science Report SC050021/SR3 Updated Technical Background to the CLEA Model.
- 4) Environment Agency, January 2009. Science Report SC050021/SR4 CLEA Software (Version 1.04) Handbook.
- 5) Environment Agency. 2008. Science Report SC050021/SR7. Compilation of Data for Priority Organic Pollutants for Derivation of Soil Guideline Values.
- 6) Land Quality Management (LQM) and Chartered Institute of Environmental Health (CIEH) 2009. The LQM/CIEH Generic Assessment Criteria for Human Health Risk Assessment. 2<sup>nd</sup> Edition.
- 7) The Soil Generic Assessment Criteria for Human Health Risk Assessment, Report ref. ISBN 978-1-905046-20-1, December 2009, published by CL:AIRE
- 8) Changes made to the CLEA framework documents after the 3 month evaluation period in 2008, released January 2009 by the Environment Agency.

### **APPENDIX F**

**HAS-WASTE Assessment** 



HASWASTE v4. Envirolab's Contaminated Land Soil Hazardous Waste Assessment Tool. Envirolab, Sandpits Business Park, Mottram Road, Hyde, Cheshire SK14 3AR.



#### Site Code and Name

241882 - Salisbury Square, Old Hatfileld

TD/MC/DLL	]	DU1	DLI1	DUI	DLII	BHO	DLIn	PLIO	BUO	PLIO	BHO	WC1	WC1	Wea	Wea
TP/WS/BH Depth (m)		BH1 0.20	BH1 0.70	BH1 1.5-1.7	BH1 2.3-2.5	BH2 0.25	BH2 0.50	BH2 0.90	BH2 1.40	BH2 3.00	BH2 4.90	WS1 0.2-0.3	WS1 0.5-0.6	WS2 0.2-0.3	WS2 0.5-0.6
Envirolab reference		mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg
Arsenic CrVI or Chromium		26 29	23 29	12 29	7 15	23 23		23 30	11 30	11 30	14 18		12 30		23 29
Copper Lead		17 14	50 278	14 21	3 5	17 14		16 43	33 46	18 35	9 16		26 68		17 17
Nickel Zinc		43 105	32 177	28 46	9	35 87		35 97	21	21	15 40		21		42 95
Cadmium		0.6	0.6	0.5	0.5	0.5		0.5	0.5	0.5	0.5		0.5		0.5
Mercury Selenium		0.2	0.2 2	0.2	0.2 1	0.2		0.2	0.2	0.2	0.2 1		0.2 1		0.2
Barium															
Beryllium Cobalt															
Manganese Molybdenum															
Total USEPA 16 PAHs															
Acenaphthene Acenaphthylene		0.01 0.01	0.01 0.05	0.02 0.01	0.01 0.01	0.01 0.01		0.04 0.14			0.02 0.01		0.02 0.01		0.01 0.01
Anthracene Benzo(a)anthracene		0.01 0.01	0.07 0.36	0.01 0.01	0.01 0.01	0.01 0.01		0.19 0.67			0.01 0.01		0.02 0.04		0.01 0.01
Benzo(a)pyrene		0.02	0.47	0.01	0.01	0.01		0.94			0.01		0.04		0.01
Benzo(b)fluoranthene Benzo(ghi)perylene		0.01	0.71	0.02	0.01 0.01	0.01		1.01			0.01		0.03		0.01 0.01
Benzo(k)fluoranthene Chrysene		0.02 0.01	0.48 0.70	0.02	0.01 0.01	0.01 0.02		0.76 1.48			0.01 0.02		0.03 0.13		0.01 0.01
Dibenzo(ah)anthracene Fluoranthene		0.01 0.01	0.10 0.78	0.01 0.07	0.01	0.01		0.14 1.90			0.01		0.01 0.14		0.01
Fluorene		0.01	0.01	0.01	0.01	0.01		0.03			0.01		0.01		0.01
Indeno(123cd)pyrene Naphthalene		0.01 0.01	0.27 0.03	0.01 0.11	0.01 0.02	0.01 0.01		0.58 0.04			0.01 0.02		0.03 0.02		0.01 0.01
Phenanthrene Pyrene		0.02 0.01	0.17 0.75	0.06 0.06	0.01 0.02	0.02 0.03		0.65 1.71			0.02 0.03		0.05 0.13		0.02 0.02
Benzo(j)fluoranthene	]										,				
Benzene Toluene		0.01 0.01													
Ethylbenzene		0.01													
Xylenes Trimethylbenzenes		0.01													
Chlorobenzene 1,2-Dichlorobenzene															
1,3-Dichlorobenzene															
1,4-Dichlorobenzene 1,2,4-Trichlorobenzene															
2-Chlorotoluene 4-Chlorotoluene															
Trichloroethene (TCE)															
Oil in Waste Carcinogenic H7 Total TPH	≥1,000mg/kg	0.1							ı			ı	10.0	10.0	
Petrol or (C6-C10)	≥1,000mg/kg	0.1											10.0	10.0	
Diesel or (C10-C25) or (conservative C10-C35)	≥10,000mg/kg														
Lube Oil or (C25+) or	≥1,000mg/kg														
(conservative C21+) 8 IARC H7 Carcinogenic PAHs marker	≥1%														
test (applicable to LRO only) Kerosene															
Kerosene															
Creosote Creosote															
pH Corrosive H8 (Irritant H4) pH (soil)	40.110.544.5														
pH (leachate)	≤2 H8 ≥11.5 ≤2 H8 ≥11.5														
Alkali Reserve (gNaOH/100g)  H4 Alkali Reserve test	≥13	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0
H8 Alkali Reserve test Produces Toxic Gases H12	≥14.5	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0
Total Sulphide	≥1,400mg/kg														
Free Cyanide Thiocyanate	≥1,200mg/kg ≥2,600mg/kg														
Elemental/Free Sulphur															
PCBs Total Phenols Total by HPLC	]														
Phenol		0.2		0.2			0.2			0.2		0.2		0.2	
Cresols Xylenols															
1-Naphthol															
Resourcinol 2,3,5,6-Tetrachlorophenol															
2,4,5-Trichlorophenol 2,4,6-Trichlorophenol															
2,4-Dichlorophenol															
4-Chloro-3-methylphenol Pentachlorophenol															
Bis(2-ethylhexyl)phthalate															
Butylbenzylphthalate Di-n-butylphthalate															
Visual Fibre Screen or Asbestos															
ID (enter Y or N)	H6≥25%														
Hazard Codes	Thresholds	%	%	%	%	%	%	%	%	%	%	%	%	%	%
Irritant H4 Irritant H4	≥10% ≥20%	0.000	0.000	0.000	0.000 0.002	0.000 0.007	0.000	0.000 0.007	0.000 0.004	0.000 0.004	0.000	0.000	0.000 0.004	0.000	0.000
Harmful H5	≥25%	0.009	0.053	0.010	0.002	0.007	0.000	0.007	0.004	0.011	0.003	0.000	0.018	0.000	0.013
Toxic H6 (Harmful H5)	≥0.1%H5<7%; H6≥7%	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
Toxic H6 (Harmful H5)	≥3%H5<25%; H6≥25%	0.013	0.010	0.008	0.003	0.011	0.000	0.011	0.006	0.006	0.005	0.000	0.006	0.000	0.012
Carcinogenic H7 Carcinogenic H7	≥0.1%	0.010	0.010 0.000	0.010 0.000	0.005 0.000	0.008	0.000	0.010 0.000	0.010 0.000	0.010	0.006 0.000	0.000	0.010 0.000	0.000	0.010
Corrosive H8 (Irritant H4)	≥5%H4<10%; H8≥10%	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
Toxic for Reproduction H10 Toxic for Reproduction H10	≥0.5%	0.009	0.041	0.006	0.002 0.001	0.007 0.002	0.000	0.007 0.006	0.007 0.007	0.005 0.005	0.003 0.002	0.000	0.010 0.010	0.000	0.008
Mutagenic H11	≥0.1%	0.000	0.000	0.000	0.000	0.002	0.000	0.000	0.007	0.000	0.002	0.000	0.000	0.000	0.002
Mutagenic H11 Ecotoxic H14	≥1% ≥1.0	0.009	0.006	0.006	0.002	0.007	0.000	0.007	0.004	0.004	0.003	0.000	0.004	0.000	0.008
New Ecotoxic H14 individual	>0.000501	0.00000	0.00004	0.00000	0.00000	0.00000	0.00000	0.00007	0.00000	0.00000	0.00000	0.00000	0.00000	0.00000	0.00000
substance specific thresholds New Ecotoxic H14 individual	≥0.0025%							0.00007	0.00000		0.00000				
substance specific thresholds	≥0.025%	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000

HASWASTE v4. Envirolab's Contaminated Land Soil Hazardous Waste Assessment Tool. Envirolab, Sandpits Business Park, Mottram Road, Hyde, Cheshire SK14 3AR.



### Site Code and Name

241882 - Salisbury Square, Old Hatfileld

TP/WS/BH Depth (m) Envirolab reference		WS3 0.2-0.3	WS3 0.5-0.6	WS4 0.2-0.3	WS4 0.5-0.6	TP1 0.10	TP1 0.40	TP2 0.50	TP3 0.10						
Arsenic	]	mg/kg	mg/kg	mg/kg 22	mg/kg	mg/kg 10	mg/kg	mg/kg	mg/kg 22	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg
CrVI or Chromium Copper				21 14	20 11	18 22		20 37	29 174						
Lead Nickel				39 30	10 33	66 14		84 18	345 33						
Zinc				73	70	80		112	306						
Cadmium Mercury				0.5 0.2	0.5 0.2	0.5 0.2		0.5 0.2	0.9 1.0						
Selenium Barium	] ]			1	1	1		1	2						
Beryllium Cobalt															
Manganese															
Molybdenum Total USEPA 16 PAHs	] ]														
Acenaphthene Acenaphthylene	j			0.01	0.01 0.01	0.02		0.13 0.01							
Anthracene				0.01	0.01	0.09		1.60							
Benzo(a)anthracene Benzo(a)pyrene				0.01 0.01	0.01 0.01	0.33 0.47		3.24 2.33							
Benzo(b)fluoranthene Benzo(ghi)perylene				0.01 0.01	0.01 0.01	0.44 0.68		2.29 1.90							
Benzo(k)fluoranthene Chrysene				0.01 0.01	0.01 0.02	0.36 0.71		2.64 5.27							
Dibenzo(ah)anthracene Fluoranthene				0.01	0.01	0.05 0.84		0.42 8.88							
Fluorene				0.01	0.01	0.01		0.17							
Indeno(123cd)pyrene Naphthalene				0.01	0.01	0.01		0.02							
Phenanthrene Pyrene				0.01 0.01	0.03 0.02	0.22 0.79		3.49 6.11							
Benzo(j)fluoranthene Benzene	]	0.04													
Toluene		0.01													
Ethylbenzene Xylenes		0.01 0.01													
Trimethylbenzenes Chlorobenzene	] ]														
1,2-Dichlorobenzene 1,3-Dichlorobenzene															
1,4-Dichlorobenzene															
1,2,4-Trichlorobenzene 2-Chlorotoluene															
4-Chlorotoluene Trichloroethene (TCE)	] ]														
Oil in Waste Carcinogenic H7															
Total TPH Petrol or (C6-C10)	≥1,000mg/kg ≥1,000mg/kg	10.0			0.0										
Diesel or (C10-C25) or (conservative C10-C35)	≥10,000mg/kg														
Lube Oil or (C25+) or	≥1,000mg/kg														
(conservative C21+)  8 IARC H7 Carcinogenic PAHs marker test (applicable to LRO only)		#DIV/0!	#DIV/0!	#DIV/0!	#DIV/0!	#DIV/0!	#DIV/0!	#DIV/0!	#DIV/0!	#DIV/0!	#DIV/0!	#DIV/0!	#DIV/0!	#DIV/0!	#DIV/0!
test (applicable to LRO only) Kerosene															
Kerosene Creosote	] ]														
Creosote															
pH Corrosive H8 (Irritant H4) pH (soil)	≤2 H8 ≥11.5														
pH (leachate) Alkali Reserve (gNaOH/100g)	≤2 H8 ≥11.5														
H4 Alkali Reserve test H8 Alkali Reserve test		0.0	0.0 0.0	0.0	0.0 0.0	0.0 0.0	0.0	0.0	0.0	0.0 0.0	0.0 0.0	0.0 0.0	0.0 0.0	0.0 0.0	0.0
Produces Toxic Gases H12															
Total Sulphide Free Cyanide	≥1,400mg/kg ≥1,200mg/kg														
Thiocyanate Elemental/Free Sulphur	≥2,600mg/kg														
PCBs Total	j				0.005										
Phenols Total by HPLC Phenol	<u> </u> 														
Cresols															
Xylenols 1-Naphthol															
Resourcinol 2,3,5,6-Tetrachlorophenol	] ]														
2,4,5-Trichlorophenol 2,4,6-Trichlorophenol															
2,4-Dichlorophenol															
4-Chloro-3-methylphenol Pentachlorophenol															
Bis(2-ethylhexyl)phthalate	]														
Butylbenzylphthalate Di-n-butylphthalate	]														
Visual Fibre Screen or Asbestos	H7≥0.1%; H5≥3%; H6≥25%														
ID (enter Y or N)	110620%														
Hazard Codes	Thresholds	0.000	% 0.000	% 0.000	% 0.000	% 0.000	%	% 0.000	% 0.000	% 0.000	% 0.000	% 0.000	% 0.000	% 0.000	% 0.000
Irritant H4 Irritant H4	≥10%	0.000	0.000	0.006	0.007	0.003	0.000	0.004	0.007	0.000	0.000	0.000	0.000	0.000	0.000
Harmful H5 Toxic H6 (Harmful H5)	≥25% ≥0.1%H5<7%;	0.001	0.000	0.013	0.009	0.015	0.000	0.020	0.077	0.000	0.000	0.000	0.000	0.000	0.000
Toxic H6 (Harmful H5)	H6≥7% ≥3%H5<25%;	0.000	0.000	0.009	0.009	0.004	0.000	0.006	0.010	0.000	0.000	0.000	0.000	0.000	0.000
Carcinogenic H7	H6≥25% ≥0.1%	0.000	0.000	0.007	0.007	0.006	0.000	0.007	0.010	0.000	0.000	0.000	0.000	0.000	0.000
Carcinogenic H7 Corrosive H8 (Irritant H4)	≥1% ≥5%H4<10%;	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
Toxic for Reproduction H10	H8≥10% ≥0.5%	0.000	0.000	0.006	0.007	0.010	0.000	0.012	0.050	0.000	0.000	0.000	0.000	0.000	0.000
Toxic for Reproduction H10 Mutagenic H11	≥5% ≥0.1%	0.000	0.000	0.006 0.000	0.001 0.000	0.010 0.000	0.000	0.012 0.000	0.050 0.000	0.000	0.000 0.000	0.000 0.000	0.000	0.000	0.000
Mutagenic H11 Ecotoxic H14	≥1% ≥1.0	0.000	0.000	0.006 0.120	0.007 0.105	0.003 0.117	0.000	0.004 0.159	0.007 0.445	0.000	0.000	0.000	0.000	0.000	0.000
New Ecotoxic H14 individual			1		1				1	1	1			1	
substance specific thresholds New Ecotoxic H14 individual	≥0.0025%	0.00000	0.00000	0.00000	0.00000	0.00003	0.00000	0.00032	0.00000	0.00000	0.00000	0.00000	0.00000	0.00000	0.00000
substance specific thresholds	≥0.025%	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000