

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG070117\
 Data File : BG027809.D
 Acq On : 2 Jul 2017 7:58
 Operator : SJ/JU
 Sample : I3979-05MSD
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP3-MH2095-COMPMSD

Manual Integrations
 APPROVED

mohammad
 7/5/2017 11:32:00 AM

Quant Time: Jul 03 16:07:36 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	29189	20.00	ng	0.00
21) Naphthalene-d8	11.30	136	147327	20.00	ng	0.00
38) Acenaphthene-d10	15.06	164	98900	20.00	ng	0.00
63) Phenanthrene-d10	17.81	188	243400	20.00	ng	0.00
75) Chrysene-d12	22.18	240	256143	20.00	ng	0.00
86) Perylene-d12	25.78	264	232211	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	6.06	112	258020	136.12	ng	0.00
7) Phenol-d6	7.62	99	352846	121.85	ng	0.00
23) Nitrobenzene-d5	9.64	82	243175	92.00	ng	0.00
41) 2,4,6-Tribromophenol	16.55	330	236060	174.02	ng	0.00
44) 2-Fluorobiphenyl	13.69	172	660018	95.31	ng	0.00
78) Terphenyl-d14	20.39	244	1114397	102.00	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	4.06	88	20446	28.972	ng	# 90
3) Pyridine	4.96	79	30494m	14.033	ng	
4) n-Nitrosodimethylamine	4.35	42	40747	38.237	ng	88
8) 2-Chlorophenol	8.05	128	99922	47.175	ng	90
9) Benzaldehyde	7.62	77	27593	16.014	ng	85
10) Phenol	7.65	94	116687	40.729	ng	84
11) bis(2-Chloroethyl)ether	7.88	93	81063	37.250	ng	98
12) 1,3-Dichlorobenzene	8.36	146	88422	39.246	ng	# 92
13) 1,4-Dichlorobenzene	8.51	146	89313	38.258	ng	95
14) 1,2-Dichlorobenzene	8.83	146	89693	39.625	ng	93
15) Benzyl Alcohol	8.71	79	11378	5.680	ng	88
16) 2,2'-oxybis(1-Chloropropan	8.99	45	176495	38.024	ng	98
17) 2-Methylphenol	8.90	107	90348	44.575	ng	# 87
18) Hexachloroethane	9.57	117	30254	36.842	ng	# 83
19) n-Nitroso-di-n-propylamine	9.27	70	81221	38.553	ng	# 97
20) 3+4-Methylphenols	9.23	107	123211	42.153	ng	89
22) Acetophenone	9.29	105	150264	42.267	ng	# 91
24) Nitrobenzene	9.69	77	112241	40.619	ng	# 83
25) Isophorone	10.20	82	235216	40.828	ng	# 92
26) 2-Nitrophenol	10.40	139	56165	44.999	ng	# 82
27) 2,4-Dimethylphenol	10.44	122	105728	45.226	ng	91
28) bis(2-Chloroethoxy)methane	10.67	93	126721	38.797	ng	# 94
29) 2,4-Dichlorophenol	10.94	162	107034	52.244	ng	97
30) 1,2,4-Trichlorobenzene	11.15	180	92424	44.233	ng	97
31) Naphthalene	11.34	128	327873	41.848	ng	98
32) Benzoic acid	10.56	122	23295	22.155	ng	# 77
33) 4-Chloroaniline	11.73	127	16676m	4.927	ng	
34) Hexachlorobutadiene	11.61	225	54638	46.930	ng	100
35) Caprolactam	12.25	113	37297m	36.195	ng	
36) 4-Chloro-3-methylphenol	12.55	107	115973	40.228	ng	92
37) 2-Methylnaphthalene	12.92	142	265723	45.535	ng	95
39) 1,2,4,5-Tetrachlorobenzene	13.28	216	121423	46.721	ng	98
40) Hexachlorocyclopentadiene	13.25	237	150387	122.732	ng	96
42) 2,4,6-Trichlorophenol	13.51	196	94635	50.417	ng	91

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,5-Trichlorophenol	13.59	196	101587	47.764	ng	# 95
45) 1,1'-Biphenyl	13.90	154	354423	44.710	ng	97
46) 2-Chloronaphthalene	13.95	162	266971	44.582	ng	98
47) 2-Nitroaniline	14.15	65	105734	44.111	ng	# 81
48) Acenaphthylene	14.79	152	461521	42.177	ng	96
49) Dimethylphthalate	14.50	163	387656	45.948	ng	97
50) 2,6-Dinitrotoluene	14.63	165	85669	46.711	ng	# 78
51) Acenaphthene	15.13	154	285111	40.653	ng	94
52) 3-Nitroaniline	15.00	138	49921	23.137	ng	# 78
53) 2,4-Dinitrophenol	15.17	184	109717	105.896	ng	94
54) Dibenzofuran	15.46	168	448942	48.411	ng	98
55) 4-Nitrophenol	15.28	139	107910	65.517	ng	# 63
56) 2,4-Dinitrotoluene	15.41	165	125294	48.980	ng	# 89
57) Fluorene	16.11	166	394291	43.852	ng	99
58) 2,3,4,6-Tetrachlorophenol	15.68	232	89371m	48.746	ng	
59) Diethylphthalate	15.86	149	429960	46.024	ng	96
60) 4-Chlorophenyl-phenylether	16.08	204	191490	48.109	ng	93
61) 4-Nitroaniline	16.14	138	62379	27.751	ng	# 72
62) Azobenzene	16.38	77	365222	44.490	ng	94
64) 4,6-Dinitro-2-methylphenol	16.18	198	72715	47.897	ng	81
65) n-Nitrosodiphenylamine	16.31	169	341494	43.130	ng	97
66) 4-Bromophenyl-phenylether	16.98	248	131876	51.803	ng	94
67) Hexachlorobenzene	17.11	284	135345	50.202	ng	98
68) Atrazine	17.29	200	13795	5.519	ng	97
69) Pentachlorophenol	17.46	266	154665	89.094	ng	97
70) Phenanthrene	17.86	178	639002	45.476	ng	95
71) Anthracene	17.95	178	633224	44.627	ng	98
72) Carbazole	18.22	167	516687	36.974	ng	96
73) Di-n-butylphthalate	18.73	149	712518	46.399	ng	# 93
74) Fluoranthene	19.85	202	694427	45.219	ng	96
76) Benzidine	19.85	184	58	3.226	ng	# 1
77) Pyrene	20.21	202	720145	44.800	ng	94
79) Butylbenzylphthalate	21.08	149	322416	44.534	ng	# 88
80) Benzo(a)anthracene	22.15	228	696387	44.969	ng	94
81) 3,3'-Dichlorobenzidine	22.10	252	19447m	3.460	ng	
82) Chrysene	22.23	228	651003	44.849	ng	95
83) Bis(2-ethylhexyl)phthalate	22.00	149	459174	43.401	ng	# 96
84) Di-n-octyl phthalate	23.35	149	772172	44.177	ng	# 90
85) Indeno(1,2,3-cd)pyrene	29.94	276	754442	45.476	ng	# 66
87) Benzo(b)fluoranthene	24.63	252	693120	46.330	ng	# 97
88) Benzo(k)fluoranthene	24.71	252	684645	46.298	ng	# 93
89) Benzo(a)pyrene	25.62	252	652155	46.457	ng	# 94
90) Dibenzo(a,h)anthracene	30.02	278	625362	43.974	ng	# 95
91) Benzo(g,h,i)perylene	31.25	276	589180	43.109	ng	# 93

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

