

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG041416\
 Data File : BG021668.D
 Acq On : 15 Apr 2016 1:11
 Operator : UM/SJ
 Sample : H2517-05MS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 S8-0572MS

Manual Integrations
 APPROVED

Sohil
 4/15/2016 4:44:52 PM

Quant Time: Apr 15 04:09:15 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG041316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 13 15:48:13 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.21	152	30058	20.00	ng	0.00
21) Naphthalene-d8	11.03	136	140719	20.00	ng	0.00
38) Acenaphthene-d10	14.82	164	93604	20.00	ng	0.00
63) Phenanthrene-d10	17.56	188	222058	20.00	ng	0.00
75) Chrysene-d12	21.83	240	252202	20.00	ng	0.00
86) Perylene-d12	25.16	264	249873	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.76	112	135878	75.28	ng	0.00
7) Phenol-d6	7.36	99	126275	46.83	ng	0.00
23) Nitrobenzene-d5	9.38	82	281065	102.66	ng	0.00
41) 2,4,6-Tribromophenol	16.31	330	179755	178.18	ng	0.00
44) 2-Fluorobiphenyl	13.46	172	645271	101.00	ng	0.00
78) Terphenyl-d14	20.14	244	938242	92.83	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.63	88	14664	18.25	ng	87
3) Pyridine	4.04	79	36570	15.17	ng	89
4) n-Nitrosodimethylamine	3.95	42	21868	18.30	ng	# 79
6) Aniline	7.54	93	38793	10.88	ng	94
8) 2-Chlorophenol	7.78	128	90520	41.83	ng	96
9) Benzaldehyde	7.35	77	34875	22.37	ng	92
10) Phenol	7.39	94	44813	15.45	ng	99
11) bis(2-Chloroethyl)ether	7.63	93	106921	46.07	ng	95
12) 1,3-Dichlorobenzene	8.10	146	96647	39.99	ng	95
13) 1,4-Dichlorobenzene	8.25	146	99993	40.64	ng	96
14) 1,2-Dichlorobenzene	8.57	146	98216	41.61	ng	97
15) Benzyl Alcohol	8.45	79	69553	33.76	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.75	45	211278	42.50	ng	99
17) 2-Methylphenol	8.65	107	70636	35.23	ng	91
18) Hexachloroethane	9.31	117	35751	39.93	ng	# 84
19) n-Nitroso-di-n-propylamine	9.02	70	98710	47.12	ng	94
20) 3+4-Methylphenols	8.98	107	85728	31.25	ng	92
22) Acetophenone	9.04	105	176126	44.94	ng	# 96
24) Nitrobenzene	9.42	77	138792	47.35	ng	97
25) Isophorone	9.95	82	272207	45.43	ng	98
26) 2-Nitrophenol	10.14	139	59066	52.78	ng	92
27) 2,4-Dimethylphenol	10.19	122	105262	41.41	ng	98
28) bis(2-Chloroethoxy)methane	10.42	93	160022	50.26	ng	91
29) 2,4-Dichlorophenol	10.67	162	110910	51.17	ng	97
30) 1,2,4-Trichlorobenzene	10.89	180	105437	45.24	ng	92
31) Naphthalene	11.08	128	345970	45.94	ng	# 95
32) Benzoic acid	10.26	122	21832m	20.15	ng	
33) 4-Chloroaniline	11.19	127	24881	7.63	ng	98
34) Hexachlorobutadiene	11.36	225	62387	41.50	ng	95
35) Caprolactam	11.96	113	6782	6.84	ng	# 86
36) 4-Chloro-3-methylphenol	12.29	107	127368	46.73	ng	98
37) 2-Methylnaphthalene	12.67	142	268663	51.97	ng	98
39) 1,2,4,5-Tetrachlorobenzene	13.03	216	134364	45.93	ng	98
40) Hexachlorocyclopentadiene	13.01	237	152522	93.86	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.26	196	96734	51.86	ng	97
43) 2,4,5-Trichlorophenol	13.33	196	106257	52.79	ng	98
45) 1,1'-Biphenyl	13.66	154	360109	45.66	ng	97
46) 2-Chloronaphthalene	13.71	162	277987	47.68	ng	99
47) 2-Nitroaniline	13.90	65	105782	52.91	ng	97
48) Acenaphthylene	14.55	152	465935	48.33	ng	97
49) Dimethylphthalate	14.27	163	393040	48.56	ng	98
50) 2,6-Dinitrotoluene	14.39	165	84845	54.98	ng	98
51) Acenaphthene	14.89	154	330552	53.64	ng	98
52) 3-Nitroaniline	14.72	138	19278	11.26	ng	100
53) 2,4-Dinitrophenol	14.92	184	70680	115.42	ng	97
54) Dibenzofuran	15.22	168	459151	54.57	ng	98
55) 4-Nitrophenol	15.01	139	53662	36.62	ng	94
56) 2,4-Dinitrotoluene	15.17	165	121709	58.36	ng	98
57) Fluorene	15.86	166	375419	49.96	ng	99
58) 2,3,4,6-Tetrachlorophenol	15.44	232	99832	59.78	ng	98
59) Diethylphthalate	15.62	149	404785	47.92	ng	97
60) 4-Chlorophenyl-phenylether	15.85	204	183475	49.56	ng	99
61) 4-Nitroaniline	15.88	138	85618	45.46	ng	98
62) Azobenzene	16.15	77	392659	54.24	ng	99
64) 4,6-Dinitro-2-methylphenol	15.93	198	50896	54.24	ng	91
65) n-Nitrosodiphenylamine	16.06	169	336298	50.50	ng	96
66) 4-Bromophenyl-phenylether	16.75	248	121892	51.24	ng	95
67) Hexachlorobenzene	16.86	284	126117	50.21	ng	96
68) Atrazine	17.00	200	131256	49.91	ng	98
69) Pentachlorophenol	17.20	266	159129	110.98	ng	96
70) Phenanthrene	17.60	178	610439	49.88	ng	98
71) Anthracene	17.69	178	618164	49.76	ng	97
72) Carbazole	17.96	167	576546	49.72	ng	97
73) Di-n-butylphthalate	18.50	149	710769	48.19	ng	98
74) Fluoranthene	19.60	202	702228	48.05	ng	99
76) Benzidine	19.77	184	163999	20.89	ng	99
77) Pyrene	19.95	202	723076	49.72	ng	99
79) Butylbenzylphthalate	20.82	149	337954	50.06	ng	98
80) Benzo(a)anthracene	21.81	228	725669	50.00	ng	97
81) 3,3'-Dichlorobenzidine	21.72	252	84139	7.32	ng	95
82) Chrysene	21.88	228	670927	48.12	ng	97
83) Bis(2-ethylhexyl)phthalate	21.70	149	488684	49.49	ng	99
84) Di-n-octyl phthalate	22.95	149	845843	51.14	ng	100
85) Indeno(1,2,3-cd)pyrene	28.97	276	834114	50.32	ng	98
87) Benzo(b)fluoranthene	24.10	252	739453	51.00	ng	99
88) Benzo(k)fluoranthene	24.17	252	696901	49.16	ng	99
89) Benzo(a)pyrene	25.01	252	701737	49.95	ng	98
90) Dibenzo(a,h)anthracene	29.05	278	701377	49.80	ng	98
91) Benzo(g,h,i)perylene	30.18	276	696837	50.42	ng	97

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

