

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG070117\
 Data File : BG027802.D
 Acq On : 2 Jul 2017 3:20
 Operator : SJ/JU
 Sample : I4007-05MSD
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP2-MH2094-COMPMSD

Quant Time: Jul 03 07:42:16 2017
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG062817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 28 19:26:37 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.48	152	31078	20.00	ng	0.00
21) Naphthalene-d8	11.30	136	139488	20.00	ng	0.00
38) Acenaphthene-d10	15.06	164	8575	20.00	ng	0.00
63) Phenanthrene-d10	17.81	188	59641	20.00	ng	0.00
75) Chrysene-d12	22.17	240	70	20.00	ng	0.00
86) Perylene-d12	0.00	264	0	0.00	ng	-25.79

System Monitoring Compounds

5) 2-Fluorophenol	6.06	112	252504	125.12	ng	0.00
7) Phenol-d6	7.62	99	68817	22.32	ng	0.00
23) Nitrobenzene-d5	9.65	82	236962	94.69	ng	0.00
41) 2,4,6-Tribromophenol	16.55	330	193131	1642.05	ng	0.00
44) 2-Fluorobiphenyl	13.69	172	574998	957.66	ng	0.00
78) Terphenyl-d14	20.39	244	13359	4474.06	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	4.06	88	27358	36.410	ng	96
4) n-Nitrosodimethylamine	4.35	42	5829	5.138	ng	# 86
6) Aniline	7.89	93	85997	24.063	ng	# 57
8) 2-Chlorophenol	8.05	128	100556	44.589	ng	88
9) Benzaldehyde	7.62	77	32290	17.600	ng	# 71
10) Phenol	7.65	94	45326	14.859	ng	88
11) bis(2-Chloroethyl)ether	7.89	93	85997	37.116	ng	93
12) 1,3-Dichlorobenzene	8.37	146	91970	38.340	ng	# 92
13) 1,4-Dichlorobenzene	8.52	146	96593	38.861	ng	96
14) 1,2-Dichlorobenzene	8.83	146	97434	40.429	ng	92
16) 2,2'-oxybis(1-Chloropropan	8.99	45	183521	37.135	ng	98
17) 2-Methylphenol	8.90	107	24322	11.270	ng	# 87
18) Hexachloroethane	9.57	117	32339	36.988	ng	85
20) 3+4-Methylphenols	9.22	107	30776	9.889	ng	86
22) Acetophenone	9.29	105	172114	51.134	ng	# 90
24) Nitrobenzene	9.69	77	119187	45.557	ng	# 84
25) Isophorone	10.20	82	11545	2.117	ng	# 91
26) 2-Nitrophenol	10.40	139	57920	49.013	ng	# 84
27) 2,4-Dimethylphenol	10.43	122	36932	16.686	ng	87
29) 2,4-Dichlorophenol	10.93	162	104541	53.895	ng	97
30) 1,2,4-Trichlorobenzene	11.16	180	94849	47.945	ng	98
31) Naphthalene	11.35	128	329598	44.433	ng	99
32) Benzoic acid	10.56	122	64620	47.809	ng	# 75
34) Hexachlorobutadiene	11.62	225	56756	51.489	ng	96
36) 4-Chloro-3-methylphenol	12.52	107	70670	25.891	ng	88
37) 2-Methylnaphthalene	12.91	142	231372	41.877	ng	96
39) 1,2,4,5-Tetrachlorobenzene	13.28	216	115317	511.762	ng	98
40) Hexachlorocyclopentadiene	13.25	237	154047	1449.977	ng	95
42) 2,4,6-Trichlorophenol	13.51	196	78967	485.217	ng	93
43) 2,4,5-Trichlorophenol	13.58	196	83715	453.973	ng	# 95
45) 1,1'-Biphenyl	13.90	154	326547	475.109	ng	98
46) 2-Chloronaphthalene	13.95	162	244140	470.213	ng	98
48) Acenaphthylene	14.79	152	10829	11.414	ng	93
49) Dimethylphthalate	14.50	163	28102	38.417	ng	# 96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
50) 2,6-Dinitrotoluene	14.63	165	73997	465.340	ng	# 83
51) Acenaphthene	15.13	154	33800	55.585	ng	97
53) 2,4-Dinitrophenol	15.17	184	95384	984.404	ng	94
54) Dibenzofuran	15.46	168	338824	421.399	ng	98
55) 4-Nitrophenol	15.30	139	50522	353.784	ng	# 60
56) 2,4-Dinitrotoluene	15.41	165	106884	481.905	ng	# 87
57) Fluorene	16.11	166	168465	216.094	ng	99
58) 2,3,4,6-Tetrachlorophenol	15.69	232	77056	484.739	ng	95
60) 4-Chlorophenyl-phenylether	16.09	204	138808	402.214	ng	95
64) 4,6-Dinitro-2-methylphenol	16.17	198	59815	146.678	ng	84
66) 4-Bromophenyl-phenylether	16.99	248	85976	137.829	ng	92
67) Hexachlorobenzene	17.12	284	119489	180.875	ng	90
69) Pentachlorophenol	17.46	266	117595	263.482	ng	97
70) Phenanthrene	17.86	178	378575	109.952	ng	94
71) Anthracene	17.95	178	92365	26.566	ng	94
74) Fluoranthene	19.84	202	29438	7.823	ng	95

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

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