

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122216\
 Data File : BG025392.D
 Acq On : 22 Dec 2016 17:01
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040EC

Quant Time: Dec 23 00:17:24 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG122116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 21 18:42:01 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.21	152	69003	20.00	ng	0.00
21) Naphthalene-d8	11.04	136	292269	20.00	ng	0.00
38) Acenaphthene-d10	14.84	164	225507	20.00	ng	0.00
63) Phenanthrene-d10	17.59	188	531235	20.00	ng	0.00
75) Chrysene-d12	21.88	240	717434	20.00	ng	0.00
86) Perylene-d12	25.29	264	757078	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.75	112	323570	85.23	ng	0.00
7) Phenol-d6	7.37	99	470615	88.02	ng	0.00
23) Nitrobenzene-d5	9.40	82	545429	82.44	ng	0.00
41) 2,4,6-Tribromophenol	16.33	330	304235	83.80	ng	0.00
44) 2-Fluorobiphenyl	13.46	172	1180592	84.76	ng	0.00
78) Terphenyl-d14	20.17	244	2178262	80.50	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.60	88	66562	41.09	ng	95
3) Pyridine	4.01	79	208087	44.32	ng	97
4) n-Nitrosodimethylamine	3.93	42	119289	42.80	ng	93
6) Aniline	7.53	93	312897	43.19	ng	95
8) 2-Chlorophenol	7.77	128	178882	41.77	ng	96
9) Benzaldehyde	7.35	77	154830	39.58	ng	99
10) Phenol	7.40	94	252453	42.60	ng	97
11) bis(2-Chloroethyl)ether	7.62	93	189886	41.58	ng	88
12) 1,3-Dichlorobenzene	8.10	146	217548	41.98	ng	95
13) 1,4-Dichlorobenzene	8.25	146	223101	41.54	ng	97
14) 1,2-Dichlorobenzene	8.57	146	211824	41.33	ng	92
15) Benzyl Alcohol	8.46	79	221124	43.32	ng	95
16) 2,2'-oxybis(1-Chloropropan	8.73	45	355137	41.99	ng	97
17) 2-Methylphenol	8.66	107	169710	43.60	ng	98
18) Hexachloroethane	9.30	117	90052	43.66	ng	# 83
19) n-Nitroso-di-n-propylamine	9.02	70	192756	42.68	ng	92
20) 3+4-Methylphenols	8.99	107	233516	43.35	ng	95
22) Acetophenone	9.05	105	335549	41.32	ng	# 95
24) Nitrobenzene	9.44	77	294164	40.54	ng	98
25) Isophorone	9.95	82	498907	41.65	ng	98
26) 2-Nitrophenol	10.15	139	115122	41.48	ng	99
27) 2,4-Dimethylphenol	10.20	122	193427	42.39	ng	96
28) bis(2-Chloroethoxy)methane	10.43	93	247818	39.84	ng	98
29) 2,4-Dichlorophenol	10.69	162	208662	42.73	ng	97
30) 1,2,4-Trichlorobenzene	10.90	180	240438	40.76	ng	93
31) Naphthalene	11.09	128	596668	40.88	ng	100
32) Benzoic acid	10.36	122	165036	44.98	ng	95
33) 4-Chloroaniline	11.20	127	262073	42.69	ng	96
34) Hexachlorobutadiene	11.35	225	197001	40.91	ng	100
35) Caprolactam	12.00	113	73627	40.13	ng	94
36) 4-Chloro-3-methylphenol	12.32	107	246838	40.96	ng	93
37) 2-Methylnaphthalene	12.68	142	450198	41.62	ng	94
39) 1,2,4,5-Tetrachlorobenzene	13.04	216	328464	44.12	ng	96
40) Hexachlorocyclopentadiene	13.00	237	109472	44.48	ng	91

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.28	196	198026	42.24	nq	98
43) 2,4,5-Trichlorophenol	13.37	196	216531	44.13	nq	93
45) 1,1'-Biphenyl	13.67	154	649059	42.79	nq	96
46) 2-Chloronaphthalene	13.73	162	499379	42.14	nq	98
47) 2-Nitroaniline	13.94	65	223506	44.68	nq	96
48) Acenaphthylene	14.57	152	776100	42.26	nq	97
49) Dimethylphthalate	14.28	163	688219	41.86	nq	97
50) 2,6-Dinitrotoluene	14.42	165	151298	42.13	nq	89
51) Acenaphthene	14.90	154	515291	42.90	nq	99
52) 3-Nitroaniline	14.77	138	151641	43.92	nq	95
53) 2,4-Dinitrophenol	14.98	184	72054	33.73	nq	97
54) Dibenzofuran	15.24	168	800772	42.17	nq	96
55) 4-Nitrophenol	15.09	139	114578	44.43	nq	# 80
56) 2,4-Dinitrotoluene	15.22	165	228005	43.60	nq	# 95
57) Fluorene	15.88	166	667613	41.99	nq	99
58) 2,3,4,6-Tetrachlorophenol	15.46	232	216021	41.10	nq	95
59) Diethylphthalate	15.63	149	702520	40.75	nq	98
60) 4-Chlorophenyl-phenylether	15.86	204	378369	41.99	nq	97
61) 4-Nitroaniline	15.93	138	152745	43.96	nq	# 82
62) Azobenzene	16.16	77	704916	42.40	nq	98
64) 4,6-Dinitro-2-methylphenol	15.98	198	152128	41.35	nq	87
65) n-Nitrosodiphenylamine	16.08	169	595162	41.45	nq	99
66) 4-Bromophenyl-phenylether	16.76	248	273596	41.75	nq	98
67) Hexachlorobenzene	16.89	284	311913	41.47	nq	# 89
68) Atrazine	17.02	200	250922	39.88	nq	97
69) Pentachlorophenol	17.24	266	155703	39.94	nq	93
70) Phenanthrene	17.63	178	1131527	41.33	nq	95
71) Anthracene	17.71	178	1135397	40.82	nq	99
72) Carbazole	17.99	167	1038115	41.72	nq	97
73) Di-n-butylphthalate	18.51	149	1244233	40.65	nq	96
74) Fluoranthene	19.62	202	1519013	42.01	nq	94
76) Benzidine	19.80	184	644825	36.03	nq	99
77) Pyrene	19.99	202	1546449	41.24	nq	97
79) Butylbenzylphthalate	20.83	149	632942	42.04	nq	99
80) Benzo(a)anthracene	21.86	228	1682784	42.05	nq	98
81) 3,3'-Dichlorobenzidine	21.76	252	652222	41.96	nq	99
82) Chrysene	21.93	228	1592322	41.49	nq	99
83) Bis(2-ethylhexyl)phthalate	21.70	149	893437	42.64	nq	98
84) Di-n-octyl phthalate	22.96	149	1483282	42.64	nq	96
85) Indeno(1,2,3-cd)pyrene	29.23	276	2077833	42.59	nq	# 73
87) Benzo(b)fluoranthene	24.20	252	1743796	42.22	nq	# 95
88) Benzo(k)fluoranthene	24.27	252	1652665	41.43	nq	# 96
89) Benzo(a)pyrene	25.13	252	1668112	41.81	nq	98
90) Dibenzo(a,h)anthracene	29.28	278	1780799	42.12	nq	# 98
91) Benzo(g,h,i)perylene	30.46	276	1732198	41.22	nq	97

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

