

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071017\
 Data File : BG027896.D
 Acq On : 10 Jul 2017 20:47
 Operator : SJ/JU
 Sample : I4071-03MSD
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 HR-06-070617-AMSD

Manual Integrations
 APPROVED

mohammad
 7/11/2017 3:34:43 PM

Quant Time: Jul 11 13:21:21 2017
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG062817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 07 17:43:21 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.42	152	39703	20.00	ng	0.00
21) Naphthalene-d8	11.23	136	186821	20.00	ng	-0.01
38) Acenaphthene-d10	15.02	164	124351	20.00	ng	0.00
63) Phenanthrene-d10	17.76	188	313786	20.00	ng	-0.01
75) Chrysene-d12	22.11	240	334038	20.00	ng	-0.01
86) Perylene-d12	25.67	264	297204	20.00	ng	-0.03

mohammad

System Monitoring Compounds

5) 2-Fluorophenol	6.00	112	317372	123.09	ng	0.00
7) Phenol-d6	7.57	99	421088	106.91	ng	0.00
23) Nitrobenzene-d5	9.58	82	247113	73.73	ng	-0.01
41) 2,4,6-Tribromophenol	16.50	330	291189	170.72	ng	-0.01
44) 2-Fluorobiphenyl	13.64	172	716740	82.32	ng	-0.01
78) Terphenyl-d14	20.35	244	1209718	84.90	ng	0.00

7/11/2017 3:34:43 PM

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.97	88	32093	33.433	ng	# 70
3) Pyridine	4.45	79	59969m	20.289	ng	
4) n-Nitrosodimethylamine	4.26	42	28473	19.644	ng	# 1
8) 2-Chlorophenol	7.99	128	122208	42.417	ng	# 79
9) Benzaldehyde	7.56	77	62235	26.553	ng	# 77
10) Phenol	7.60	94	145263	37.277	ng	# 65
11) bis(2-Chloroethyl)ether	7.82	93	172623	58.318	ng	# 72
12) 1,3-Dichlorobenzene	8.31	146	102707	33.515	ng	# 86
13) 1,4-Dichlorobenzene	8.45	146	107567	33.875	ng	95
14) 1,2-Dichlorobenzene	8.78	146	105804	34.365	ng	92
15) Benzyl Alcohol	8.65	79	75600	27.746	ng	# 77
16) 2,2'-oxybis(1-Chloropropan	8.93	45	128221	20.309	ng	83
17) 2-Methylphenol	8.85	107	108446	39.335	ng	# 81
18) Hexachloroethane	9.50	117	35373	31.669	ng	80
19) n-Nitroso-di-n-propylamine	9.21	70	91360	31.882	ng	# 69
20) 3+4-Methylphenols	9.18	107	151768	38.173	ng	84
22) Acetophenone	9.23	105	179376	39.789	ng	# 72
24) Nitrobenzene	9.63	77	123000	35.103	ng	# 76
25) Isophorone	10.14	82	271549	37.170	ng	# 91
26) 2-Nitrophenol	10.34	139	65427	41.338	ng	# 61
27) 2,4-Dimethylphenol	10.38	122	137376	46.341	ng	87
28) bis(2-Chloroethoxy)methane	10.61	93	164134	39.628	ng	97
29) 2,4-Dichlorophenol	10.88	162	121096	46.612	ng	97
30) 1,2,4-Trichlorobenzene	11.09	180	102342	38.626	ng	98
31) Naphthalene	11.29	128	383121	38.563	ng	98
32) Benzoic acid	10.52	122	64904	38.069	ng	# 73
33) 4-Chloroaniline	11.46	127	94722	22.069	ng	# 82
34) Hexachlorobutadiene	11.56	225	50882	34.465	ng	97
35) Caprolactam	12.18	113	45787m	35.040	ng	
36) 4-Chloro-3-methylphenol	12.48	107	156604	42.839	ng	# 84
37) 2-Methylnaphthalene	12.86	142	303987	41.080	ng	# 89
39) 1,2,4,5-Tetrachlorobenzene	13.22	216	132030	40.405	ng	100
40) Hexachlorocyclopentadiene	13.19	237	140850	91.422	ng	95
42) 2,4,6-Trichlorophenol	13.45	196	105302	44.618	ng	96

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071017\
 Data File : BG027896.D
 Acq On : 10 Jul 2017 20:47
 Operator : SJ/JU
 Sample : I4071-03MSD
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 HR-06-070617-AMSD

Manual Integrations
 APPROVED

mohammad
 7/11/2017 3:34:43 PM

Quant Time: Jul 11 13:21:21 2017
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG062817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 07 17:43:21 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,5-Trichlorophenol	13.54	196	118866	44.450	ng	92
45) 1,1'-Biphenyl	13.85	154	409271	41.062	ng	98
46) 2-Chloronaphthalene	13.90	162	299633	39.795	ng	97
47) 2-Nitroaniline	14.10	65	100465	33.335	ng	# 51
48) Acenaphthylene	14.74	152	547170	39.770	ng	97
49) Dimethylphthalate	14.45	163	455434	42.933	ng	98 mohammad
50) 2,6-Dinitrotoluene	14.58	165	102865	44.608	ng	# 60
51) Acenaphthene	15.08	154	338192	38.352	ng	96
52) 3-Nitroaniline	14.92	138	88967	32.794	ng	# 73
53) 2,4-Dinitrophenol	15.12	184	125612	97.191	ng	# 70
54) Dibenzofuran	15.41	168	534310	45.824	ng	90 7/11/2017 3:34:43 PM
55) 4-Nitrophenol	15.22	139	156459	75.552	ng	# 37
56) 2,4-Dinitrotoluene	15.37	165	148967	46.315	ng	# 72
57) Fluorene	16.06	166	489082	43.261	ng	99
58) 2,3,4,6-Tetrachlorophenol	15.63	232	116594	50.578	ng	# 98
59) Diethylphthalate	15.80	149	493449	42.009	ng	97
60) 4-Chlorophenyl-phenylether	16.04	204	216693	43.298	ng	92
61) 4-Nitroaniline	16.09	138	114686	40.578	ng	# 49
62) Azobenzene	16.33	77	421384	40.825	ng	80
64) 4,6-Dinitro-2-methylphenol	16.13	198	80010	41.758	ng	83
65) n-Nitrosodiphenylamine	16.26	169	432049	42.327	ng	99
66) 4-Bromophenyl-phenylether	16.94	248	149007	45.403	ng	93
67) Hexachlorobenzene	17.07	284	168169	48.385	ng	# 86
68) Atrazine	17.20	200	104602	32.458	ng	93
69) Pentachlorophenol	17.41	266	191789	85.933	ng	96
70) Phenanthrene	17.81	178	788426	43.524	ng	95
71) Anthracene	17.90	178	783828	42.850	ng	97
72) Carbazole	18.17	167	657073	36.473	ng	95
73) Di-n-butylphthalate	18.69	149	768042	38.795	ng	# 92
74) Fluoranthene	19.80	202	785147	39.658	ng	92
76) Benzidine	20.01	184	4301	3.722	ng	96
77) Pyrene	20.16	202	801082	38.214	ng	96
79) Butylbenzylphthalate	21.03	149	348505	36.912	ng	85
80) Benzo(a)anthracene	22.09	228	792749	39.254	ng	95
81) 3,3'-Dichlorobenzidine	21.99	252	180946	24.686	ng	# 98
82) Chrysene	22.16	228	745904	39.404	ng	95
83) Bis(2-ethylhexyl)phthalate	21.94	149	503270	36.476	ng	100
84) Di-n-octyl phthalate	23.27	149	866816	38.027	ng	# 82
85) Indeno(1,2,3-cd)pyrene	29.77	276	945703	43.711	ng	# 86
87) Benzo(b)fluoranthene	24.53	252	808790	42.240	ng	# 94
88) Benzo(k)fluoranthene	24.61	252	799876	42.261	ng	# 92
89) Benzo(a)pyrene	25.50	252	764525	42.552	ng	# 91
90) Dibenzo(a,h)anthracene	29.84	278	815431	44.800	ng	# 92
91) Benzo(g,h,i)perylene	31.06	276	758094	43.338	ng	# 87

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

