

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM020915\
 Data File : BM000463.D
 Acq On : 10 Feb 2015 15:56
 Operator : TP/IZ
 Sample : G1249-11
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4FL9

Quant Time: Feb 11 04:01:18 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-BM020615.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Feb 11 03:41:38 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	105647	20.00	ng/ul	0.00
18) Naphthalene-d8	10.52	136	413245	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.37	164	259017	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.12	188	595562	20.00	ng/ul	0.00
75) Chrysene-d12	21.32	240	632034	20.00	ng/ul	0.00
83) Perylene-d12	23.56	264	590618	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0	0.00	ng/uL	
5) Phenol-d5	6.89	99	214892	30.62	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.06	67	130901	30.32	ng/ul	0.00
9) 2-Chlorophenol-d4	7.26	132	186391	30.54	ng/ul	0.00
13) 4-Methylphenol-d8	8.43	113	178260	30.66	ng/ul	0.00
19) Nitrobenzene-d5	8.88	128	87734	33.89	ng/ul	0.00
22) 2-Nitrophenol-d4	9.60	143	95758	32.72	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.13	165	188691	29.93	ng/ul	0.00
29) 4-Chloroaniline-d4	10.65	131	39871	6.88	ng/ul	0.00
43) Dimethylphthalate-d6	13.79	166	603568	33.82	ng/ul	0.00
46) Acenaphthylene-d8	14.07	160	725826	29.77	ng/ul	0.00
51) 4-Nitrophenol-d4	14.57	143	97654	33.44	ng/ul	0.00
57) Fluorene-d10	15.37	176	535537	29.81	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.49	200	112242	35.61	ng/ul	0.00
70) Anthracene-d10	17.22	188	835594	30.85	ng/ul	0.00
76) Pyrene-d10	19.52	212	909994	30.55	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.42	264	822982	31.20	ng/ul	0.00

Target Compounds

						Qvalue
6) Phenol	6.92	94	131225	18.09	ng/ul	96
8) Bis(2-Chloroethyl)ether	7.15	93	297500	50.88	ng/ul	97
11) 2-Methylphenol	8.16	108	313055	55.37	ng/ul	98
15) N-Nitroso-di-n-propylamine	8.53	70	251133	56.34	ng/ul	97
16) 4-Methylphenol	8.49	108	330243	53.92	ng/ul	93
17) Hexachloroethane	8.79	117	92650	33.35	ng/ul	93
21) Isophorone	9.45	82	445358	37.70	ng/ul	99
24) 2,4-Dimethylphenol	9.69	107	305180	40.69	ng/ul	97
25) Bis(2-Chloroethoxy)methane	9.93	93	320469	44.34	ng/ul	99
27) 2,4-Dichlorophenol	10.16	162	92309	15.00	ng/ul	96
28) Naphthalene	10.56	128	435906	20.94	ng/ul	99
31) Hexachlorobutadiene	10.85	225	247036	52.03	ng/ul	99
32) Caprolactam	11.42	113	60552	44.92	ng/ul	87
36) 1,2,4,5-Tetrachlorobenzene	12.56	216	288654	33.83	ng/ul	98
37) Hexachlorocyclopentadiene	12.53	237	166165	33.09	ng/ul	97
39) 2,4,5-Trichlorophenol	12.87	196	186545	35.75	ng/ul	99
41) 2-Chloronaphthalene	13.24	162	279435	17.36	ng/ul	99
44) Dimethylphthalate	13.83	163	356771	18.19	ng/ul	99
47) Acenaphthylene	14.09	152	513871	20.19	ng/ul	99
48) 3-Nitroaniline	14.28	138	13426	4.11	ng/ul#	98
49) Acenaphthene	14.44	153	498402	29.62	ng/ul	95
50) 2,4-Dinitrophenol	14.49	184	86026	48.76	ng/ul	92
53) Dibenzofuran	14.77	168	426268	17.36	ng/ul	94

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58) Fluorene	15.43	166	143740	7.25	ng/ul	96
59) 4-Chlorophenyl-phenylether	15.42	204	290007	28.15	ng/ul	93
60) 4-Nitroaniline	15.45	138	217552	59.65	ng/ul	99
63) 4,6-Dinitro-2-methylphenol	15.51	198	230656	69.54	ng/ul#	96
68) Pentachlorophenol	16.77	266	278834	82.00	ng/ul	99
71) Anthracene	17.25	178	951909	29.25	ng/ul	99
72) Carbazole	17.53	167	1007612	36.52	ng/ul	98
74) Fluoranthene	19.19	202	633014	17.81	ng/ul	100
78) Butylbenzylphthalate	20.45	149	298793	23.08	ng/ul	94
81) Bis(2-ethylhexyl)phthalate	21.23	149	420131	22.90	ng/ul	98
82) Chrysene	21.36	228	991160	28.82	ng/ul	99
86) Benzo(k)fluoranthene	22.93	252	1201585	35.53	ng/ul	99
88) Benzo(a)pyrene	23.47	252	1158233	34.82	ng/ul	99
89) Indeno(1,2,3-cd)pyrene	25.82	276	503955	13.67	ng/ul	98

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

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