

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG062616\
 Data File : BG022599.D
 Acq On : 26 Jun 2016 7:25
 Operator : UM/SJ
 Sample : H3663-13MS
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP-7MS

Quant Time: Jun 26 08:51:34 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG062516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 24 20:14:59 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.16	152	130854	20.00	ng	-0.01
21) Naphthalene-d8	10.99	136	549251	20.00	ng	0.00
38) Acenaphthene-d10	14.80	164	398300	20.00	ng	0.00
63) Phenanthrene-d10	17.56	188	1041824	20.00	ng	0.00
75) Chrysene-d12	21.85	240	1159801	20.00	ng	0.00
86) Perylene-d12	25.21	264	1183510	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.71	112	1229248	150.68	ng	0.00
7) Phenol-d6	7.31	99	1847233	160.64	ng	0.01
23) Nitrobenzene-d5	9.34	82	1293831	99.70	ng	0.00
41) 2,4,6-Tribromophenol	16.29	330	1025442	163.69	ng	0.00
44) 2-Fluorobiphenyl	13.43	172	2366571	94.23	ng	0.00
78) Terphenyl-d14	20.16	244	3362097	76.80	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.58	88	123739	37.42	ng	95
3) Pyridine	3.99	79	395863	42.80	ng	91
4) n-Nitrosodimethylamine	3.90	42	230356	43.54	ng	# 92
6) Aniline	7.48	93	691828	45.38	ng	98
8) 2-Chlorophenol	7.73	128	441184	52.21	ng	99
9) Benzaldehyde	7.30	77	17272	4.27	ng	# 77
10) Phenol	7.34	94	657325	54.08	ng	95
11) bis(2-Chloroethyl)ether	7.58	93	469970	46.97	ng	90
12) 1,3-Dichlorobenzene	8.05	146	463793	45.10	ng	99
13) 1,4-Dichlorobenzene	8.20	146	473304	45.91	ng	99
14) 1,2-Dichlorobenzene	8.52	146	456989	46.61	ng	95
15) Benzyl Alcohol	8.40	79	589418	55.40	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.70	45	535673	48.41	ng	97
17) 2-Methylphenol	8.60	107	425957	53.17	ng	97
18) Hexachloroethane	9.26	117	195484	45.81	ng	95
19) n-Nitroso-di-n-propylamine	8.97	70	471004	49.18	ng	99
20) 3+4-Methylphenols	8.93	107	602246	54.57	ng	97
22) Acetophenone	8.99	105	790683	49.80	ng	95
24) Nitrobenzene	9.38	77	689341	48.06	ng	95
25) Isophorone	9.90	82	1235353	49.04	ng	99
26) 2-Nitrophenol	10.09	139	274811	53.19	ng	88
27) 2,4-Dimethylphenol	10.14	122	480699	48.70	ng	92
28) bis(2-Chloroethoxy)methane	10.38	93	686465	49.91	ng	99
29) 2,4-Dichlorophenol	10.63	162	530690	55.18	ng	99
30) 1,2,4-Trichlorobenzene	10.85	180	550492	48.21	ng	98
31) Naphthalene	11.04	128	1329619	48.16	ng	99
32) Benzoic acid	10.24	122	158997	27.13	ng	90
33) 4-Chloroaniline	11.15	127	417305	35.09	ng	94
34) Hexachlorobutadiene	11.32	225	410125	46.21	ng	97
35) Caprolactam	11.93	113	187104m	61.76	ng	
36) 4-Chloro-3-methylphenol	12.25	107	650534	55.10	ng	98
37) 2-Methylnaphthalene	12.63	142	1046372	48.87	ng	97
39) 1,2,4,5-Tetrachlorobenzene	13.00	216	704720	46.87	ng	99
40) Hexachlorocyclopentadiene	12.98	237	934125	77.82	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.23	196	504910	52.36	ng	99
43) 2,4,5-Trichlorophenol	13.31	196	540644	53.58	ng	96
45) 1,1'-Biphenyl	13.63	154	1432981	47.23	ng	96
46) 2-Chloronaphthalene	13.68	162	1188904	47.62	ng	97
47) 2-Nitroaniline	13.88	65	515632	53.51	ng	95
48) Acenaphthylene	14.52	152	1728634	47.74	ng	96
49) Dimethylphthalate	14.25	163	1931633	58.46	ng	# 96
50) 2,6-Dinitrotoluene	14.37	165	383634	53.43	ng	92
51) Acenaphthene	14.86	154	1157149	43.38	ng	98
52) 3-Nitroaniline	14.70	138	305649	45.50	ng	94
53) 2,4-Dinitrophenol	14.90	184	393095	88.89	ng	95
54) Dibenzofuran	15.20	168	1835397	47.50	ng	100
55) 4-Nitrophenol	15.01	139	632916	111.41	ng	98
56) 2,4-Dinitrotoluene	15.15	165	567690	57.08	ng	96
57) Fluorene	15.85	166	1510718	49.00	ng	98
58) 2,3,4,6-Tetrachlorophenol	15.42	232	572195	54.70	ng	99
59) Diethylphthalate	15.61	149	1672570	49.24	ng	98
60) 4-Chlorophenyl-phenylether	15.83	204	905893	49.35	ng	99
61) 4-Nitroaniline	15.87	138	383097	55.27	ng	96
62) Azobenzene	16.13	77	1691713	45.99	ng	95
64) 4,6-Dinitro-2-methylphenol	15.92	198	367137	52.55	ng	95
65) n-Nitrosodiphenylamine	16.05	169	1444917	47.66	ng	96
66) 4-Bromophenyl-phenylether	16.73	248	630515	46.65	ng	98
67) Hexachlorobenzene	16.85	284	671072	46.96	ng	100
68) Atrazine	17.00	200	641493	50.12	ng	97
69) Pentachlorophenol	17.20	266	969418	98.57	ng	100
70) Phenanthrene	17.60	178	2405970	45.29	ng	94
71) Anthracene	17.69	178	2433204	44.50	ng	94
72) Carbazole	17.96	167	2240536	48.34	ng	97
73) Di-n-butylphthalate	18.51	149	2511189	46.52	ng	96
74) Fluoranthene	19.60	202	2830554	46.24	ng	96
76) Benzidine	19.78	184	1799597	145.74	ng	98
77) Pyrene	19.96	202	2807855	45.45	ng	93
79) Butylbenzylphthalate	20.84	149	1307792	48.91	ng	97
80) Benzo(a)anthracene	21.83	228	3026407	47.16	ng	95
81) 3,3'-Dichlorobenzidine	21.74	252	888976	33.54	ng	97
82) Chrysene	21.90	228	2853491	47.32	ng	91
83) Bis(2-ethylhexyl)phthalate	21.72	149	1766308	46.91	ng	96
84) Di-n-octyl phthalate	22.98	149	2896342	46.67	ng	100
85) Indeno(1,2,3-cd)pyrene	29.07	276	3849205	48.57	ng	# 100
87) Benzo(b)fluoranthene	24.14	252	3391692	47.66	ng	93
88) Benzo(k)fluoranthene	24.21	252	3148689	46.80	ng	# 95
89) Benzo(a)pyrene	25.05	252	3307018	47.67	ng	# 94
90) Dibenzo(a,h)anthracene	29.14	278	3197358	48.96	ng	# 93
91) Benzo(g,h,i)perylene	30.28	276	3222329	48.47	ng	# 97

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

