

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052517\
 Data File : BM010215.D
 Acq On : 25 May 2017 13:56
 Operator : SJ/MA
 Sample : PB99246BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB99246BS

Quant Time: May 26 05:12:32 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM051917.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat May 20 04:25:49 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.95	152	141257	20.00	ng	-0.03
21) Naphthalene-d8	10.75	136	480884	20.00	ng	-0.04
38) Acenaphthene-d10	14.58	164	293672	20.00	ng	-0.03
63) Phenanthrene-d10	17.32	188	620290	20.00	ng	-0.03
75) Chrysene-d12	21.49	240	770905	20.00	ng	-0.02
86) Perylene-d12	23.84	264	941374	20.00	ng	-0.05

System Monitoring Compounds

5) 2-Fluorophenol	5.53	112	1185277	151.53	ng	-0.03
7) Phenol-d6	7.14	99	1406064	139.69	ng	-0.03
23) Nitrobenzene-d5	9.13	82	969583	108.78	ng	-0.04
41) 2,4,6-Tribromophenol	16.07	330	619841	138.97	ng	-0.02
44) 2-Fluorobiphenyl	13.20	172	2464260	107.87	ng	-0.04
78) Terphenyl-d14	19.93	244	3079639	90.83	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.43	88	160053	44.53	ng	# 100
3) Pyridine	3.84	79	418931	43.54	ng	90
4) n-Nitrosodimethylamine	3.76	42	219579	63.10	ng	# 75
6) Aniline	7.30	93	262366	21.04	ng	93
8) 2-Chlorophenol	7.52	128	377006	44.30	ng	92
9) Benzaldehyde	7.11	77	96985	15.48	ng	90
10) Phenol	7.16	94	473262	44.62	ng	95
11) bis(2-Chloroethyl)ether	7.38	93	340620	43.08	ng	97
12) 1,3-Dichlorobenzene	7.83	146	461399	42.66	ng	# 89
13) 1,4-Dichlorobenzene	7.98	146	468132	42.14	ng	96
14) 1,2-Dichlorobenzene	8.30	146	453969	42.57	ng	97
15) Benzyl Alcohol	8.21	79	342009	44.91	ng	# 86
16) 2,2'-oxybis(1-Chloropropan	8.47	45	275575	35.45	ng	75
17) 2-Methylphenol	8.40	107	308450	40.87	ng	# 83
18) Hexachloroethane	9.02	117	162169	45.03	ng	# 75
19) n-Nitroso-di-n-propylamine	8.75	70	318342	44.81	ng	93
20) 3+4-Methylphenols	8.73	107	412097	40.43	ng	87
22) Acetophenone	8.79	105	580580	47.10	ng	# 98
24) Nitrobenzene	9.18	77	486816	53.02	ng	92
25) Isophorone	9.68	82	743125	48.25	ng	# 93
26) 2-Nitrophenol	9.88	139	190393	49.19	ng	# 61
27) 2,4-Dimethylphenol	9.93	122	319861	45.77	ng	# 79
28) bis(2-Chloroethoxy)methane	10.16	93	426112	43.59	ng	100
29) 2,4-Dichlorophenol	10.40	162	401394	49.88	ng	99
30) 1,2,4-Trichlorobenzene	10.60	180	497200	47.76	ng	99
31) Naphthalene	10.80	128	1082110	42.07	ng	100
32) Benzoic acid	10.08	122	199284	41.84	ng	# 73
33) 4-Chloroaniline	10.96	127	144124	14.54	ng	# 83
34) Hexachlorobutadiene	11.05	225	348184	50.14	ng	97
35) Caprolactam	11.76	113	83216	36.04	ng	# 82
36) 4-Chloro-3-methylphenol	12.06	107	362866	41.64	ng	# 83
37) 2-Methylnaphthalene	12.40	142	843811	45.56	ng	# 94
39) 1,2,4,5-Tetrachlorobenzene	12.76	216	566568	50.17	ng	# 100
40) Hexachlorocyclopentadiene	12.72	237	867373	133.02	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)

42) 2,4,6-Trichlorophenol	13.02	196	340958	51.26	ng	99
43) 2,4,5-Trichlorophenol	13.09	196	357073	48.58	ng #	93
45) 1,1'-Biphenyl	13.42	154	1138459	47.14	ng	99
46) 2-Chloronaphthalene	13.46	162	901216	46.09	ng	95
47) 2-Nitroaniline	13.70	65	237380	47.21	ng #	80
48) Acenaphthylene	14.31	152	1337104	46.25	ng	100
49) Dimethylphthalate	14.05	163	1070467	40.86	ng #	98
50) 2,6-Dinitrotoluene	14.17	165	222691	44.77	ng #	82
51) Acenaphthene	14.65	154	813602	45.27	ng	98
52) 3-Nitroaniline	14.53	138	176351	38.14	ng #	72
53) 2,4-Dinitrophenol	14.72	184	281564	84.78	ng #	85
54) Dibenzofuran	14.98	168	1363156	48.23	ng	96
55) 4-Nitrophenol	14.84	139	277142	74.49	ng #	23
56) 2,4-Dinitrotoluene	14.96	165	297335	42.25	ng #	87
57) Fluorene	15.63	166	1090363	44.36	ng	99
58) 2,3,4,6-Tetrachlorophenol	15.20	232	324369	51.81	ng #	100
59) Diethylphthalate	15.39	149	988169	39.90	ng	98
60) 4-Chlorophenyl-phenylether	15.62	204	615173	44.58	ng	97
61) 4-Nitroaniline	15.69	138	165670	35.30	ng #	17
62) Azobenzene	15.91	77	1017746	55.89	ng	88
64) 4,6-Dinitro-2-methylphenol	15.71	198	179471	47.36	ng	80
65) n-Nitrosodiphenylamine	15.84	169	899390	48.38	ng	99
66) 4-Bromophenyl-phenylether	16.51	248	391812	49.34	ng	94
67) Hexachlorobenzene	16.60	284	433454	44.93	ng	94
68) Atrazine	16.78	200	335783	47.64	ng	98
69) Pentachlorophenol	16.97	266	508565	95.98	ng	98
70) Phenanthrene	17.37	178	1596755	46.01	ng	99
71) Anthracene	17.46	178	1621740	47.09	ng	100
72) Carbazole	17.75	167	1323780	43.41	ng	99
73) Di-n-butylphthalate	18.26	149	1446986	40.02	ng #	97
74) Fluoranthene	19.37	202	1840647	40.85	ng	97
76) Benzidine	19.58	184	1134461	79.78	ng	99
77) Pyrene	19.74	202	1879062	45.46	ng	99
79) Butylbenzylphthalate	20.61	149	617689	40.37	ng #	84
80) Benzo(a)anthracene	21.47	228	1976680	46.53	ng	99
81) 3,3'-Dichlorobenzidine	21.41	252	600553	35.14	ng #	96
82) Chrysene	21.52	228	1772933	44.00	ng	100
83) Bis(2-ethylhexyl)phthalate	21.35	149	888732	38.84	ng #	98
84) Di-n-octyl phthalate	22.26	149	1584596	38.68	ng #	98
85) Indeno(1,2,3-cd)pyrene	26.26	276	3687251	62.40	ng #	100
87) Benzo(b)fluoranthene	23.12	252	2396557	43.50	ng #	92
88) Benzo(k)fluoranthene	23.17	252	2350269	43.54	ng #	96
89) Benzo(a)pyrene	23.74	252	2467487	47.00	ng #	97
90) Dibenzo(a,h)anthracene	26.27	278	3076616	52.41	ng #	92
91) Benzo(g,h,i)perylene	27.01	276	3029426	52.66	ng #	85

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

