

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG062215\
 Data File : BG017747.D
 Acq On : 22 Jun 2015 16:17
 Operator : TP/IZ
 Sample : G2653-03
 Misc : LOD-SVOC-SOIL-1PPM
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LOD-SOIL2015-01

Quant Time: Jun 23 03:47:45 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG061815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 19 06:05:07 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	74439	20.00	ng	0.00
21) Naphthalene-d8	10.39	136	325202	20.00	ng	-0.01
38) Acenaphthene-d10	14.26	164	195018	20.00	ng	-0.01
63) Phenanthrene-d10	17.01	188	442569	20.00	ng	-0.02
75) Chrysene-d12	21.21	240	471383	20.00	ng	-0.02
86) Perylene-d12	23.44	264	434124	20.00	ng	-0.02

System Monitoring Compounds						
5) 2-Fluorophenol	5.22	112	514219	121.23	ng	0.00
7) Phenol-d6	6.79	99	688757	115.53	ng	-0.01
23) Nitrobenzene-d5	8.76	82	422652	74.76	ng	-0.01
41) 2,4,6-Tribromophenol	15.76	330	190123	105.14	ng	-0.01
44) 2-Fluorobiphenyl	12.88	172	845072	70.38	ng	-0.02
78) Terphenyl-d14	19.66	244	1393335	67.16	ng	-0.02

Target Compounds						Qvalue
6) Aniline	6.95	93	2171	0.26	ng	# 96
8) 2-Chlorophenol	7.18	128	4763	0.90	ng	# 81
9) Benzaldehyde	6.75	77	3655	1.06	ng	# 84
11) bis(2-Chloroethyl)ether	7.05	93	3316	0.67	ng	# 98
12) 1,3-Dichlorobenzene	7.50	146	3809	0.65	ng	# 88
13) 1,4-Dichlorobenzene	7.64	146	3881	0.64	ng	# 84
14) 1,2-Dichlorobenzene	7.95	146	3911	0.68	ng	# 75
15) Benzyl Alcohol	7.85	79	2923	0.57	ng	# 83
16) 2,2'-oxybis(1-Chloropropan	8.15	45	6165	0.82	ng	# 87
17) 2-Methylphenol	8.05	107	3079	0.64	ng	# 93
18) Hexachloroethane	8.68	117	1523	0.65	ng	# 67
19) n-Nitroso-di-n-propylamine	8.41	70	2807	0.56	ng	# 93
20) 3+4-Methylphenols	8.38	107	4033	0.62	ng	# 89
22) Acetophenone	8.43	105	4569	0.62	ng	# 81
24) Nitrobenzene	8.80	77	4035	0.66	ng	# 75
25) Isophorone	9.32	82	7731	0.61	ng	# 93
26) 2-Nitrophenol	9.51	139	1851	0.67	ng	# 75
27) 2,4-Dimethylphenol	9.57	122	3280	0.62	ng	# 81
28) bis(2-Chloroethoxy)methane	9.82	93	4073	0.64	ng	# 93
29) 2,4-Dichlorophenol	10.04	162	2691	0.58	ng	# 88
30) 1,2,4-Trichlorobenzene	10.25	180	3086	0.59	ng	# 90
31) Naphthalene	10.44	128	11431	0.65	ng	# 93
33) 4-Chloroaniline	10.55	127	2997	0.39	ng	# 77
34) Hexachlorobutadiene	10.74	225	1781	0.60	ng	# 81
36) 4-Chloro-3-methylphenol	11.69	107	2864	0.50	ng	# 78
37) 2-Methylnaphthalene	12.06	142	7604m	0.60	ng	# 91
39) 1,2,4,5-Tetrachlorobenzene	12.43	216	3768	0.72	ng	# 91
42) 2,4,6-Trichlorophenol	12.69	196	1550m	0.44	ng	# 77
43) 2,4,5-Trichlorophenol	12.75	196	2227	0.58	ng	# 91
45) 1,1'-Biphenyl	13.09	154	9165	0.64	ng	# 85
46) 2-Chloronaphthalene	13.13	162	6902	0.60	ng	# 84
47) 2-Nitroaniline	13.34	65	2153	0.62	ng	# 97
48) Acenaphthylene	13.98	152	11961	0.58	ng	# 93
49) Dimethylphthalate	13.73	163	9325	0.63	ng	# 93

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
50) 2,6-Dinitrotoluene	13.84	165	1611	0.52	ng	93
51) Acenaphthene	14.33	154	7809	0.62	ng	95
52) 3-Nitroaniline	14.16	138	1620	0.45	ng	# 70
54) Dibenzofuran	14.66	168	11041	0.63	ng	97
55) 4-Nitrophenol	14.48	139	1287	0.45	ng	# 84
56) 2,4-Dinitrotoluene	14.63	165	1962	3.35	ng	# 82
57) Fluorene	15.31	166	9795	0.62	ng	90
58) 2,3,4,6-Tetrachlorophenol	14.90	232	1489	0.46	ng	# 89
59) Diethylphthalate	15.10	149	9938	0.61	ng	# 90
60) 4-Chlorophenyl-phenylether	15.32	204	4558	0.62	ng	91
61) 4-Nitroaniline	15.34	138	1971	0.49	ng	94
62) Azobenzene	15.61	77	10510	0.62	ng	91
65) n-Nitrosodiphenylamine	15.53	169	7169	0.51	ng	92
66) 4-Bromophenyl-phenylether	16.21	248	2354	0.51	ng	85
67) Hexachlorobenzene	16.32	284	2881	0.58	ng	# 73
68) Atrazine	16.49	200	2633	0.51	ng	95
69) Pentachlorophenol	16.66	266	944	0.38	ng	# 34
70) Phenanthrene	17.06	178	15785	0.63	ng	96
71) Anthracene	17.15	178	16323m	0.65	ng	
72) Carbazole	17.42	167	14152	0.58	ng	93
73) Di-n-butylphthalate	18.00	149	18381	0.60	ng	# 93
74) Fluoranthene	19.08	202	17849	0.63	ng	92
77) Pyrene	19.44	202	18951	0.63	ng	95
79) Butylbenzylphthalate	20.37	149	7496	0.53	ng	95
80) Benzo(a)anthracene	21.19	228	17749	0.65	ng	96
81) 3,3'-Dichlorobenzidine	21.14	252	3671	0.36	ng	86
82) Chrysene	21.25	228	15398	0.62	ng	95
83) Bis(2-ethylhexyl)phthalate	21.15	149	10650	0.54	ng	# 89
84) Di-n-octyl phthalate	22.02	149	15914	0.49	ng	# 85
85) Indeno(1,2,3-cd)pyrene	25.70	276	17358	0.58	ng	95
87) Benzo(b)fluoranthene	22.76	252	16588m	0.63	ng	
88) Benzo(k)fluoranthene	22.81	252	16387	0.66	ng	98
89) Benzo(a)pyrene	23.33	252	14718	0.61	ng	# 92
90) Dibenzo(a,h)anthracene	25.71	278	14445	0.59	ng	# 92
91) Benzo(g,h,i)perylene	26.38	276	13849	0.59	ng	# 85

(#) = qualifier out of range (m) = manual integration (+) = signals summed

