

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG061815\
 Data File : BG017718.D
 Acq On : 18 Jun 2015 11:29
 Operator : TP/IZ
 Sample : SSTDICC010
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC010

Quant Time: Jun 19 04:27:04 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG061815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 19 04:20:56 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	65398	20.00	ng	0.00
21) Naphthalene-d8	10.40	136	307559	20.00	ng	0.00
38) Acenaphthene-d10	14.27	164	192442	20.00	ng	0.00
63) Phenanthrene-d10	17.03	188	432127	20.00	ng	0.00
75) Chrysene-d12	21.23	240	451921	20.00	ng	0.00
86) Perylene-d12	23.47	264	426311	20.00	ng	0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.22	112	75073	16.98	ng	0.00
7) Phenol-d6	6.79	99	105555	19.77	ng	0.00
23) Nitrobenzene-d5	8.77	82	106186	21.06	ng	0.00
41) 2,4,6-Tribromophenol	15.77	330	34379	22.51	ng	0.00
44) 2-Fluorobiphenyl	12.89	172	250937	16.12	ng	0.00
78) Terphenyl-d14	19.68	244	443977	20.03	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.18	88	16922	9.75	ng	# 92
3) Pyridine	3.57	79	44448	9.82	ng	94
4) n-Nitrosodimethylamine	3.49	42	19154	9.42	ng	# 87
6) Aniline	6.95	93	71131	10.62	ng	97
8) 2-Chlorophenol	7.18	128	47226	9.69	ng	94
9) Benzaldehyde	6.77	77	36874	19.21	ng	93
10) Phenol	6.82	94	57334	10.16	ng	97
11) bis(2-Chloroethyl)ether	7.05	93	42870	9.76	ng	97
12) 1,3-Dichlorobenzene	7.51	146	52830	10.07	ng	98
13) 1,4-Dichlorobenzene	7.65	146	55320	10.26	ng	93
14) 1,2-Dichlorobenzene	7.96	146	52335	10.39	ng	98
15) Benzyl Alcohol	7.86	79	44934	14.30	ng	95
16) 2,2'-oxybis(1-Chloropropan	8.16	45	68195	8.42	ng	99
17) 2-Methylphenol	8.06	107	40665	11.40	ng	97
18) Hexachloroethane	8.69	117	19991	11.38	ng	97
19) n-Nitroso-di-n-propylamine	8.43	70	44871	15.13	ng	98
20) 3+4-Methylphenols	8.39	107	55942	11.75	ng	96
22) Acetophenone	8.44	105	69254	9.27	ng	# 94
24) Nitrobenzene	8.81	77	57516	11.62	ng	97
25) Isophorone	9.34	82	119520	12.13	ng	99
26) 2-Nitrophenol	9.52	139	23502	7.80	ng	99
27) 2,4-Dimethylphenol	9.59	122	48866	9.09	ng	99
28) bis(2-Chloroethoxy)methane	9.83	93	58852	9.36	ng	94
29) 2,4-Dichlorophenol	10.05	162	43268	9.95	ng	97
30) 1,2,4-Trichlorobenzene	10.27	180	49401	10.52	ng	97
31) Naphthalene	10.45	128	169086	9.58	ng	98
32) Benzoic acid	9.66	122	16863	4.96	ng	95
33) 4-Chloroaniline	10.57	127	72104	10.37	ng	98
34) Hexachlorobutadiene	10.74	225	27772	13.15	ng	93
35) Caprolactam	11.32	113	24283	14.30	ng	97
36) 4-Chloro-3-methylphenol	11.69	107	53531	12.94	ng	96
37) 2-Methylnaphthalene	12.08	142	120675	11.02	ng	97
39) 1,2,4,5-Tetrachlorobenzene	12.45	216	51564	9.60	ng	98
40) Hexachlorocyclopentadiene	12.43	237	33595	10.39	ng	99

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42) 2,4,6-Trichlorophenol	12.69	196	33540	9.35	ng	97
43) 2,4,5-Trichlorophenol	12.76	196	37233	9.95	ng	97
45) 1,1'-Biphenyl	13.10	154	146089	8.13	ng	98
46) 2-Chloronaphthalene	13.14	162	115376	9.24	ng	99
47) 2-Nitroaniline	13.35	65	31879	9.49	ng	99
48) Acenaphthylene	13.99	152	211488	9.56	ng	98
49) Dimethylphthalate	13.75	163	151651	11.40	ng	97
50) 2,6-Dinitrotoluene	13.86	165	28012	9.14	ng	99
51) Acenaphthene	14.34	154	125769	9.27	ng	94
52) 3-Nitroaniline	14.18	138	32985	8.74	ng	99
53) 2,4-Dinitrophenol	14.39	184	8957	5.62	ng	97
54) Dibenzofuran	14.68	168	180307	10.28	ng	95
55) 4-Nitrophenol	14.49	139	25608	8.47	ng	94
56) 2,4-Dinitrotoluene	14.64	165	33002	8.64	ng	# 98
57) Fluorene	15.33	166	162817	10.58	ng	96
58) 2,3,4,6-Tetrachlorophenol	14.90	232	30406	10.54	ng	98
59) Diethylphthalate	15.12	149	167514	12.41	ng	97
60) 4-Chlorophenyl-phenylether	15.33	204	71846	11.47	ng	97
61) 4-Nitroaniline	15.35	138	37141	9.66	ng	93
62) Azobenzene	15.63	77	175054	13.79	ng	97
64) 4,6-Dinitro-2-methylphenol	15.40	198	15317	6.02	ng	90
65) n-Nitrosodiphenylamine	15.55	169	138813	8.85	ng	97
66) 4-Bromophenyl-phenylether	16.23	248	42579	10.36	ng	99
67) Hexachlorobenzene	16.33	284	46488	10.52	ng	98
68) Atrazine	16.51	200	52711	13.43	ng	99
69) Pentachlorophenol	16.68	266	21604	7.93	ng	96
70) Phenanthrene	17.07	178	254011	9.62	ng	98
71) Anthracene	17.16	178	250947	9.56	ng	98
72) Carbazole	17.43	167	250879	10.27	ng	96
73) Di-n-butylphthalate	18.02	149	326423	11.12	ng	98
74) Fluoranthene	19.10	202	298787	12.60	ng	97
76) Benzidine	19.29	184	163189	14.05	ng	99
77) Pyrene	19.46	202	307851	8.85	ng	97
79) Butylbenzylphthalate	20.39	149	137734	8.60	ng	100
80) Benzo(a)anthracene	21.22	228	275487	10.18	ng	94
81) 3,3'-Dichlorobenzidine	21.16	252	96946	11.28	ng	98
82) Chrysene	21.27	228	250835	9.63	ng	95
83) Bis(2-ethylhexyl)phthalate	21.18	149	194672	8.60	ng	96
84) Di-n-octyl phthalate	22.06	149	333284	9.72	ng	99
85) Indeno(1,2,3-cd)pyrene	25.74	276	284009	12.21	ng	100
87) Benzo(b)fluoranthene	22.80	252	253842	9.54	ng	96
88) Benzo(k)fluoranthene	22.85	252	258472	9.59	ng	97
89) Benzo(a)pyrene	23.37	252	237591	9.56	ng	98
90) Dibenzo(a,h)anthracene	25.77	278	236806	10.72	ng	98
91) Benzo(g,h,i)perylene	26.44	276	230688	10.34	ng	98

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

