

Data Path : Z:\HPCHEM1\BNA_G\Data\BG121317\
 Data File : BG031539.D
 Acq On : 13 Dec 2017 22:14
 Operator : SJ/JU
 Sample : I6676-06MSD
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 OK-03-121217MSD

Manual Integrations
 APPROVED

Quant Time: Dec 25 04:33:46 2017
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG121117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 11 17:25:55 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.11	152	51662	20.00	ng	0.00
21) Naphthalene-d8	10.92	136	220916	20.00	ng	0.00
38) Acenaphthene-d10	14.73	164	162736	20.00	ng	0.00
63) Phenanthrene-d10	17.47	188	471994	20.00	ng	0.00
75) Chrysene-d12	21.75	240	526876	20.00	ng	0.00
86) Perylene-d12	25.03	264	270887	20.00	ng	0.00

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System Monitoring Compounds

5) 2-Fluorophenol	5.67	112	410714	140.92	ng	0.00
7) Phenol-d6	7.28	99	523058	129.27	ng	0.01
23) Nitrobenzene-d5	9.27	82	340504	95.00	ng	0.00
41) 2,4,6-Tribromophenol	16.22	330	330725	173.47	ng	0.00
44) 2-Fluorobiphenyl	13.35	172	968477	87.36	ng	0.00
78) Terphenyl-d14	20.07	244	2171088	93.75	ng	0.00

12/15/2017 7:48:02 PM

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.52	88	45619	32.80	ng	# 81
3) Pyridine	3.95	79	150356	41.06	ng	92
4) n-Nitrosodimethylamine	3.83	42	61420	35.61	ng	# 95
6) Aniline	7.44	93	11975m	2.25	ng	
8) 2-Chlorophenol	7.68	128	143525	44.15	ng	89
9) Benzaldehyde	7.24	77	79212	33.77	ng	91
10) Phenol	7.31	94	166568	40.07	ng	93
11) bis(2-Chloroethyl)ether	7.52	93	130989	38.61	ng	89
12) 1,3-Dichlorobenzene	7.99	146	150876	39.02	ng	97
13) 1,4-Dichlorobenzene	8.14	146	155011	38.58	ng	95
14) 1,2-Dichlorobenzene	8.46	146	148420	38.78	ng	97
15) Benzyl Alcohol	8.35	79	72276	22.83	ng	95
16) 2,2'-oxybis(1-Chloropropan	8.63	45	192550	50.52	ng	90
17) 2-Methylphenol	8.56	107	122357	43.18	ng	97
18) Hexachloroethane	9.19	117	52542	38.79	ng	96
19) n-Nitroso-di-n-propylamine	8.91	70	106172	33.34	ng	# 96
20) 3+4-Methylphenols	8.89	107	169906	40.25	ng	94
22) Acetophenone	8.93	105	220348	41.58	ng	# 93
24) Nitrobenzene	9.32	77	158271	43.09	ng	92
25) Isophorone	9.84	82	303571	44.82	ng	# 92
26) 2-Nitrophenol	10.03	139	85195	46.91	ng	# 88
27) 2,4-Dimethylphenol	10.09	122	147093	53.32	ng	91
28) bis(2-Chloroethoxy)methane	10.31	93	164405	37.60	ng	95
29) 2,4-Dichlorophenol	10.58	162	166683	51.28	ng	90
30) 1,2,4-Trichlorobenzene	10.78	180	175269	42.21	ng	93
31) Naphthalene	10.97	128	441342	40.67	ng	99
32) Benzoic acid	10.26	122	70850	44.82	ng	# 82
33) 4-Chloroaniline	11.11	127	7855	1.67	ng	# 88
34) Hexachlorobutadiene	11.25	225	109345	39.70	ng	96
35) Caprolactam	11.86	113	67819m	53.14	ng	
36) 4-Chloro-3-methylphenol	12.21	107	186438	50.25	ng	85
37) 2-Methylnaphthalene	12.57	142	355668	42.33	ng	94
39) 1,2,4,5-Tetrachlorobenzene	12.94	216	232416	36.54	ng	98
40) Hexachlorocyclopentadiene	12.91	237	133853	66.79	ng	92

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.18	196	156066	48.03	ng	97
43) 2,4,5-Trichlorophenol	13.27	196	179299	51.22	ng #	91
45) 1,1'-Biphenyl	13.56	154	513592	38.42	ng	98
46) 2-Chloronaphthalene	13.61	162	403469	41.24	ng	96
47) 2-Nitroaniline	13.82	65	104172	37.89	ng #	74
48) Acenaphthylene	14.46	152	612651	38.92	ng	99
49) Dimethylphthalate	14.18	163	611230	45.37	ng	98
50) 2,6-Dinitrotoluene	14.30	165	138294	48.02	ng #	75
51) Acenaphthene	14.79	154	393803	39.56	ng	96
52) 3-Nitroaniline	14.66	138	18132	6.18	ng #	84
53) 2,4-Dinitrophenol	14.86	184	129387	104.53	ng #	50
54) Dibenzofuran	15.13	168	675487	42.53	ng	97
55) 4-Nitrophenol	14.97	139	189714	94.95	ng #	81
56) 2,4-Dinitrotoluene	15.10	165	206412	50.46	ng #	74
57) Fluorene	15.77	166	555656	43.66	ng	99
58) 2,3,4,6-Tetrachlorophenol	15.36	232	181848	55.27	ng #	88
59) Diethylphthalate	15.53	149	624245	46.22	ng	97
60) 4-Chlorophenyl-phenylether	15.76	204	323713	41.57	ng	92
61) 4-Nitroaniline	15.81	138	56553	18.36	ng	85
62) Azobenzene	16.06	77	457149	40.86	ng	93
64) 4,6-Dinitro-2-methylphenol	15.87	198	108690	39.06	ng	80
65) n-Nitrosodiphenylamine	15.98	169	71988	5.21	ng	96
66) 4-Bromophenyl-phenylether	16.65	248	226782	41.32	ng	96
67) Hexachlorobenzene	16.78	284	230088	40.60	ng	96
68) Atrazine	16.92	200	9698	1.92	ng	97
69) Pentachlorophenol	17.13	266	207527	76.45	ng	99
70) Phenanthrene	17.52	178	1033433	41.92	ng	98
71) Anthracene	17.61	178	1060951	43.52	ng	99
72) Carbazole	17.88	167	172952	7.84	ng	98
73) Di-n-butylphthalate	18.41	149	1176493	46.18	ng	98
74) Fluoranthene	19.52	202	1376386	44.62	ng	97
77) Pyrene	19.88	202	1376641	42.05	ng	97
79) Butylbenzylphthalate	20.74	149	527989	45.97	ng #	83
80) Benzo(a)anthracene	21.72	228	1347513	42.37	ng	97
82) Chrysene	21.80	228	1310249	42.98	ng	98
83) Bis(2-ethylhexyl)phthalate	21.60	149	745154	45.10	ng	99
84) Di-n-octyl phthalate	22.82	149	1254027	46.01	ng #	94
85) Indeno(1,2,3-cd)pyrene	28.81	276	1143766	35.10	ng	99
87) Benzo(b)fluoranthene	23.99	252	1300387	80.30	ng	98
88) Benzo(k)fluoranthene	24.05	252	1239103	74.97	ng	97
89) Benzo(a)pyrene	24.88	252	975860	63.52	ng	98
90) Dibenzo(a,h)anthracene	28.85	278	1111147	75.63	ng	98
91) Benzo(g,h,i)perylene	29.99	276	837610	57.99	ng	98

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

