

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG041615\
 Data File : BG016462.D
 Acq On : 16 Apr 2015 19:06
 Operator : TP/IZ
 Sample : G1860-03MSD
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-1MSD

Quant Time: Apr 17 04:40:57 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG040615.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 17 03:57:14 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	61030	20.00	ng	0.00
21) Naphthalene-d8	10.45	136	265469	20.00	ng	0.00
38) Acenaphthene-d10	14.31	164	157559	20.00	ng	0.00
63) Phenanthrene-d10	17.05	188	320800	20.00	ng	0.00
75) Chrysene-d12	21.23	240	263177	20.00	ng	0.00
86) Perylene-d12	23.45	264	251913	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.26	112	256147	66.24	ng	0.00
7) Phenol-d6	6.85	99	223695	38.53	ng	0.00
23) Nitrobenzene-d5	8.82	82	407241	88.87	ng	0.00
41) 2,4,6-Tribromophenol	15.80	330	164758	154.62	ng	0.00
44) 2-Fluorobiphenyl	12.93	172	883028	95.42	ng	0.00
78) Terphenyl-d14	19.68	244	931725	82.86	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.14	88	30665	17.44	ng	95
3) Pyridine	3.54	79	72264	13.75	ng	93
4) n-Nitrosodimethylamine	3.45	42	24842	15.91	ng	99
6) Aniline	6.99	93	131352	15.84	ng	98
8) 2-Chlorophenol	7.23	128	185133	39.87	ng	94
9) Benzaldehyde	6.80	77	48237	14.19	ng	95
10) Phenol	6.88	94	88863	14.31	ng	97
11) bis(2-Chloroethyl)ether	7.09	93	203745	41.14	ng	99
12) 1,3-Dichlorobenzene	7.55	146	199085	42.45	ng	95
13) 1,4-Dichlorobenzene	7.70	146	202298	41.58	ng	97
14) 1,2-Dichlorobenzene	8.01	146	195682	42.06	ng	97
15) Benzyl Alcohol	7.91	79	109135	26.97	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.20	45	210295	37.70	ng	96
17) 2-Methylphenol	8.12	107	128520	30.86	ng	97
18) Hexachloroethane	8.73	117	73780	39.64	ng	96
19) n-Nitroso-di-n-propylamine	8.47	70	156300	37.58	ng	97
20) 3+4-Methylphenols	8.45	107	158679	27.50	ng	97
22) Acetophenone	8.48	105	331392	53.84	ng	98
24) Nitrobenzene	8.86	77	216133	44.10	ng	92
25) Isophorone	9.38	82	438578	43.10	ng	97
26) 2-Nitrophenol	9.57	139	126785	55.48	ng	94
27) 2,4-Dimethylphenol	9.64	122	193679	45.50	ng	98
28) bis(2-Chloroethoxy)methane	9.87	93	265862	45.77	ng	98
29) 2,4-Dichlorophenol	10.11	162	182569	54.21	ng	97
30) 1,2,4-Trichlorobenzene	10.31	180	172482	49.05	ng	99
31) Naphthalene	10.50	128	638083	46.13	ng	99
32) Benzoic acid	9.75	122	42897	21.37	ng	89
33) 4-Chloroaniline	10.61	127	107624	16.58	ng	95
34) Hexachlorobutadiene	10.78	225	72722	47.08	ng	96
35) Caprolactam	11.43	113	16956m	8.00	ng	
36) 4-Chloro-3-methylphenol	11.76	107	202710	45.56	ng	93
37) 2-Methylnaphthalene	12.11	142	480080	48.25	ng	99
39) 1,2,4,5-Tetrachlorobenzene	12.48	216	169737	48.93	ng	99
40) Hexachlorocyclopentadiene	12.47	237	144702	91.13	ng	97

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42) 2,4,6-Trichlorophenol	12.74	196	141850	55.03	ng	98
43) 2,4,5-Trichlorophenol	12.81	196	154954	56.26	ng	96
45) 1,1'-Biphenyl	13.14	154	590923	48.89	ng	98
46) 2-Chloronaphthalene	13.18	162	446227	48.47	ng	100
47) 2-Nitroaniline	13.39	65	140292	47.27	ng	90
48) Acenaphthylene	14.03	152	775261	47.35	ng	99
49) Dimethylphthalate	13.77	163	579976	50.22	ng	99
50) 2,6-Dinitrotoluene	13.89	165	146246	52.36	ng	93
51) Acenaphthene	14.37	154	474103	48.46	ng	100
52) 3-Nitroaniline	14.22	138	71128	19.94	ng	87
53) 2,4-Dinitrophenol	14.44	184	167458	101.17	ng	91
54) Dibenzofuran	14.71	168	685572	50.50	ng	98
55) 4-Nitrophenol	14.55	139	103709	35.18	ng	91
56) 2,4-Dinitrotoluene	14.68	165	201285	52.91	ng	91
57) Fluorene	15.35	166	591643	49.40	ng	99
58) 2,3,4,6-Tetrachlorophenol	14.94	232	120157	56.53	ng	93
59) Diethylphthalate	15.13	149	609442	49.58	ng	100
60) 4-Chlorophenyl-phenylether	15.35	204	247166	50.45	ng	97
61) 4-Nitroaniline	15.39	138	171617	42.39	ng	97
62) Azobenzene	15.64	77	589206	45.04	ng	98
64) 4,6-Dinitro-2-methylphenol	15.45	198	112992	49.88	ng	90
65) n-Nitrosodiphenylamine	15.57	169	535712	49.40	ng	99
66) 4-Bromophenyl-phenylether	16.25	248	132496	51.73	ng	94
67) Hexachlorobenzene	16.36	284	133132	50.01	ng	99
68) Atrazine	16.52	200	166317	55.03	ng	100
69) Pentachlorophenol	16.71	266	164328	101.09	ng	97
70) Phenanthrene	17.09	178	877793	48.65	ng	99
71) Anthracene	17.18	178	880716	49.25	ng	99
72) Carbazole	17.46	167	892821	49.75	ng	99
73) Di-n-butylphthalate	18.01	149	1054175	48.00	ng	99
74) Fluoranthene	19.11	202	892434	47.99	ng	97
76) Benzidine	19.29	184	259209	23.23	ng	97
77) Pyrene	19.47	202	917461	50.06	ng	99
79) Butylbenzylphthalate	20.37	149	479508	53.26	ng	97
80) Benzo(a)anthracene	21.22	228	778543	50.05	ng	100
81) 3,3'-Dichlorobenzidine	21.15	252	158852	31.06	ng	97
82) Chrysene	21.26	228	730125	48.62	ng	98
83) Bis(2-ethylhexyl)phthalate	21.14	149	655164	53.75	ng	97
84) Di-n-octyl phthalate	22.01	149	1097820	55.28	ng	97
85) Indeno(1,2,3-cd)pyrene	25.73	276	813122	52.12	ng	98
87) Benzo(b)fluoranthene	22.78	252	746663	50.61	ng	98
88) Benzo(k)fluoranthene	22.83	252	704969	48.83	ng	98
89) Benzo(a)pyrene	23.36	252	713671	51.61	ng	100
90) Dibenzo(a,h)anthracene	25.74	278	684240	50.85	ng	99
91) Benzo(g,h,i)perylene	26.42	276	699519	52.05	ng	98

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

