

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081415\
 Data File : BG018438.D
 Acq On : 14 Aug 2015 21:11
 Operator : UM/MA
 Sample : MDL-04-W
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Quant Time: Aug 17 04:03:08 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080915.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Aug 17 03:15:20 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.02	152	89315	20.00	ng/ul	0.00
18) Naphthalene-d8	10.82	136	394775	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.64	164	251963	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.39	188	629210	20.00	ng/ul	0.00
78) Chrysene-d12	21.65	240	625600	20.00	ng/ul	0.00
86) Perylene-d12	24.85	264	615775	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.45	96	13542	6.70	ng/uL	0.00
5) Phenol-d5	7.17	99	273438	33.73	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.34	67	174033	34.64	ng/ul	0.00
9) 2-Chlorophenol-d4	7.55	132	205104	33.04	ng/ul	0.00
13) 4-Methylphenol-d8	8.72	113	229258	33.66	ng/ul	0.00
19) Nitrobenzene-d5	9.17	128	112259	36.48	ng/ul	0.00
22) 2-Nitrophenol-d4	9.90	143	117355	37.44	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.44	165	221384	33.28	ng/ul	0.00
29) 4-Chloroaniline-d4	10.95	131	320541	42.64	ng/ul	0.00
44) Dimethylphthalate-d6	14.05	166	773465	36.48	ng/ul	0.00
47) Acenaphthylene-d8	14.34	160	941833	35.28	ng/ul	0.00
52) 4-Nitrophenol-d4	14.83	143	153394	40.23	ng/ul	0.00
58) Fluorene-d10	15.64	176	678797	36.76	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.74	200	117779	38.06	ng/ul	0.00
71) Anthracene-d10	17.49	188	1097960	36.49	ng/ul	0.00
79) Pyrene-d10	19.77	212	1189797	36.66	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.63	264	1233858	37.74	ng/ul	0.00

Target Compounds

					Qvalue
2) 1,4-Dioxane	3.48	88	1674	0.77 ng/uL	95
4) Benzaldehyde	7.15	77	16930	3.56 ng/ul	91
6) Phenol	7.20	94	23461	2.84 ng/ul	98
8) Bis(2-Chloroethyl)ether	7.43	93	19656	3.09 ng/ul	98
10) 2-Chlorophenol	7.58	128	18073	2.86 ng/ul	93
11) 2-Methylphenol	8.46	108	17513	2.73 ng/ul	93
12) 2,2'-oxybis(1-Chloropropan	8.55	45	25049	2.45 ng/ul#	92
14) Acetophenone	8.83	105	29859	3.04 ng/ul	94
15) N-Nitroso-di-n-propylamine	8.81	70	16937	3.00 ng/ul#	85
16) 4-Methylphenol	8.78	108	20023	2.86 ng/ul	97
17) Hexachloroethane	9.10	117	7212	2.64 ng/ul	93
20) Nitrobenzene	9.22	77	22260	2.87 ng/ul#	90
21) Isophorone	9.74	82	47640	3.19 ng/ul#	94
23) 2-Nitrophenol	9.93	139	11478	3.30 ng/ul	94
24) 2,4-Dimethylphenol	9.98	107	20910	2.68 ng/ul	97
25) Bis(2-Chloroethoxy)methane	10.22	93	27694	2.97 ng/ul	97
27) 2,4-Dichlorophenol	10.46	162	18150	2.91 ng/ul	92
28) Naphthalene	10.87	128	61979	2.97 ng/ul	98
30) 4-Chloroaniline	10.98	127	29042	3.80 ng/ul#	77
31) Hexachlorobutadiene	11.16	225	11432	2.90 ng/ul	88
32) Caprolactam	11.75	113	7228m	2.88 ng/ul	
33) 4-Chloro-3-methylphenol	12.10	107	23376	3.24 ng/ul#	93
34) 2-Methylnaphthalene	12.47	142	46832	3.10 ng/ul	88

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
37) 1,2,4,5-Tetrachlorobenzene	12.84	216	25288	3.13	ng/ul	97
38) Hexachlorocyclopentadiene	12.83	237	11860	2.47	ng/ul#	70
39) 2,4,6-Trichlorophenol	13.08	196	14990	2.72	ng/ul	96
40) 2,4,5-Trichlorophenol	13.14	196	15687	2.65	ng/ul	96
41) 1,1'-Biphenyl	13.48	154	63650	3.03	ng/ul	93
42) 2-Chloronaphthalene	13.52	162	50118	3.07	ng/ul	98
43) 2-Nitroaniline	13.71	65	15351	2.91	ng/ul	98
45) Dimethylphthalate	14.09	163	63850	3.05	ng/ul	96
46) 2,6-Dinitrotoluene	14.21	165	12153	2.92	ng/ul	88
48) Acenaphthylene	14.37	152	82991	3.10	ng/ul	98
49) 3-Nitroaniline	14.54	138	14961m	3.52	ng/ul	
50) Acenaphthene	14.71	153	56591	3.01	ng/ul	96
51) 2,4-Dinitrophenol	14.75	184	4384	2.16	ng/ul#	73
53) 4-Nitrophenol	14.84	109	10542	3.18	ng/ul#	82
54) Dibenzofuran	15.04	168	78777	3.21	ng/ul	94
55) 2,4-Dinitrotoluene	14.99	165	17627	2.97	ng/ul#	97
56) 2,3,4,6-Tetrachlorophenol	15.26	232	15281	2.96	ng/ul#	96
57) Diethylphthalate	15.45	149	72780	3.28	ng/ul	99
59) Fluorene	15.69	166	66155	3.13	ng/ul	94
60) 4-Chlorophenyl-phenylether	15.68	204	31438	3.15	ng/ul	98
61) 4-Nitroaniline	15.69	138	15358	3.27	ng/ul	97
64) 4,6-Dinitro-2-methylphenol	15.76	198	10592	3.21	ng/ul#	63
65) N-Nitrosodiphenylamine	15.89	169	58335	3.08	ng/ul	95
66) 4-Bromophenyl-phenylether	16.58	248	19728	2.90	ng/ul	95
67) Hexachlorobenzene	16.69	284	22633	3.15	ng/ul	94
68) Atrazine	16.83	200	22815	3.22	ng/ul	98
69) Pentachlorophenol	17.04	266	10952	2.27	ng/ul#	87
70) Phenanthrene	17.43	178	114957	3.37	ng/ul	95
72) Anthracene	17.52	178	115560	3.32	ng/ul	96
73) 1,2,3,4-Tetrachlorobenzene	13.45	216	23287	2.67	ng/uL	95
74) Pentachlorobenzene	14.97	250	23664	2.85	ng/uL	91
75) Carbazole	17.79	167	105998	3.47	ng/ul	96
76) Di-n-butylphthalate	18.34	149	134117	3.37	ng/ul#	98
77) Fluoranthene	19.43	202	135442	3.71	ng/ul	99
80) Pyrene	19.79	202	143375	3.39	ng/ul	96
81) Butylbenzylphthalate	20.68	149	57477	3.21	ng/ul	95
82) 3,3'-Dichlorobenzidine	21.54	252	47717	3.67	ng/ul	93
83) Benzo(a)anthracene	21.63	228	127898	3.30	ng/ul	99
84) Bis(2-ethylhexyl)phthalate	21.54	149	83725	3.31	ng/ul	98
85) Chrysene	21.69	228	121779	3.36	ng/ul	94
87) Di-n-octyl phthalate	22.75	149	142266	3.58	ng/ul	100
88) Benzo(b)fluoranthene	23.83	252	122800m	3.17	ng/ul	
89) Benzo(k)fluoranthene	23.90	252	117920	3.17	ng/ul	99
91) Benzo(a)pyrene	24.70	252	116794	3.18	ng/ul	97
92) Indeno(1,2,3-cd)pyrene	28.48	276	132553m	3.11	ng/ul	
93) Dibenzo(a,h)anthracene	28.55	278	109251	3.09	ng/ul	96
94) Benzo(g,h,i)perylene	29.62	276	108276m	3.03	ng/ul	

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

