

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG090116\
 Data File : BG023939.D
 Acq On : 1 Sep 2016 19:11
 Operator : UM/SJ
 Sample : PB93192BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB93192BS

Quant Time: Sep 02 01:19:51 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG083116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 31 20:05:47 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.09	152	24263	20.00	ng	0.00
21) Naphthalene-d8	10.92	136	126760	20.00	ng	-0.01
38) Acenaphthene-d10	14.76	164	107601	20.00	ng	0.00
63) Phenanthrene-d10	17.52	188	294734	20.00	ng	0.00
75) Chrysene-d12	21.84	240	316296	20.00	ng	0.00
86) Perylene-d12	25.21	264	314098	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.64	112	209392	151.52	ng	0.00
7) Phenol-d6	7.26	99	315770	148.44	ng	0.00
23) Nitrobenzene-d5	9.27	82	234000	82.41	ng	0.00
41) 2,4,6-Tribromophenol	16.26	330	239419	123.50	ng	0.00
44) 2-Fluorobiphenyl	13.37	172	582020	77.55	ng	0.00
78) Terphenyl-d14	20.13	244	1256131	90.44	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.49	88	18285	32.08	ng	# 91
3) Pyridine	3.90	79	62021	36.21	ng	94
4) n-Nitrosodimethylamine	3.81	42	48320	53.52	ng	# 60
6) Aniline	7.41	93	89032	31.53	ng	96
8) 2-Chlorophenol	7.65	128	77767	48.57	ng	96
9) Benzaldehyde	7.23	77	12179	8.06	ng	88
10) Phenol	7.29	94	110056	49.18	ng	96
11) bis(2-Chloroethyl)ether	7.50	93	67975	38.60	ng	91
12) 1,3-Dichlorobenzene	7.97	146	80698	43.07	ng	96
13) 1,4-Dichlorobenzene	8.12	146	80236	40.42	ng	99
14) 1,2-Dichlorobenzene	8.44	146	75517	40.53	ng	# 91
15) Benzyl Alcohol	8.34	79	97870	52.18	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.62	45	94883	40.20	ng	91
17) 2-Methylphenol	8.55	107	74881	49.66	ng	94
18) Hexachloroethane	9.17	117	32856	42.98	ng	92
19) n-Nitroso-di-n-propylamine	8.90	70	81363	43.29	ng	# 89
20) 3+4-Methylphenols	8.88	107	105112	50.02	ng	95
22) Acetophenone	8.92	105	117266	35.73	ng	# 95
24) Nitrobenzene	9.31	77	123835	40.56	ng	99
25) Isophorone	9.84	82	212993	38.68	ng	99
26) 2-Nitrophenol	10.03	139	47434	40.87	ng	97
27) 2,4-Dimethylphenol	10.09	122	94733	48.02	ng	87
28) bis(2-Chloroethoxy)methane	10.31	93	113310	40.76	ng	97
29) 2,4-Dichlorophenol	10.58	162	99803	47.75	ng	99
30) 1,2,4-Trichlorobenzene	10.78	180	93731	40.89	ng	94
31) Naphthalene	10.97	128	268321	41.08	ng	97
32) Benzoic acid	10.26	122	50202	31.34	ng	# 80
33) 4-Chloroaniline	11.10	127	63892	23.43	ng	# 92
34) Hexachlorobutadiene	11.24	225	71656	41.62	ng	96
35) Caprolactam	11.90	113	33001m	44.80	ng	
36) 4-Chloro-3-methylphenol	12.23	107	126575	45.32	ng	90
37) 2-Methylnaphthalene	12.58	142	224726	45.23	ng	98
39) 1,2,4,5-Tetrachlorobenzene	12.94	216	116982	32.75	ng	95
40) Hexachlorocyclopentadiene	12.91	237	160949	75.25	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.20	196	95044	39.88	ng	97
43) 2,4,5-Trichlorophenol	13.28	196	100022	41.51	ng	87
45) 1,1'-Biphenyl	13.58	154	293635	33.93	ng	97
46) 2-Chloronaphthalene	13.63	162	244437	38.47	ng	94
47) 2-Nitroaniline	13.85	65	114157	45.61	ng	98
48) Acenaphthylene	14.48	152	421267	39.78	ng	97
49) Dimethylphthalate	14.21	163	373953	40.53	ng	98
50) 2,6-Dinitrotoluene	14.34	165	80167	41.16	ng	98
51) Acenaphthene	14.82	154	263674	40.27	ng	94
52) 3-Nitroaniline	14.68	138	45965	25.23	ng	# 82
53) 2,4-Dinitrophenol	14.89	184	108437	76.46	ng	# 81
54) Dibenzofuran	15.16	168	449043	43.94	ng	99
55) 4-Nitrophenol	15.01	139	117048	80.75	ng	84
56) 2,4-Dinitrotoluene	15.13	165	125906	43.98	ng	# 81
57) Fluorene	15.81	166	378978	42.76	ng	95
58) 2,3,4,6-Tetrachlorophenol	15.39	232	116764m	47.75	ng	
59) Diethylphthalate	15.57	149	419160	42.55	ng	97
60) 4-Chlorophenyl-phenylether	15.79	204	209760	43.97	ng	96
61) 4-Nitroaniline	15.86	138	79178	42.99	ng	92
62) Azobenzene	16.09	77	462032	48.43	ng	97
64) 4,6-Dinitro-2-methylphenol	15.91	198	86618	42.33	ng	96
65) n-Nitrosodiphenylamine	16.02	169	357975	42.76	ng	97
66) 4-Bromophenyl-phenylether	16.70	248	154956	43.21	ng	98
67) Hexachlorobenzene	16.82	284	162263	38.58	ng	96
68) Atrazine	16.97	200	154977	43.03	ng	98
69) Pentachlorophenol	17.18	266	183389	82.10	ng	97
70) Phenanthrene	17.57	178	692841	45.19	ng	95
71) Anthracene	17.66	178	708959	45.33	ng	95
72) Carbazole	17.94	167	614716	43.31	ng	96
73) Di-n-butylphthalate	18.47	149	754727	44.57	ng	97
74) Fluoranthene	19.58	202	868502	47.05	ng	94
76) Benzidine	19.76	184	346400	35.37	ng	96
77) Pyrene	19.94	202	885306	45.51	ng	92
79) Butylbenzylphthalate	20.82	149	331985	44.27	ng	97
80) Benzo(a)anthracene	21.81	228	835556	47.19	ng	92
81) 3,3'-Dichlorobenzidine	21.73	252	169533	23.95	ng	# 97
82) Chrysene	21.88	228	778533	46.19	ng	95
83) Bis(2-ethylhexyl)phthalate	21.68	149	449821	43.81	ng	95
84) Di-n-octyl phthalate	22.94	149	750768	44.58	ng	99
85) Indeno(1,2,3-cd)pyrene	29.07	276	876775	46.99	ng	# 100
87) Benzo(b)fluoranthene	24.13	252	828132	46.41	ng	98
88) Benzo(k)fluoranthene	24.20	252	799983	45.22	ng	97
89) Benzo(a)pyrene	25.05	252	781866	46.15	ng	98
90) Dibenzo(a,h)anthracene	29.14	278	723649	43.43	ng	94
91) Benzo(g,h,i)perylene	30.29	276	726433	45.33	ng	98

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

