

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061720\
 Data File : BG045793.D
 Acq On : 17 Jun 2020 17:11
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040EC

Quant Time: Jun 17 17:48:27 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG061520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 15 18:34:19 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	44947	20.00	ng	0.00
21) Naphthalene-d8	10.72	136	170133	20.00	ng	0.00
39) Acenaphthene-d10	14.57	164	121165	20.00	ng	0.00
64) Phenanthrene-d10	17.32	188	282911	20.00	ng	0.00
76) Chrysene-d12	21.58	240	264180	20.00	ng	0.00
86) Perylene-d12	24.71	264	287998	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.53	112	224806	84.49	ng	0.00
7) Phenol-d6	7.16	99	342200	84.55	ng	0.00
23) Nitrobenzene-d5	9.08	82	323611	86.89	ng	0.00
42) 2,4,6-Tribromophenol	16.07	330	157194	85.93	ng	0.00
45) 2-Fluorobiphenyl	13.19	172	710598	86.19	ng	0.00
79) Terphenyl-d14	19.94	244	1188144	91.15	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.37	88	47957	39.352	ng	100
3) Pyridine	3.77	79	147112	40.737	ng	99
4) n-Nitrosodimethylamine	3.68	42	49753	41.044	ng	# 97
6) Aniline	7.25	93	200410	42.054	ng	97
8) 2-Chlorophenol	7.51	128	118519	41.614	ng	92
9) Benzaldehyde	7.05	77	100365	41.287	ng	95
10) Phenol	7.19	94	163282	41.637	ng	99
11) bis(2-Chloroethyl)ether	7.34	93	123243	41.375	ng	98
12) 1,3-Dichlorobenzene	7.81	146	143058	42.149	ng	99
13) 1,4-Dichlorobenzene	7.96	146	143584	42.353	ng	95
14) 1,2-Dichlorobenzene	8.27	146	136051	41.919	ng	97
15) Benzyl Alcohol	8.18	79	132681	42.363	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.45	45	114431	40.762	ng	83
17) 2-Methylphenol	8.41	107	123790	42.416	ng	96
18) Hexachloroethane	9.00	117	55300	42.949	ng	94
19) n-Nitroso-di-n-propylamine	8.72	70	113870	42.795	ng	99
20) 3+4-Methylphenols	8.75	107	163481	42.593	ng	# 73
22) Acetophenone	8.74	105	197539	42.303	ng	# 99
24) Nitrobenzene	9.12	77	168033	42.316	ng	99
25) Isophorone	9.65	82	310994	43.136	ng	99
26) 2-Nitrophenol	9.84	139	72740	43.970	ng	97
27) 2,4-Dimethylphenol	9.93	122	108558	42.981	ng	98
28) bis(2-Chloroethoxy)methane	10.13	93	160726	42.678	ng	97
29) 2,4-Dichlorophenol	10.41	162	129374	43.903	ng	97
30) 1,2,4-Trichlorobenzene	10.59	180	148136	42.464	ng	98
31) Naphthalene	10.77	128	397625	42.749	ng	99
32) Benzoic acid	10.13	122	76735	47.062	ng	92
33) 4-Chloroaniline	10.90	127	166341	42.705	ng	97
34) Hexachlorobutadiene	11.07	225	108782	43.846	ng	98
35) Caprolactam	11.66	113	45377	47.586	ng	93
36) 4-Chloro-3-methylphenol	12.07	107	149442	43.923	ng	97
37) 2-Methylnaphthalene	12.39	142	296819	42.854	ng	97
38) 1-Methylnaphthalene	12.61	142	280225	42.754	ng	95
40) 1,2,4,5-Tetrachlorobenzene	12.76	216	169230	42.417	ng	98

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41) Hexachlorocyclopentadiene	12.74	237	72851	37.997	nq	98
43) 2,4,6-Trichlorophenol	13.02	196	115788	42.805	nq	92
44) 2,4,5-Trichlorophenol	13.12	196	130565	42.351	nq	98
46) 1,1'-Biphenyl	13.40	154	369892	42.640	nq	100
47) 2-Chloronaphthalene	13.44	162	296376	42.295	nq	98
48) 2-Nitroaniline	13.66	65	99661	42.757	nq	96
49) Acenaphthylene	14.29	152	474619	42.286	nq	99
50) Dimethylphthalate	14.03	163	412970	43.507	nq	99
51) 2,6-Dinitrotoluene	14.15	165	93439	43.828	nq	98
52) Acenaphthene	14.63	154	286774	42.152	nq	98
53) 3-Nitroaniline	14.49	138	92631	43.500	nq	99
54) 2,4-Dinitrophenol	14.69	184	46415	44.161	nq	98
55) Dibenzofuran	14.97	168	472534	42.755	nq	99
56) 4-Nitrophenol	14.88	139	63309	43.408	nq	97
57) 2,4-Dinitrotoluene	14.94	165	133479	43.698	nq	96
58) Fluorene	15.62	166	397710	43.276	nq	98
59) 2,3,4,6-Tetrachlorophenol	15.22	232	115287	43.085	nq	97
60) Diethylphthalate	15.39	149	426395	43.943	nq	97
61) 4-Chlorophenyl-phenylether	15.61	204	226931	42.843	nq	98
62) 4-Nitroaniline	15.66	138	101436	46.414	nq	96
63) Azobenzene	15.90	77	391858	42.058	nq	97
65) 4,6-Dinitro-2-methylphenol	15.72	198	79987	45.749	nq	100
66) n-Nitrosodiphenylamine	15.83	169	344953	43.048	nq	98
67) 4-Bromophenyl-phenylether	16.51	248	159062	43.270	nq	98
68) Hexachlorobenzene	16.63	284	167196	43.862	nq	95
69) Atrazine	16.78	200	138315	44.482	nq	99
70) Pentachlorophenol	16.99	266	69389	41.021	nq	97
71) Phenanthrene	17.36	178	633573	43.422	nq	99
72) Anthracene	17.45	178	629863	43.350	nq	100
73) Carbazole	17.74	167	621958	44.957	nq	98
74) Di-n-butylphthalate	18.28	149	735552	44.903	nq	100
75) Fluoranthene	19.37	202	778340	43.381	nq	100
77) Benzidine	19.56	184	325293	47.364	nq	100
78) Pyrene	19.74	202	797707	45.092	nq	99
80) Butylbenzylphthalate	20.63	149	322652	43.574	nq	98
81) Benzo(a)anthracene	21.56	228	740163	43.214	nq	99
82) 3,3'-Dichlorobenzidine	21.48	252	280181	43.246	nq	100
83) Chrysene	21.62	228	720008	43.462	nq	100
84) Bis(2-ethylhexyl)phthalate	21.47	149	444442	43.151	nq	# 98
85) Di-n-octyl phthalate	22.65	149	756130	42.560	nq	99
87) Indeno(1,2,3-cd)pyrene	28.27	276	908665	43.521	nq	# 95
88) Benzo(b)fluoranthene	23.72	252	781466	43.277	nq	99
89) Benzo(k)fluoranthene	23.79	252	757211	43.814	nq	99
90) Benzo(a)pyrene	24.56	252	736955	43.690	nq	100
91) Dibenzo(a,h)anthracene	28.33	278	730451	43.628	nq	99
92) Benzo(g,h,i)perylene	29.37	276	731849	43.672	nq	98

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

