

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG110117\
 Data File : BG030626.D
 Acq On : 1 Nov 2017 9:17
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Quant Time: Nov 01 17:41:07 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG101617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 01 17:41:24 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.16	152	74826	20.00	ng	0.00
21) Naphthalene-d8	10.97	136	358921	20.00	ng	0.00
38) Acenaphthene-d10	14.77	164	234150	20.00	ng	0.00
63) Phenanthrene-d10	17.51	188	551637	20.00	ng	0.00
75) Chrysene-d12	21.79	240	606110	20.00	ng	0.00
86) Perylene-d12	25.10	264	626086	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.71	112	346843	77.44	ng	0.00
7) Phenol-d6	7.32	99	483543	76.91	ng	0.00
23) Nitrobenzene-d5	9.33	82	436052	77.20	ng	0.00
41) 2,4,6-Tribromophenol	16.26	330	253921	83.99	ng	0.00
44) 2-Fluorobiphenyl	13.40	172	1191490	74.63	ng	0.00
78) Terphenyl-d14	20.10	244	1874022	70.34	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.58	88	73180	40.268	ng	# 81
3) Pyridine	3.99	79	212803	40.210	ng	89
4) n-Nitrosodimethylamine	3.90	42	90273	36.491	ng	# 81
6) Aniline	7.48	93	303430	38.975	ng	93
8) 2-Chlorophenol	7.72	128	188358	36.688	ng	89
9) Benzaldehyde	7.29	77	139710	44.180	ng	89
10) Phenol	7.34	94	243607	35.940	ng	90
11) bis(2-Chloroethyl)ether	7.57	93	195560	39.600	ng	88
12) 1,3-Dichlorobenzene	8.05	146	215154	37.327	ng	97
13) 1,4-Dichlorobenzene	8.19	146	222671	37.837	ng	95
14) 1,2-Dichlorobenzene	8.52	146	216804	38.195	ng	99
15) Benzyl Alcohol	8.39	79	178171	40.214	ng	90
16) 2,2'-oxybis(1-Chloropropan	8.68	45	216598	25.826	ng	94
17) 2-Methylphenol	8.60	107	166162	36.903	ng	96
18) Hexachloroethane	9.25	117	75968	37.880	ng	99
19) n-Nitroso-di-n-propylamine	8.96	70	172432	40.336	ng	# 94
20) 3+4-Methylphenols	8.93	107	241534	38.071	ng	94
22) Acetophenone	8.98	105	328852	38.018	ng	# 94
24) Nitrobenzene	9.37	77	221973	34.954	ng	# 88
25) Isophorone	9.89	82	414884	35.186	ng	# 94
26) 2-Nitrophenol	10.08	139	108369	35.704	ng	# 85
27) 2,4-Dimethylphenol	10.13	122	177492	32.321	ng	93
28) bis(2-Chloroethoxy)methane	10.37	93	276998	38.311	ng	95
29) 2,4-Dichlorophenol	10.62	162	211846	36.176	ng	95
30) 1,2,4-Trichlorobenzene	10.84	180	245313	38.775	ng	97
31) Naphthalene	11.03	128	659564	36.872	ng	98
32) Benzoic acid	10.27	122	139739	30.391	ng	84
33) 4-Chloroaniline	11.13	127	292974	38.936	ng	94
34) Hexachlorobutadiene	11.31	225	154772	39.347	ng	98
35) Caprolactam	11.90	113	77770	39.960	ng	# 81
36) 4-Chloro-3-methylphenol	12.24	107	224900	34.919	ng	# 86
37) 2-Methylnaphthalene	12.62	142	504305	38.682	ng	93
39) 1,2,4,5-Tetrachlorobenzene	12.98	216	332184	38.857	ng	97
40) Hexachlorocyclopentadiene	12.96	237	128287	41.966	ng	88

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.22	196	189688	35.346	ng	99
43) 2,4,5-Trichlorophenol	13.29	196	204134	37.039	ng	93
45) 1,1'-Biphenyl	13.61	154	735557	36.547	ng	97
46) 2-Chloronaphthalene	13.66	162	531683	36.218	ng	99
47) 2-Nitroaniline	13.86	65	147927	36.686	ng	# 83
48) Acenaphthylene	14.50	152	863711	37.593	ng	99
49) Dimethylphthalate	14.22	163	695696	38.081	ng	100
50) 2,6-Dinitrotoluene	14.35	165	147011	38.387	ng	# 79
51) Acenaphthene	14.84	154	540787	36.177	ng	98
52) 3-Nitroaniline	14.67	138	159386	38.568	ng	# 82
53) 2,4-Dinitrophenol	14.89	184	50085	28.412	ng	# 47
54) Dibenzofuran	15.17	168	841882	39.451	ng	99
55) 4-Nitrophenol	14.98	139	121126	35.552	ng	# 79
56) 2,4-Dinitrotoluene	15.13	165	204203	39.388	ng	# 76
57) Fluorene	15.82	166	671147	38.071	ng	96
58) 2,3,4,6-Tetrachlorophenol	15.40	232	188043	40.895	ng	# 93
59) Diethylphthalate	15.57	149	705386	38.743	ng	99
60) 4-Chlorophenyl-phenylether	15.80	204	382668	40.713	ng	95
61) 4-Nitroaniline	15.84	138	168408	39.686	ng	81
62) Azobenzene	16.10	77	587423	38.261	ng	95
64) 4,6-Dinitro-2-methylphenol	15.90	198	109723	36.334	ng	83
65) n-Nitrosodiphenylamine	16.02	169	642712	38.894	ng	99
66) 4-Bromophenyl-phenylether	16.70	248	247214	39.055	ng	99
67) Hexachlorobenzene	16.82	284	264141	36.839	ng	94
68) Atrazine	16.96	200	236836	40.167	ng	99
69) Pentachlorophenol	17.17	266	156857	38.083	ng	98
70) Phenanthrene	17.56	178	1067396	37.455	ng	98
71) Anthracene	17.65	178	1082107	37.816	ng	99
72) Carbazole	17.92	167	997755	36.297	ng	99
73) Di-n-butylphthalate	18.46	149	1178134	37.265	ng	99
74) Fluoranthene	19.56	202	1299986	39.001	ng	96
76) Benzidine	19.73	184	720368	39.363	ng	98
77) Pyrene	19.91	202	1336496	36.350	ng	97
79) Butylbenzylphthalate	20.78	149	562873	35.933	ng	86
80) Benzo(a)anthracene	21.77	228	1362940	38.300	ng	97
81) 3,3'-Dichlorobenzidine	21.68	252	565190	38.479	ng	98
82) Chrysene	21.84	228	1283530	37.662	ng	97
83) Bis(2-ethylhexyl)phthalate	21.65	149	802853	35.713	ng	99
84) Di-n-octyl phthalate	22.89	149	1368167	34.920	ng	96
85) Indeno(1,2,3-cd)pyrene	28.89	276	1663408	39.674	ng	98
87) Benzo(b)fluoranthene	24.05	252	1416190	39.510	ng	99
88) Benzo(k)fluoranthene	24.12	252	1376036	38.534	ng	98
89) Benzo(a)pyrene	24.94	252	1326407	38.493	ng	99
90) Dibenzo(a,h)anthracene	28.95	278	1406989	40.331	ng	98
91) Benzo(g,h,i)perylene	30.09	276	1353201	41.748	ng	97

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

