

Data Path : Z:\HPCHEM1\BNA_G\Data\BG101417\
 Data File : BG029190.D
 Acq On : 13 Oct 2017 11:18
 Operator : SJ/JU
 Sample : SSTD00557
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTD00557

Quant Time: Oct 13 16:54:58 2017
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG101417.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 13 16:52:26 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.17	152	35985	20.00	ng/ul	0.00
18) Naphthalene-d8	10.99	136	150179	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.79	164	156812	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.52	188	279533	20.00	ng/ul	0.00
75) Chrysene-d12	21.79	240	304730	20.00	ng/ul	0.00
83) Perylene-d12	25.10	264	299491	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.54	96	1485	4.60	ng/uL	0.00
5) Phenol-d5	7.31	99	17207	4.93	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.49	67	11368	4.93	ng/ul	0.00
9) 2-Chlorophenol-d4	7.70	132	11371	4.67	ng/ul	0.00
13) 4-Methylphenol-d8	8.87	113	12777	3.90	ng/ul	0.00
19) Nitrobenzene-d5	9.33	128	4780	4.10	ng/ul	0.00
22) 2-Nitrophenol-d4	10.05	143	4713	4.34	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.60	165	12909	4.80	ng/ul	0.00
29) 4-Chloroaniline-d4	11.11	131	12101	3.94	ng/ul	0.00
43) Dimethylphthalate-d6	14.19	166	70155	5.36	ng/ul	0.00
46) Acenaphthylene-d8	14.48	160	84450	5.15	ng/ul	0.00
51) 4-Nitrophenol-d4	14.96	143	9148	3.38	ng/ul	0.00
57) Fluorene-d10	15.77	176	48673	4.21	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.88	200	5163	3.68	ng/ul	0.00
70) Anthracene-d10	17.62	188	67315	5.02	ng/ul	0.00
76) Pyrene-d10	19.89	212	78865	4.79	ng/ul	0.00
87) Benzo(a)pyrene-d12	24.87	264	74337	4.84	ng/ul	-0.01

Target Compounds

					Qvalue
2) 1,4-Dioxane	3.59	88	2304	5.536 ng/uL#	83
4) Benzaldehyde	7.30	77	13762	5.451 ng/ul	97
6) Phenol	7.35	94	17702	4.923 ng/ul	92
8) Bis(2-Chloroethyl)ether	7.58	93	13198	4.683 ng/ul	96
10) 2-Chlorophenol	7.73	128	11680	4.720 ng/ul	95
11) 2-Methylphenol	8.60	108	11402	4.386 ng/ul	89
12) 2,2'-oxybis(1-Chloropropan	8.70	45	24881	4.938 ng/ul	96
14) Acetophenone	8.99	105	21180	3.676 ng/ul	95
15) N-Nitroso-di-n-propylamine	8.97	70	12179	4.188 ng/ul#	96
16) 4-Methylphenol	8.93	108	12415	3.348 ng/ul	97
17) Hexachloroethane	9.26	117	4952	4.139 ng/ul#	78
20) Nitrobenzene	9.38	77	13727	4.083 ng/ul#	97
21) Isophorone	9.90	82	32859	4.821 ng/ul	95
23) 2-Nitrophenol	10.09	139	5015	4.088 ng/ul#	89
24) 2,4-Dimethylphenol	10.14	107	15653	4.681 ng/ul	92
25) Bis(2-Chloroethoxy)methane	10.38	93	18895	4.877 ng/ul	97
27) 2,4-Dichlorophenol	10.62	162	11995	4.672 ng/ul	93
28) Naphthalene	11.04	128	39991	4.913 ng/ul	99
30) 4-Chloroaniline	11.14	127	11126	3.671 ng/ul	98
31) Hexachlorobutadiene	11.32	225	9895	4.829 ng/ul	91
32) Caprolactam	11.88	113	4865	3.708 ng/ul#	77
33) 4-Chloro-3-methylphenol	12.25	107	14736	4.677 ng/ul	89
34) 2-Methylnaphthalene	12.63	142	47930	6.538 ng/ul	90

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36) 1,2,4,5-Tetrachlorobenzene	13.00	216	21842	5.063	ng/ul#	94
37) Hexachlorocyclopentadiene	12.97	237	10493	4.201	ng/ul#	97
38) 2,4,6-Trichlorophenol	13.23	196	15223	5.201	ng/ul	95
39) 2,4,5-Trichlorophenol	13.30	196	16169	5.200	ng/ul	97
40) 1,1'-Biphenyl	13.63	154	68210	6.171	ng/ul	99
41) 2-Chloronaphthalene	13.67	162	50434	5.938	ng/ul	94
42) 2-Nitroaniline	13.86	65	13198	4.939	ng/ul	96
44) Dimethylphthalate	14.23	163	69893	5.281	ng/ul	98
45) 2,6-Dinitrotoluene	14.34	165	9005	4.200	ng/ul#	91
47) Acenaphthylene	14.51	152	87793	5.327	ng/ul	99
48) 3-Nitroaniline	14.68	138	8847	3.790	ng/ul#	83
49) Acenaphthene	14.85	153	57414	4.975	ng/ul	100
50) 2,4-Dinitrophenol	14.88	184	3832	3.080	ng/ul	89
52) 4-Nitrophenol	14.97	109	7920	3.471	ng/ul#	80
53) Dibenzofuran	15.18	168	82578	4.942	ng/ul	98
54) 2,4-Dinitrotoluene	15.14	165	12968	3.757	ng/ul#	98
55) 2,3,4,6-Tetrachlorophenol	15.40	232	13682	4.198	ng/ul#	99
56) Diethylphthalate	15.59	149	70784	4.575	ng/ul	99
58) Fluorene	15.83	166	51159	4.046	ng/ul	95
59) 4-Chlorophenyl-phenylether	15.82	204	26472	4.215	ng/ul	96
60) 4-Nitroaniline	15.83	138	9833	3.041	ng/ul#	84
63) 4,6-Dinitro-2-methylphenol	15.89	198	5012	3.472	ng/ul#	83
64) N-Nitrosodiphenylamine	16.02	169	39378	4.704	ng/ul	98
65) 4-Bromophenyl-phenylether	16.71	248	14924	4.931	ng/ul	97
66) Hexachlorobenzene	16.83	284	16241	5.093	ng/ul	95
67) Atrazine	16.96	200	17948	5.054	ng/ul	93
68) Pentachlorophenol	17.17	266	8390	4.449	ng/ul	88
69) Phenanthrene	17.56	178	74604	4.956	ng/ul	98
71) Anthracene	17.66	178	77858	5.003	ng/ul	98
72) Carbazole	17.92	167	71149	5.000	ng/ul	99
73) Di-n-butylphthalate	18.47	149	81475	4.798	ng/ul	98
74) Fluoranthene	19.56	202	96007	4.361	ng/ul	97
77) Pyrene	19.92	202	104284	4.912	ng/ul	99
78) Butylbenzylphthalate	20.80	149	37395	4.314	ng/ul	98
79) 3,3'-Dichlorobenzidine	21.68	252	30166	4.733	ng/ul	95
80) Benzo(a)anthracene	21.78	228	101851	5.094	ng/ul	97
81) Bis(2-ethylhexyl)phthalate	21.67	149	57306	4.576	ng/ul	99
82) Chrysene	21.84	228	91013	5.049	ng/ul	97
84) Di-n-octyl phthalate	22.92	149	94810	4.019	ng/ul	100
85) Benzo(b)fluoranthene	24.11	252	91245	4.662	ng/ul	97
86) Benzo(k)fluoranthene	24.11	252	91245	4.831	ng/ul	97
88) Benzo(a)pyrene	24.94	252	92172	4.991	ng/ul#	95
89) Indeno(1,2,3-cd)pyrene	28.87	276	103993	5.477	ng/ul#	96
90) Dibenzo(a,h)anthracene	28.94	278	88718	5.525	ng/ul#	93
91) Benzo(g,h,i)perylene	30.05	276	88498	5.566	ng/ul	94

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

