

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052717\
 Data File : BM010273.D
 Acq On : 27 May 2017 13:42
 Operator : SJ/MA
 Sample : SSTD08005
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD08005

Quant Time: May 27 14:43:36 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM052717.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat May 27 14:42:28 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.93	152	301730	20.00	ng/ul	0.00
18) Naphthalene-d8	10.74	136	1147322	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.57	164	717271	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.31	188	1668468	20.00	ng/ul	0.00
75) Chrysene-d12	21.47	240	2329088	20.00	ng/ul	0.00
83) Perylene-d12	23.83	264	2358800	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.38	96	247905	33.71	ng/uL	0.00
5) Phenol-d5	7.12	99	1927163	86.40	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.28	67	1068387	84.76	ng/ul	0.00
9) 2-Chlorophenol-d4	7.48	132	1624636	89.15	ng/ul	0.00
13) 4-Methylphenol-d8	8.66	113	1543363	85.72	ng/ul	0.00
19) Nitrobenzene-d5	9.12	128	739082	87.94	ng/ul	0.00
22) 2-Nitrophenol-d4	9.83	143	893473	91.78	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.36	165	1748624	87.43	ng/ul	0.00
29) 4-Chloroaniline-d4	10.90	131	1696786	89.72	ng/ul	0.00
43) Dimethylphthalate-d6	13.99	166	5134592	86.16	ng/ul	0.00
46) Acenaphthylene-d8	14.27	160	6092206	87.63	ng/ul	0.00
51) 4-Nitrophenol-d4	14.82	143	780095	93.55	ng/ul	0.01
57) Fluorene-d10	15.56	176	4496480	85.83	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.69	200	1042007	93.71	ng/ul	0.01
70) Anthracene-d10	17.42	188	7047535	86.81	ng/ul	0.00
76) Pyrene-d10	19.70	212	8328057	86.47	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.68	264	9599001	87.41	ng/ul	0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.42	88	264159	33.70	ng/uL#	53
4) Benzaldehyde	7.09	77	1278910	80.95	ng/ul	96
6) Phenol	7.15	94	1962164	85.89	ng/ul	95
8) Bis(2-Chloroethyl)ether	7.37	93	1372089	86.08	ng/ul	94
10) 2-Chlorophenol	7.50	128	1615219	88.32	ng/ul	95
11) 2-Methylphenol	8.39	108	1457485	87.77	ng/ul	98
12) 2,2'-oxybis(1-Chloropropan	8.46	45	616475	85.42	ng/ul#	62
14) Acetophenone	8.78	105	2540303	88.05	ng/ul#	96
15) N-Nitroso-di-n-propylamine	8.75	70	1369069	90.98	ng/ul#	84
16) 4-Methylphenol	8.73	108	1577730	86.68	ng/ul	92
17) Hexachloroethane	9.00	117	731838	88.12	ng/ul	91
20) Nitrobenzene	9.16	77	2112161	88.00	ng/ul	96
21) Isophorone	9.68	82	3320393	89.90	ng/ul	96
23) 2-Nitrophenol	9.87	139	918730	89.00	ng/ul	89
24) 2,4-Dimethylphenol	9.92	107	2054841	85.62	ng/ul	96
25) Bis(2-Chloroethoxy)methane	10.15	93	1782514	85.87	ng/ul	94
27) 2,4-Dichlorophenol	10.39	162	1707000	87.11	ng/ul	93
28) Naphthalene	10.79	128	4830355	83.94	ng/ul	100
30) 4-Chloroaniline	10.93	127	1613094	88.55	ng/ul	96
31) Hexachlorobutadiene	11.03	225	1449355	80.74	ng/ul	94
32) Caprolactam	11.75	113	224060	47.01	ng/ul#	72
33) 4-Chloro-3-methylphenol	12.05	107	1695511	89.90	ng/ul	99
34) 2-Methylnaphthalene	12.39	142	3752708	86.01	ng/ul	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
36) 1,2,4,5-Tetrachlorobenzene	12.75	216	2475058	82.95	ng/ul	98
37) Hexachlorocyclopentadiene	12.71	237	1670548	85.33	ng/ul	98
38) 2,4,6-Trichlorophenol	13.00	196	1499448	88.41	ng/ul	97
39) 2,4,5-Trichlorophenol	13.08	196	1490069	87.10	ng/ul	98
40) 1,1'-Biphenyl	13.40	154	4962215	85.28	ng/ul	100
41) 2-Chloronaphthalene	13.45	162	3900455	85.33	ng/ul	97
42) 2-Nitroaniline	13.69	65	1140774	97.98	ng/ul	96
44) Dimethylphthalate	14.03	163	4974281	84.87	ng/ul	99
45) 2,6-Dinitrotoluene	14.17	165	1002083	93.28	ng/ul	91
47) Acenaphthylene	14.30	152	5959568	87.72	ng/ul	99
48) 3-Nitroaniline	14.52	138	916057	91.31	ng/ul	94
49) Acenaphthene	14.63	153	4072202	85.47	ng/ul	97
50) 2,4-Dinitrophenol	14.70	184	718509	93.54	ng/ul	89
52) 4-Nitrophenol	14.83	109	1068760	95.01	ng/ul	94
53) Dibenzofuran	14.97	168	6054190	85.93	ng/ul	99
54) 2,4-Dinitrotoluene	14.95	165	1490939	94.51	ng/ul	99
55) 2,3,4,6-Tetrachlorophenol	15.19	232	1488393	90.57	ng/ul	93
56) Diethylphthalate	15.38	149	4968923	87.65	ng/ul	95
58) Fluorene	15.62	166	5086578	88.01	ng/ul	97
59) 4-Chlorophenyl-phenylether	15.60	204	2869760	86.88	ng/ul	94
60) 4-Nitroaniline	15.69	138	985116	96.16	ng/ul	96
63) 4,6-Dinitro-2-methylphenol	15.70	198	1079480	91.11	ng/ul#	98
64) N-Nitrosodiphenylamine	15.83	169	4200635	86.37	ng/ul	99
65) 4-Bromophenyl-phenylether	16.50	248	1772616	86.35	ng/ul	94
66) Hexachlorobenzene	16.59	284	1820536	84.68	ng/ul#	84
67) Atrazine	16.77	200	1796642	88.20	ng/ul	96
68) Pentachlorophenol	16.96	266	1198193	91.97	ng/ul	97
69) Phenanthrene	17.36	178	7636218	85.20	ng/ul	99
71) Anthracene	17.45	178	8145545	87.60	ng/ul	98
72) Carbazole	17.73	167	6781063	87.06	ng/ul	98
73) Di-n-butylphthalate	18.25	149	8294813	93.36	ng/ul	99
74) Fluoranthene	19.36	202	9851726	86.74	ng/ul	99
77) Pyrene	19.73	202	10371120	86.41	ng/ul	98
78) Butylbenzylphthalate	20.60	149	3952955	95.30	ng/ul	98
79) 3,3'-Dichlorobenzidine	21.40	252	3969980	85.97	ng/ul	95
80) Benzo(a)anthracene	21.46	228	11038893	84.05	ng/ul	96
81) Bis(2-ethylhexyl)phthalate	21.34	149	5960764	96.58	ng/ul#	99
82) Chrysene	21.52	228	9966060	83.94	ng/ul	98
84) Di-n-octyl phthalate	22.24	149	10304030	90.96	ng/ul#	95
85) Benzo(b)fluoranthene	23.11	252	11873269	86.51	ng/ul#	99
86) Benzo(k)fluoranthene	23.16	252	11522956	87.88	ng/ul#	98
88) Benzo(a)pyrene	23.73	252	11791398	86.67	ng/ul#	98
89) Indeno(1,2,3-cd)pyrene	26.24	276	14990018	87.21	ng/ul#	95
90) Dibenzo(a,h)anthracene	26.25	278	12970322	87.62	ng/ul#	97
91) Benzo(g,h,i)perylene	26.99	276	12041918	85.04	ng/ul#	98

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

