

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG061215\
 Data File : BG017664.D
 Acq On : 13 Jun 2015 8:23
 Operator : TP/IZ
 Sample : G2522-03MSD
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-19MSD

Manual Integrations
 APPROVED

apatel
 6/15/2015 4:40:26 PM

Quant Time: Jun 15 04:44:53 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG061215.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jun 13 02:22:44 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.52	152	19322	20.00	ng	0.00
21) Naphthalene-d8	10.32	136	73085	20.00	ng	0.00
38) Acenaphthene-d10	14.20	164	58219	20.00	ng	0.00
63) Phenanthrene-d10	16.97	188	159071	20.00	ng	0.00
75) Chrysene-d12	21.17	240	235659	20.00	ng	0.00
86) Perylene-d12	23.36	264	223581	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.11	112	82796	82.99	ng	0.00
7) Phenol-d6	6.74	99	76365	52.24	ng	0.00
23) Nitrobenzene-d5	8.69	82	178828	102.71	ng	0.00
41) 2,4,6-Tribromophenol	15.71	330	165245	164.43	ng	0.00
44) 2-Fluorobiphenyl	12.82	172	415711	98.76	ng	0.00
78) Terphenyl-d14	19.61	244	1072527	94.94	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.95	88	10729	21.59	ng	# 90
3) Pyridine	3.35	79	22904	17.81	ng	93
4) n-Nitrosodimethylamine	3.28	42	15674	21.19	ng	# 95
6) Aniline	6.86	93	37130	19.07	ng	# 85
8) 2-Chlorophenol	7.10	128	51009	42.12	ng	94
9) Benzaldehyde	6.67	77	3910	3.96	ng	92
10) Phenol	6.77	94	28969	19.05	ng	98
11) bis(2-Chloroethyl)ether	6.96	93	52944	46.96	ng	93
12) 1,3-Dichlorobenzene	7.41	146	57977m	40.98	ng	
13) 1,4-Dichlorobenzene	7.56	146	61680	41.39	ng	92
14) 1,2-Dichlorobenzene	7.88	146	58546	42.27	ng	93
15) Benzyl Alcohol	7.78	79	51363	31.88	ng	90
16) 2,2'-oxybis(1-Chloropropan	8.07	45	21436	47.46	ng	92
17) 2-Methylphenol	8.01	107	39094	35.48	ng	95
18) Hexachloroethane	8.59	117	25633	41.42	ng	# 83
19) n-Nitroso-di-n-propylamine	8.34	70	60112	45.84	ng	# 95
20) 3+4-Methylphenols	8.33	107	50127	33.05	ng	# 87
22) Acetophenone	8.36	105	100357	52.10	ng	# 94
24) Nitrobenzene	8.73	77	105854	56.13	ng	97
25) Isophorone	9.26	82	154317	47.19	ng	95
26) 2-Nitrophenol	9.44	139	32776	45.22	ng	92
27) 2,4-Dimethylphenol	9.53	122	52838	45.31	ng	# 95
28) bis(2-Chloroethoxy)methane	9.74	93	72233	50.99	ng	92
29) 2,4-Dichlorophenol	9.99	162	67554	54.33	ng	96
30) 1,2,4-Trichlorobenzene	10.19	180	77114	47.98	ng	96
31) Naphthalene	10.37	128	189143	47.95	ng	# 96
32) Benzoic acid	9.68	122	7444	10.74	ng	# 63
33) 4-Chloroaniline	10.50	127	29638	18.35	ng	98
34) Hexachlorobutadiene	10.66	225	61342	44.49	ng	97
35) Caprolactam	11.27	113	5029m	8.81	ng	
36) 4-Chloro-3-methylphenol	11.65	107	78211	49.19	ng	98
37) 2-Methylnaphthalene	12.00	142	151942	48.18	ng	97
39) 1,2,4,5-Tetrachlorobenzene	12.38	216	115644	49.34	ng	97
40) Hexachlorocyclopentadiene	12.35	237	89149	84.40	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	12.64	196	70407	48.44	ng	99
43) 2,4,5-Trichlorophenol	12.73	196	75990	49.98	ng	91
45) 1,1'-Biphenyl	13.03	154	216116	50.62	ng	98
46) 2-Chloronaphthalene	13.07	162	169610	47.88	ng	96
47) 2-Nitroaniline	13.29	65	64063	51.22	ng	95
48) Acenaphthylene	13.93	152	279178	46.37	ng	99
49) Dimethylphthalate	13.68	163	256978	51.55	ng	97
50) 2,6-Dinitrotoluene	13.80	165	58529	53.54	ng	95
51) Acenaphthene	14.27	154	171445	47.95	ng	98
52) 3-Nitroaniline	14.13	138	19694	20.16	ng	# 95
53) 2,4-Dinitrophenol	14.35	184	76026	100.99	ng	94
54) Dibenzofuran	14.61	168	303108	51.34	ng	97
55) 4-Nitrophenol	14.50	139	27162	43.59	ng	92
56) 2,4-Dinitrotoluene	14.60	165	82767	52.17	ng	97
57) Fluorene	15.26	166	256411	50.62	ng	98
58) 2,3,4,6-Tetrachlorophenol	14.85	232	82186	49.64	ng	95
59) Diethylphthalate	15.05	149	273995	49.53	ng	98
60) 4-Chlorophenyl-phenylether	15.26	204	170324	53.70	ng	98
61) 4-Nitroaniline	15.30	138	33619	32.29	ng	# 78
62) Azobenzene	15.56	77	301864	50.28	ng	98
64) 4,6-Dinitro-2-methylphenol	15.37	198	65049	50.90	ng	93
65) n-Nitrosodiphenylamine	15.48	169	238606	51.48	ng	97
66) 4-Bromophenyl-phenylether	16.15	248	113973	53.87	ng	97
67) Hexachlorobenzene	16.27	284	117653	51.10	ng	93
68) Atrazine	16.44	200	114846	50.25	ng	100
69) Pentachlorophenol	16.62	266	116879	107.81	ng	94
70) Phenanthrene	17.01	178	439068	49.46	ng	98
71) Anthracene	17.09	178	453330	51.43	ng	97
72) Carbazole	17.38	167	420764	52.06	ng	100
73) Di-n-butylphthalate	17.94	149	531873	53.77	ng	98
74) Fluoranthene	19.03	202	663191	52.21	ng	99
76) Benzidine	19.23	184	80674	14.65	ng	98
77) Pyrene	19.40	202	680214	51.12	ng	99
79) Butylbenzylphthalate	20.31	149	258671	53.85	ng	98
80) Benzo(a)anthracene	21.15	228	729626	50.19	ng	98
81) 3,3'-Dichlorobenzidine	21.09	252	99263	19.08	ng	# 96
82) Chrysene	21.20	228	686671	52.37	ng	100
83) Bis(2-ethylhexyl)phthalate	21.08	149	370474	52.50	ng	97
84) Di-n-octyl phthalate	21.95	149	607655	51.69	ng	99
85) Indeno(1,2,3-cd)pyrene	25.60	276	834205	48.92	ng	98
87) Benzo(b)fluoranthene	22.71	252	726743	50.43	ng	100
88) Benzo(k)fluoranthene	22.75	252	704422	50.81	ng	99
89) Benzo(a)pyrene	23.27	252	696940	52.36	ng	99
90) Dibenzo(a,h)anthracene	25.60	278	714137	51.85	ng	# 97
91) Benzo(g,h,i)perylene	26.28	276	667885	50.99	ng	98

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

