

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG040615\
 Data File : BG016233.D
 Acq On : 6 Apr 2015 20:50
 Operator : TP/IZ
 Sample : G1614-06
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LOQ-WATER2015-02

Quant Time: Apr 07 05:11:22 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG040615.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 07 04:17:17 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	65533	20.00	ng	0.00
21) Naphthalene-d8	10.49	136	303583	20.00	ng	0.00
38) Acenaphthene-d10	14.34	164	178168	20.00	ng	0.00
63) Phenanthrene-d10	17.08	188	386265	20.00	ng	0.00
75) Chrysene-d12	21.26	240	380806	20.00	ng	0.00
86) Perylene-d12	23.51	264	343375	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.29	112	713506	171.83	ng	0.00
7) Phenol-d6	6.88	99	1014537	162.76	ng	0.00
23) Nitrobenzene-d5	8.86	82	535365	102.16	ng	0.00
41) 2,4,6-Tribromophenol	15.83	330	200629	166.51	ng	0.00
44) 2-Fluorobiphenyl	12.96	172	1070906	102.34	ng	0.00
78) Terphenyl-d14	19.71	244	1431905	88.00	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.18	88	36869	19.53	ng	97
3) Pyridine	3.57	79	100503	17.81	ng	99
4) n-Nitrosodimethylamine	3.49	42	34313	20.46	ng	100
6) Aniline	7.03	93	96148	10.80	ng	98
8) 2-Chlorophenol	7.27	128	121863	24.44	ng	96
9) Benzaldehyde	6.84	77	92021	25.20	ng	96
10) Phenol	6.90	94	161930	24.28	ng	99
11) bis(2-Chloroethyl)ether	7.13	93	112421	21.14	ng	100
12) 1,3-Dichlorobenzene	7.59	146	108074	21.46	ng	98
13) 1,4-Dichlorobenzene	7.74	146	111115	21.27	ng	98
14) 1,2-Dichlorobenzene	8.05	146	105763	21.17	ng	99
15) Benzyl Alcohol	7.94	79	94932	21.85	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.24	45	128201	21.40	ng	98
17) 2-Methylphenol	8.14	107	102379	22.89	ng	98
18) Hexachloroethane	8.77	117	41168	20.60	ng	99
19) n-Nitroso-di-n-propylamine	8.50	70	85718	19.19	ng	99
20) 3+4-Methylphenols	8.47	107	142150	22.95	ng	99
22) Acetophenone	8.52	105	155355	22.07	ng	99
24) Nitrobenzene	8.90	77	121305	21.65	ng	95
25) Isophorone	9.42	82	242203	20.81	ng	100
26) 2-Nitrophenol	9.60	139	60047	22.98	ng	97
27) 2,4-Dimethylphenol	9.67	122	114037	23.43	ng	99
28) bis(2-Chloroethoxy)methane	9.90	93	147144	22.15	ng	100
29) 2,4-Dichlorophenol	10.14	162	98073	25.46	ng	96
30) 1,2,4-Trichlorobenzene	10.35	180	90000	22.38	ng	100
31) Naphthalene	10.54	128	345712	21.85	ng	99
32) Benzoic acid	9.77	122	48088	21.12	ng	97
33) 4-Chloroaniline	10.65	127	53058	7.15	ng	98
34) Hexachlorobutadiene	10.82	225	39419	22.32	ng	97
35) Caprolactam	11.40	113	54327	22.40	ng	97
36) 4-Chloro-3-methylphenol	11.78	107	118561	23.30	ng	97
37) 2-Methylnaphthalene	12.15	142	251776	22.13	ng	98
39) 1,2,4,5-Tetrachlorobenzene	12.52	216	85642	21.83	ng	99
40) Hexachlorocyclopentadiene	12.50	237	37632	20.96	ng	95

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG040615\
 Data File : BG016233.D
 Acq On : 6 Apr 2015 20:50
 Operator : TP/IZ
 Sample : G1614-06
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LOQ-WATER2015-02

Quant Time: Apr 07 05:11:22 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG040615.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 07 04:17:17 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	12.77	196	65880	22.60	ng	97
43) 2,4,5-Trichlorophenol	12.84	196	72826	23.38	ng	99
45) 1,1'-Biphenyl	13.17	154	305332	22.34	ng	98
46) 2-Chloronaphthalene	13.21	162	230530	22.14	ng	100
47) 2-Nitroaniline	13.42	65	74865	22.31	ng	98
48) Acenaphthylene	14.06	152	403236	21.78	ng	99
49) Dimethylphthalate	13.80	163	292145	22.37	ng	99
50) 2,6-Dinitrotoluene	13.92	165	70128	22.21	ng	98
51) Acenaphthene	14.40	154	242375	21.91	ng	98
52) 3-Nitroaniline	14.25	138	46834	11.61	ng	97
53) 2,4-Dinitrophenol	14.46	184	21410	18.77	ng	93
54) Dibenzofuran	14.74	168	354725	23.11	ng	98
55) 4-Nitrophenol	14.56	139	75103	22.53	ng	98
56) 2,4-Dinitrotoluene	14.70	165	98572	22.92	ng	97
57) Fluorene	15.39	166	307974	22.74	ng	98
58) 2,3,4,6-Tetrachlorophenol	14.96	232	57385	23.88	ng	97
59) Diethylphthalate	15.16	149	314668	22.64	ng	98
60) 4-Chlorophenyl-phenylether	15.39	204	126688	22.87	ng	96
61) 4-Nitroaniline	15.41	138	98804	21.58	ng	97
62) Azobenzene	15.67	77	334367	22.60	ng	98
64) 4,6-Dinitro-2-methylphenol	15.47	198	44279	18.85	ng	98
65) n-Nitrosodiphenylamine	15.60	169	277208	21.23	ng	98
66) 4-Bromophenyl-phenylether	16.27	248	66344	21.51	ng	99
67) Hexachlorobenzene	16.39	284	70026	21.85	ng	94
68) Atrazine	16.55	200	87901	24.15	ng	99
69) Pentachlorophenol	16.73	266	40849	20.87	ng	96
70) Phenanthrene	17.12	178	474690	21.85	ng	99
71) Anthracene	17.21	178	487916	22.66	ng	99
72) Carbazole	17.48	167	511905	23.69	ng	99
73) Di-n-butylphthalate	18.05	149	636649	24.08	ng	98
74) Fluoranthene	19.14	202	544004	24.29	ng	95
76) Benzidine	19.32	184	87038	5.39	ng	99
77) Pyrene	19.50	202	584127	22.03	ng	99
79) Butylbenzylphthalate	20.40	149	299944	23.02	ng	98
80) Benzo(a)anthracene	21.24	228	506176	22.49	ng	97
81) 3,3'-Dichlorobenzidine	21.18	252	83145	11.23	ng	98
82) Chrysene	21.30	228	480755	22.13	ng	98
83) Bis(2-ethylhexyl)phthalate	21.17	149	399542	22.65	ng	99
84) Di-n-octyl phthalate	22.05	149	662099	23.04	ng	99
85) Indeno(1,2,3-cd)pyrene	25.79	276	492161	21.80	ng	99
87) Benzo(b)fluoranthene	22.82	252	459369	22.84	ng	98
88) Benzo(k)fluoranthene	22.87	252	440839	22.40	ng	99
89) Benzo(a)pyrene	23.41	252	436169	23.14	ng	99
90) Dibenzo(a,h)anthracene	25.80	278	404977	22.08	ng	99
91) Benzo(g,h,i)perylene	26.50	276	410424	22.40	ng	95

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

