

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062215\
 Data File : BF080097.D
 Acq On : 22 Jun 2015 16:39
 Operator : TP/IZ
 Sample : PB84131BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB84131BS

Quant Time: Jun 23 01:19:12 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 23 00:53:56 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.33	152	50531	20.00	ng	0.00
21) Naphthalene-d8	8.61	136	200546	20.00	ng	-0.01
38) Acenaphthene-d10	10.39	164	104842	20.00	ng	0.00
63) Phenanthrene-d10	11.89	188	219817	20.00	ng	-0.02
75) Chrysene-d12	14.92	240	249678	20.00	ng	-0.08
86) Perylene-d12	16.83	264	237405	20.00	ng	-0.10

System Monitoring Compounds

5) 2-Fluorophenol	5.93	112	231147	80.97	ng	0.00
7) Phenol-d6	6.92	99	286966	75.54	ng	-0.01
23) Nitrobenzene-d5	7.88	82	313988	91.15	ng	0.00
41) 2,4,6-Tribromophenol	11.18	330	82875	80.84	ng	0.00
44) 2-Fluorobiphenyl	9.69	172	623205	84.77	ng	0.00
78) Terphenyl-d14	13.61	244	812314	75.98	ng	-0.06

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.21	88	36198	26.68	ng	86
3) Pyridine	4.02	79	106122	29.41	ng	# 80
4) n-Nitrosodimethylamine	3.93	42	46602	27.18	ng	98
6) Aniline	6.97	93	157486	30.13	ng	92
8) 2-Chlorophenol	7.10	128	108770	32.97	ng	82
9) Benzaldehyde	6.87	77	53983	24.62	ng	# 88
10) Phenol	6.94	94	140973	34.01	ng	85
11) bis(2-Chloroethyl)ether	7.04	93	103046	31.33	ng	87
12) 1,3-Dichlorobenzene	7.27	146	119013	30.03	ng	96
13) 1,4-Dichlorobenzene	7.34	146	135451m	35.94	ng	
14) 1,2-Dichlorobenzene	7.50	146	124914	34.06	ng	97
15) Benzyl Alcohol	7.44	79	96288	31.00	ng	# 73
16) 2,2'-oxybis(1-Chloropropan	7.59	45	181132	37.10	ng	85
17) 2-Methylphenol	7.56	107	89311	31.69	ng	# 90
18) Hexachloroethane	7.84	117	41580	33.13	ng	90
19) n-Nitroso-di-n-propylamine	7.72	70	88416	33.53	ng	# 77
20) 3+4-Methylphenols	7.70	107	124549	34.24	ng	93
22) Acetophenone	7.72	105	157511	33.70	ng	# 79
24) Nitrobenzene	7.90	77	126179	35.49	ng	# 72
25) Isophorone	8.13	82	207264	29.90	ng	# 82
26) 2-Nitrophenol	8.22	139	58819	39.52	ng	# 70
27) 2,4-Dimethylphenol	8.24	122	112456	36.24	ng	# 85
28) bis(2-Chloroethoxy)methane	8.33	93	128768	31.61	ng	# 91
29) 2,4-Dichlorophenol	8.46	162	105088	39.54	ng	98
30) 1,2,4-Trichlorobenzene	8.55	180	107744	35.70	ng	99
31) Naphthalene	8.63	128	315308m	30.07	ng	
32) Benzoic acid	8.30	122	41557	22.15	ng	# 61
33) 4-Chloroaniline	8.68	127	107038m	23.85	ng	
34) Hexachlorobutadiene	8.76	225	63948	34.41	ng	99
35) Caprolactam	9.01	113	28913	33.52	ng	87
36) 4-Chloro-3-methylphenol	9.14	107	94095	31.39	ng	96
37) 2-Methylnaphthalene	9.33	142	234547	34.58	ng	# 91
39) 1,2,4,5-Tetrachlorobenzene	9.50	216	111571	36.57	ng	99
40) Hexachlorocyclopentadiene	9.49	237	107755	69.52	ng	97

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42) 2,4,6-Trichlorophenol	9.60	196	72805	35.07	ng	98
43) 2,4,5-Trichlorophenol	9.64	196	70663	29.54	ng #	83
45) 1,1'-Biphenyl	9.80	154	284676	33.37	ng	96
46) 2-Chloronaphthalene	9.82	162	209365	30.25	ng #	92
47) 2-Nitroaniline	9.91	65	60719	31.96	ng	96
48) Acenaphthylene	10.24	152	354264	30.67	ng	98
49) Dimethylphthalate	10.08	163	264600	33.57	ng #	97
50) 2,6-Dinitrotoluene	10.15	165	50530	31.98	ng #	88
51) Acenaphthene	10.41	154	228722	32.36	ng	90
52) 3-Nitroaniline	10.31	138	49756	27.29	ng #	40
53) 2,4-Dinitrophenol	10.42	184	51075	79.45	ng #	1
54) Dibenzofuran	10.58	168	311243	31.04	ng #	85
55) 4-Nitrophenol	10.46	139	99919	68.57	ng #	55
56) 2,4-Dinitrotoluene	10.55	165	74385	37.10	ng #	79
57) Fluorene	10.94	166	271132	34.07	ng	93
58) 2,3,4,6-Tetrachlorophenol	10.71	232	62026	30.72	ng	99
59) Diethylphthalate	10.78	149	247143	31.14	ng	95
60) 4-Chlorophenyl-phenylether	10.92	204	116026	30.34	ng #	79
61) 4-Nitroaniline	10.93	138	58090	30.85	ng #	44
62) Azobenzene	11.08	77	243206	30.10	ng #	84
64) 4,6-Dinitro-2-methylphenol	10.96	198	34156	34.70	ng	93
65) n-Nitrosodiphenylamine	11.03	169	214222	29.93	ng	91
66) 4-Bromophenyl-phenylether	11.42	248	76750	32.84	ng	93
67) Hexachlorobenzene	11.50	284	82798	31.87	ng #	84
68) Atrazine	11.56	200	73532	32.52	ng	84
69) Pentachlorophenol	11.69	266	81648	55.47	ng	99
70) Phenanthrene	11.92	178	395423	29.71	ng	97
71) Anthracene	11.97	178	392011	30.59	ng	98
72) Carbazole	12.13	167	348313	28.47	ng	96
73) Di-n-butylphthalate	12.46	149	410794	29.06	ng #	98
74) Fluoranthene	13.19	202	424712	29.97	ng	95
76) Benzidine	13.32	184	313345	44.90	ng #	97
77) Pyrene	13.45	202	443535	28.85	ng	97
79) Butylbenzylphthalate	14.17	149	181381	29.38	ng #	96
80) Benzo(a)anthracene	14.91	228	433445	28.50	ng	97
81) 3,3'-Dichlorobenzidine	14.85	252	117437	22.39	ng #	93
82) Chrysene	14.95	228	408931	29.15	ng	96
83) Bis(2-ethylhexyl)phthalate	14.88	149	277344	28.43	ng #	93
84) Di-n-octyl phthalate	15.72	149	464617	27.79	ng	95
85) Indeno(1,2,3-cd)pyrene	18.68	276	450589	25.48	ng #	89
87) Benzo(b)fluoranthene	16.29	252	403207	28.05	ng #	86
88) Benzo(k)fluoranthene	16.32	252	438484	29.96	ng #	88
89) Benzo(a)pyrene	16.75	252	396651	29.06	ng #	88
90) Dibenzo(a,h)anthracene	18.71	278	374996	26.84	ng #	79
91) Benzo(g,h,i)perylene	19.24	276	383309	27.17	ng #	77

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

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