

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021119\
 Data File : BG039431.D
 Acq On : 11 Feb 2019 12:58
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Quant Time: Feb 12 00:12:47 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG012919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 29 15:41:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.17	152	61776	20.00	ng	-0.01
21) Naphthalene-d8	10.99	136	287224	20.00	ng	-0.01
39) Acenaphthene-d10	14.80	164	198746	20.00	ng	-0.01
64) Phenanthrene-d10	17.54	188	480211	20.00	ng	-0.01
76) Chrysene-d12	21.83	240	462726	20.00	ng	-0.01
87) Perylene-d12	25.21	264	481197	20.00	ng	-0.03

System Monitoring Compounds

5) 2-Fluorophenol	5.71	112	290580	80.51	ng	0.00
7) Phenol-d6	7.31	99	445993	83.94	ng	0.00
23) Nitrobenzene-d5	9.35	82	425191	92.48	ng	-0.01
42) 2,4,6-Tribromophenol	16.28	330	192527	89.96	ng	0.00
45) 2-Fluorobiphenyl	13.42	172	1039079	75.00	ng	-0.01
79) Terphenyl-d14	20.14	244	1597333	73.07	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.60	88	59665	37.790	ng	95
3) Pyridine	4.01	79	181631	43.450	ng	93
4) n-Nitrosodimethylamine	3.93	42	77034	36.848	ng	90
6) Aniline	7.50	93	273936	41.163	ng	99
8) 2-Chlorophenol	7.73	128	165707	41.999	ng	98
9) Benzaldehyde	7.31	77	112018	43.175	ng	95
10) Phenol	7.34	94	230248	41.573	ng	98
11) bis(2-Chloroethyl)ether	7.59	93	177376	40.530	ng	97
12) 1,3-Dichlorobenzene	8.06	146	191903	40.150	ng	98
13) 1,4-Dichlorobenzene	8.21	146	194676	40.864	ng	97
14) 1,2-Dichlorobenzene	8.53	146	188885	40.956	ng	99
15) Benzyl Alcohol	8.41	79	177308	42.508	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.69	45	352482	35.666	ng	99
17) 2-Methylphenol	8.60	107	150616	42.024	ng	97
18) Hexachloroethane	9.26	117	67735	41.327	ng	96
19) n-Nitroso-di-n-propylamine	8.97	70	154289	40.915	ng	96
20) 3+4-Methylphenols	8.93	107	213688	43.297	ng	98
22) Acetophenone	9.00	105	301452	39.072	ng	# 94
24) Nitrobenzene	9.39	77	216980	43.684	ng	95
25) Isophorone	9.91	82	397938	39.058	ng	98
26) 2-Nitrophenol	10.10	139	101860	51.252	ng	95
27) 2,4-Dimethylphenol	10.14	122	165183	40.079	ng	97
28) bis(2-Chloroethoxy)methane	10.39	93	247568	39.450	ng	97
29) 2,4-Dichlorophenol	10.63	162	191200	42.447	ng	96
30) 1,2,4-Trichlorobenzene	10.85	180	205734	40.682	ng	99
31) Naphthalene	11.05	128	548062	38.463	ng	99
32) Benzoic acid	10.27	122	117345	44.329	ng	92
33) 4-Chloroaniline	11.15	127	267185	40.441	ng	98
34) Hexachlorobutadiene	11.31	225	138804	41.272	ng	96
35) Caprolactam	11.95	113	83873	44.997	ng	# 90
36) 4-Chloro-3-methylphenol	12.25	107	217422	41.736	ng	97
37) 2-Methylnaphthalene	12.63	142	409779	38.828	ng	98
38) 1-Methylnaphthalene	12.85	142	399085	39.229	ng	97
40) 1,2,4,5-Tetrachlorobenzene	12.99	216	251461	39.766	ng	98

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41) Hexachlorocyclopentadiene	12.96	237	149914	43.507	ng	96
43) 2,4,6-Trichlorophenol	13.23	196	169163	43.509	ng	99
44) 2,4,5-Trichlorophenol	13.30	196	176045	43.201	ng	99
46) 1,1'-Biphenyl	13.63	154	618561	37.407	ng	99
47) 2-Chloronaphthalene	13.68	162	483393	38.859	ng	99
48) 2-Nitroaniline	13.89	65	165243	46.545	ng	96
49) Acenaphthylene	14.53	152	716254	38.158	ng	99
50) Dimethylphthalate	14.24	163	634126	39.506	ng	99
51) 2,6-Dinitrotoluene	14.37	165	139371	46.668	ng	97
52) Acenaphthene	14.86	154	458227	37.608	ng	99
53) 3-Nitroaniline	14.71	138	148713	46.190	ng	# 93
54) 2,4-Dinitrophenol	14.91	184	84004	68.252	ng	91
55) Dibenzofuran	15.20	168	757752	38.788	ng	99
56) 4-Nitrophenol	14.99	139	122690	45.006	ng	94
57) 2,4-Dinitrotoluene	15.15	165	197600	50.864	ng	89
58) Fluorene	15.84	166	556561	38.217	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.41	232	172851	45.303	ng	97
60) Diethylphthalate	15.60	149	634881	38.255	ng	97
61) 4-Chlorophenyl-phenylether	15.83	204	333676	40.465	ng	96
62) 4-Nitroaniline	15.87	138	153585	45.836	ng	91
63) Azobenzene	16.12	77	585486	36.094	ng	97
65) 4,6-Dinitro-2-methylphenol	15.91	198	108417	54.395	ng	96
66) n-Nitrosodiphenylamine	16.04	169	554910	38.003	ng	98
67) 4-Bromophenyl-phenylether	16.72	248	219443	41.191	ng	96
68) Hexachlorobenzene	16.83	284	222352	41.600	ng	95
69) Atrazine	16.98	200	197645	37.040	ng	96
70) Pentachlorophenol	17.17	266	159876	43.037	ng	98
71) Phenanthrene	17.59	178	931770	37.489	ng	98
72) Anthracene	17.67	178	925871	37.819	ng	97
73) Carbazole	17.94	167	906049	37.084	ng	99
74) Di-n-butylphthalate	18.48	149	1059327	37.165	ng	99
75) Fluoranthene	19.58	202	1082390	37.520	ng	99
77) Benzidine	19.77	184	585401	38.587	ng	99
78) Pyrene	19.95	202	1111954	38.602	ng	99
80) Butylbenzylphthalate	20.82	149	483444	39.785	ng	93
81) Benzo(a)anthracene	21.81	228	1045387	38.233	ng	100
82) 3,3'-Dichlorobenzidine	21.72	252	408626	40.305	ng	98
83) Chrysene	21.88	228	1006635	38.675	ng	100
84) Bis(2-ethylhexyl)phthalate	21.69	149	677575	39.086	ng	98
85) Di-n-octyl phthalate	22.95	149	1117106	39.005	ng	97
86) Indeno(1,2,3-cd)pyrene	29.11	276	1164069	41.684	ng	98
88) Benzo(b)fluoranthene	24.13	252	1018067	38.795	ng	98
89) Benzo(k)fluoranthene	24.21	252	1008724	38.286	ng	99
90) Benzo(a)pyrene	25.05	252	996262	39.419	ng	99
91) Dibenzo(a,h)anthracene	29.18	278	953289	40.277	ng	99
92) Benzo(g,h,i)perylene	30.33	276	964785	40.896	ng	99

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

