

Data Path : Z:\HPCHEM1\BNA G\DATA\BG051115\
 Data File : BG017056.D
 Acq On : 11 May 2015 16:51
 Operator : TP/IZ
 Sample : G2176-20MS
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PS-26(3-5)MS

Quant Time: May 12 04:46:59 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG050515.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 12 04:00:00 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.75	152	52312	20.00	ng	0.00
21) Naphthalene-d8	10.55	136	213907	20.00	ng	0.00
38) Acenaphthene-d10	14.40	164	136234	20.00	ng	0.00
63) Phenanthrene-d10	17.15	188	308077	20.00	ng	0.00
75) Chrysene-d12	21.34	240	348041	20.00	ng	0.00
86) Perylene-d12	23.63	264	336761	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	218474	72.72	ng	0.00
7) Phenol-d6	6.95	99	305365	71.14	ng	0.00
23) Nitrobenzene-d5	8.91	82	195766	44.71	ng	0.00
41) 2,4,6-Tribromophenol	15.90	330	117312	85.78	ng	0.00
44) 2-Fluorobiphenyl	13.02	172	401332	44.45	ng	0.00
78) Terphenyl-d14	19.78	244	656296	44.21	ng	0.00

Target Compounds						Qvalue
2) 1,4-Dioxane	3.23	88	22597	17.95	ng	# 1
3) Pyridine	3.63	79	67177	17.26	ng	96
4) n-Nitrosodimethylamine	3.55	42	52122	22.43	ng	90
6) Aniline	7.08	93	108274	17.96	ng	98
8) 2-Chlorophenol	7.33	128	79186	22.75	ng	94
9) Benzaldehyde	6.89	77	45560	13.06	ng	98
10) Phenol	6.97	94	106591	23.16	ng	94
11) bis(2-Chloroethyl)ether	7.18	93	75320	21.16	ng	99
12) 1,3-Dichlorobenzene	7.64	146	77834	20.28	ng	96
13) 1,4-Dichlorobenzene	7.79	146	82593	21.25	ng	97
14) 1,2-Dichlorobenzene	8.11	146	75050	20.12	ng	# 88
15) Benzyl Alcohol	8.00	79	83778	21.37	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.29	45	123424	18.74	ng	98
17) 2-Methylphenol	8.22	107	71741	22.13	ng	93
18) Hexachloroethane	8.83	117	33167	20.76	ng	94
19) n-Nitroso-di-n-propylamine	8.56	70	70643	18.77	ng	97
20) 3+4-Methylphenols	8.54	107	96985	21.71	ng	91
22) Acetophenone	8.58	105	116625	22.89	ng	# 98
24) Nitrobenzene	8.96	77	99727	22.55	ng	98
25) Isophorone	9.47	82	189111	21.39	ng	96
26) 2-Nitrophenol	9.66	139	45016	25.00	ng	99
27) 2,4-Dimethylphenol	9.74	122	82124	25.20	ng	97
28) bis(2-Chloroethoxy)methane	9.96	93	96778	21.69	ng	# 95
29) 2,4-Dichlorophenol	10.21	162	74911	24.34	ng	98
30) 1,2,4-Trichlorobenzene	10.41	180	74622	22.86	ng	95
31) Naphthalene	10.60	128	238190	22.39	ng	99
33) 4-Chloroaniline	10.71	127	102843	20.93	ng	96
34) Hexachlorobutadiene	10.88	225	44748	21.90	ng	92
35) Caprolactam	11.48	113	39031	24.41	ng	88
36) 4-Chloro-3-methylphenol	11.87	107	97386	24.71	ng	97
37) 2-Methylnaphthalene	12.22	142	178074	21.97	ng	96
39) 1,2,4,5-Tetrachlorobenzene	12.59	216	83858	22.77	ng	# 100
40) Hexachlorocyclopentadiene	12.56	237	102978	43.30	ng	94
42) 2,4,6-Trichlorophenol	12.83	196	66460	25.47	ng	98

Data Path : Z:\HPCHEM1\BNA G\DATA\BG051115\
 Data File : BG017056.D
 Acq On : 11 May 2015 16:51
 Operator : TP/IZ
 Sample : G2176-20MS
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PS-26(3-5)MS

Quant Time: May 12 04:46:59 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG050515.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 12 04:00:00 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,5-Trichlorophenol	12.92	196	71283	25.73	nq	99
45) 1,1'-Biphenyl	13.23	154	230010	23.58	nq	99
46) 2-Chloronaphthalene	13.28	162	177898	23.03	nq	98
47) 2-Nitroaniline	13.49	65	72545	26.98	nq	96
48) Acenaphthylene	14.13	152	307184	22.83	nq	97
49) Dimethylphthalate	13.87	163	268168	26.36	nq	98
50) 2,6-Dinitrotoluene	13.99	165	55931	26.48	nq	94
51) Acenaphthene	14.47	154	183637	22.93	nq	97
52) 3-Nitroaniline	14.32	138	58810	24.47	nq	97
53) 2,4-Dinitrophenol	14.52	184	13280	13.75	nq	93
54) Dibenzofuran	14.80	168	279686	24.60	nq	93
55) 4-Nitrophenol	14.66	139	104347	51.21	nq	92
56) 2,4-Dinitrotoluene	14.77	165	79760	29.52	nq	91
57) Fluorene	15.46	166	239670	23.81	nq	98
58) 2,3,4,6-Tetrachlorophenol	15.04	232	61371	25.67	nq	# 77
59) Diethylphthalate	15.23	149	262017	24.40	nq	99
60) 4-Chlorophenyl-phenylether	15.45	204	121627	24.05	nq	98
61) 4-Nitroaniline	15.48	138	71941	25.82	nq	93
62) Azobenzene	15.74	77	274492	24.68	nq	95
64) 4,6-Dinitro-2-methylphenol	15.54	198	32671	21.39	nq	98
65) n-Nitrosodiphenylamine	15.67	169	216099	22.98	nq	97
66) 4-Bromophenyl-phenylether	16.34	248	74554	24.11	nq	98
67) Hexachlorobenzene	16.45	284	78827	24.17	nq	96
68) Atrazine	16.62	200	89081	26.32	nq	99
69) Pentachlorophenol	16.81	266	92817	43.84	nq	99
70) Phenanthrene	17.19	178	373148	23.54	nq	97
71) Anthracene	17.29	178	379034	23.70	nq	97
72) Carbazole	17.56	167	378538	24.89	nq	97
73) Di-n-butylphthalate	18.12	149	486591	24.68	nq	99
74) Fluoranthene	19.22	202	456751	24.83	nq	98
76) Benzidine	19.40	184	382044	32.31	nq	98
77) Pyrene	19.58	202	470812	22.92	nq	97
79) Butylbenzylphthalate	20.48	149	228273	23.38	nq	97
80) Benzo(a)anthracene	21.32	228	449493	24.07	nq	98
81) 3,3'-Dichlorobenzidine	21.26	252	128080	17.57	nq	97
82) Chrysene	21.37	228	418423	23.93	nq	98
83) Bis(2-ethylhexyl)phthalate	21.25	149	305775	22.90	nq	98
84) Di-n-octyl phthalate	22.14	149	521422	23.24	nq	# 100
85) Indeno(1,2,3-cd)pyrene	25.98	276	478304	22.95	nq	# 100
87) Benzo(b)fluoranthene	22.93	252	440912	23.50	nq	97
88) Benzo(k)fluoranthene	22.98	252	416892	23.10	nq	99
89) Benzo(a)pyrene	23.53	252	414789	23.71	nq	98
90) Dibenzo(a,h)anthracene	26.00	278	408416	22.86	nq	96
91) Benzo(g,h,i)perylene	26.70	276	389066	22.86	nq	94

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

Data Path : Z:\HPCHEM1\BNA G\DATA\BG051115\
 Data File : BG017056.D
 Acq On : 11 May 2015 16:51
 Operator : TP/IZ
 Sample : G2176-20MS
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PS-26(3-5)MS

Quant Time: May 12 04:46:59 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG050515.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 12 04:00:00 2015
 Response via : Initial Calibration

