

Data Path : Z:\HPCHEM1\BNA G\DATA\BG101617\
 Data File : BG029282.D
 Acq On : 16 Oct 2017 22:07
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Quant Time: Oct 17 04:02:30 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG101617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 16 17:32:15 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.17	152	62433	20.00	ng	0.00
21) Naphthalene-d8	10.99	136	293483	20.00	ng	0.00
38) Acenaphthene-d10	14.79	164	182527	20.00	ng	0.00
63) Phenanthrene-d10	17.53	188	409408	20.00	ng	0.00
75) Chrysene-d12	21.80	240	424335	20.00	ng	0.00
86) Perylene-d12	25.11	264	434521	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.72	112	320364	85.72	ng	0.00
7) Phenol-d6	7.32	99	452780	86.31	ng	0.00
23) Nitrobenzene-d5	9.34	82	398473	86.28	ng	0.00
41) 2,4,6-Tribromophenol	16.27	330	205704	87.29	ng	0.00
44) 2-Fluorobiphenyl	13.42	172	1047563	84.17	ng	0.00
78) Terphenyl-d14	20.11	244	1567227	84.02	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.59	88	64959	42.839	ng	85
3) Pyridine	3.99	79	193594	43.842	ng	90
4) n-Nitrosodimethylamine	3.90	42	88027	42.646	ng	# 94
6) Aniline	7.49	93	277604	42.735	ng	96
8) 2-Chlorophenol	7.73	128	183747	42.894	ng	91
9) Benzaldehyde	7.30	77	112068	42.473	ng	93
10) Phenol	7.35	94	241994	42.789	ng	91
11) bis(2-Chloroethyl)ether	7.58	93	178580	43.340	ng	93
12) 1,3-Dichlorobenzene	8.06	146	204284	42.476	ng	96
13) 1,4-Dichlorobenzene	8.21	146	211291	43.030	ng	94
14) 1,2-Dichlorobenzene	8.53	146	201002	42.440	ng	97
15) Benzyl Alcohol	8.40	79	160204	43.336	ng	91
16) 2,2'-oxybis(1-Chloropropan	8.70	45	295556	42.236	ng	96
17) 2-Methylphenol	8.61	107	159412	42.432	ng	98
18) Hexachloroethane	9.26	117	70599	42.190	ng	96
19) n-Nitroso-di-n-propylamine	8.97	70	152267	42.689	ng	# 97
20) 3+4-Methylphenols	8.94	107	226831	42.850	ng	93
22) Acetophenone	8.99	105	294933	41.700	ng	# 94
24) Nitrobenzene	9.38	77	220195	42.405	ng	91
25) Isophorone	9.90	82	411006	42.630	ng	# 93
26) 2-Nitrophenol	10.09	139	112387	45.284	ng	90
27) 2,4-Dimethylphenol	10.14	122	190759	42.483	ng	91
28) bis(2-Chloroethoxy)methane	10.38	93	251457	42.534	ng	97
29) 2,4-Dichlorophenol	10.63	162	205103	42.834	ng	94
30) 1,2,4-Trichlorobenzene	10.85	180	218538	42.245	ng	99
31) Naphthalene	11.04	128	617383	42.210	ng	98
32) Benzoic acid	10.29	122	163200	43.407	ng	# 83
33) 4-Chloroaniline	11.14	127	264484	42.987	ng	97
34) Hexachlorobutadiene	11.32	225	134264	41.744	ng	99
35) Caprolactam	11.91	113	70100	44.050	ng	# 78
36) 4-Chloro-3-methylphenol	12.25	107	223773	42.491	ng	91
37) 2-Methylnaphthalene	12.63	142	450721	42.281	ng	93
39) 1,2,4,5-Tetrachlorobenzene	12.99	216	280298	42.061	ng	98
40) Hexachlorocyclopentadiene	12.98	237	97851	41.163	ng	90

Data Path : Z:\HPCHEM1\BNA G\DATA\BG101617\
 Data File : BG029282.D
 Acq On : 16 Oct 2017 22:07
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Quant Time: Oct 17 04:02:30 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG101617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 16 17:32:15 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.23	196	181605	43.411	ng	99
43) 2,4,5-Trichlorophenol	13.30	196	185297	43.130	ng	95
45) 1,1'-Biphenyl	13.63	154	661917	42.190	ng	96
46) 2-Chloronaphthalene	13.68	162	481015	42.033	ng	97
47) 2-Nitroaniline	13.87	65	138928	44.199	ng	# 86
48) Acenaphthylene	14.52	152	762146	42.555	ng	99
49) Dimethylphthalate	14.24	163	608698	42.742	ng	99
50) 2,6-Dinitrotoluene	14.36	165	136399	45.689	ng	# 83
51) Acenaphthene	14.86	154	501307	43.020	ng	99
52) 3-Nitroaniline	14.69	138	144450	44.840	ng	# 81
53) 2,4-Dinitrophenol	14.89	184	61971	41.203	ng	# 53
54) Dibenzofuran	15.19	168	705310	42.399	ng	99
55) 4-Nitrophenol	14.99	139	118095	44.466	ng	# 82
56) 2,4-Dinitrotoluene	15.14	165	184273	45.597	ng	# 79
57) Fluorene	15.83	166	580999	42.278	ng	98
58) 2,3,4,6-Tetrachlorophenol	15.41	232	155711	43.441	ng	95
59) Diethylphthalate	15.59	149	603370	42.512	ng	98
60) 4-Chlorophenyl-phenylether	15.82	204	308588	42.117	ng	97
61) 4-Nitroaniline	15.84	138	143580	43.405	ng	88
62) Azobenzene	16.11	77	507860	42.434	ng	96
64) 4,6-Dinitro-2-methylphenol	15.91	198	97252	42.421	ng	84
65) n-Nitrosodiphenylamine	16.03	169	520318	42.426	ng	99
66) 4-Bromophenyl-phenylether	16.71	248	199620	42.492	ng	97
67) Hexachlorobenzene	16.84	284	223649	42.028	ng	96
68) Atrazine	16.97	200	191767	43.822	ng	99
69) Pentachlorophenol	17.18	266	135100	44.195	ng	99
70) Phenanthrene	17.57	178	893300	42.236	ng	99
71) Anthracene	17.66	178	907111	42.713	ng	100
72) Carbazole	17.92	167	862522	42.278	ng	98
73) Di-n-butylphthalate	18.48	149	1009053	43.005	ng	100
74) Fluoranthene	19.57	202	1055559	42.670	ng	97
76) Benzidine	19.74	184	561938	43.860	ng	98
77) Pyrene	19.93	202	1093778	42.492	ng	97
79) Butylbenzylphthalate	20.80	149	469645	42.825	ng	94
80) Benzo(a)anthracene	21.78	228	1048911	42.102	ng	97
81) 3,3'-Dichlorobenzidine	21.69	252	435828	42.383	ng	99
82) Chrysene	21.85	228	996207	41.753	ng	96
83) Bis(2-ethylhexyl)phthalate	21.67	149	666254	42.332	ng	99
84) Di-n-octyl phthalate	22.92	149	1180740	43.045	ng	97
85) Indeno(1,2,3-cd)pyrene	28.90	276	1261583	42.980	ng	# 89
87) Benzo(b)fluoranthene	24.06	252	1073440	43.151	ng	99
88) Benzo(k)fluoranthene	24.13	252	1080214	43.586	ng	98
89) Benzo(a)pyrene	24.96	252	1040756	43.519	ng	97
90) Dibenzo(a,h)anthracene	28.97	278	1045035	43.162	ng	96
91) Benzo(g,h,i)perylene	30.09	276	937218	41.661	ng	98

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

