

Data Path : Z:\HPCHEM1\BNA F\DATA\BF070516\
 Data File : BF088711.D
 Acq On : 5 Jul 2016 19:11
 Operator : UM/SJ
 Sample : PB91760BS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB91760BS

Quant Time: Jul 06 02:01:42 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 05 17:15:08 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.78	152	106847	20.00	ng	0.00
21) Naphthalene-d8	8.06	136	430751	20.00	ng	-0.01
38) Acenaphthene-d10	9.82	164	210665	20.00	ng	0.00
63) Phenanthrene-d10	11.29	188	397690	20.00	ng	0.00
75) Chrysene-d12	13.92	240	248932	20.00	ng	0.00
86) Perylene-d12	15.30	264	199232	20.00	ng	0.00

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System Monitoring Compounds

5) 2-Fluorophenol	5.38	112	702437	98.53	ng	0.01
7) Phenol-d6	6.42	99	992380	107.73	ng	0.01
23) Nitrobenzene-d5	7.35	82	713113	86.76	ng	0.00
41) 2,4,6-Tribromophenol	10.60	330	280866	143.57	ng	0.00
44) 2-Fluorobiphenyl	9.14	172	1183722	88.76	ng	0.00
78) Terphenyl-d14	12.87	244	986341	88.25	ng	0.00

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Target Compounds

						Qvalue
3) Pyridine	3.08	79	292730	28.54	ng	92
4) n-Nitrosodimethylamine	3.04	42	125205	31.95	ng	88
6) Aniline	6.43	93	333015	24.81	ng	# 1
8) 2-Chlorophenol	6.56	128	285640	37.28	ng	96
9) Benzaldehyde	6.32	77	243865	39.08	ng	87
10) Phenol	6.43	94	421087	40.05	ng	# 70
11) bis(2-Chloroethyl)ether	6.51	93	263327	33.59	ng	93
12) 1,3-Dichlorobenzene	6.72	146	315364	37.40	ng	97
13) 1,4-Dichlorobenzene	6.80	146	303863	36.31	ng	96
14) 1,2-Dichlorobenzene	6.95	146	302462m	38.61	ng	
15) Benzyl Alcohol	6.92	79	288989	42.62	ng	95
16) 2,2'-oxybis(1-Chloropropan	7.06	45	344147	30.31	ng	90
17) 2-Methylphenol	7.04	107	250613	40.09	ng	96
18) Hexachloroethane	7.29	117	124456	38.44	ng	90
19) n-Nitroso-di-n-propylamine	7.20	70	210054	33.95	ng	91
20) 3+4-Methylphenols	7.20	107	326686	43.55	ng	89
22) Acetophenone	7.19	105	416955	39.88	ng	# 88
24) Nitrobenzene	7.36	77	314274	37.92	ng	# 83
25) Isophorone	7.60	82	631122	41.45	ng	95
26) 2-Nitrophenol	7.68	139	169197	49.79	ng	93
27) 2,4-Dimethylphenol	7.72	122	299679	42.34	ng	90
28) bis(2-Chloroethoxy)methane	7.82	93	370705	37.41	ng	# 95
29) 2,4-Dichlorophenol	7.92	162	239335	41.84	ng	96
30) 1,2,4-Trichlorobenzene	8.00	180	258056	41.11	ng	98
31) Naphthalene	8.08	128	842463	39.67	ng	100
32) Benzoic acid	7.80	122	43269	8.42	ng	91
33) 4-Chloroaniline	8.14	127	189672	20.07	ng	# 92
34) Hexachlorobutadiene	8.20	225	137208	40.82	ng	98
35) Caprolactam	8.50	113	81517m	41.96	ng	
36) 4-Chloro-3-methylphenol	8.62	107	297197	43.93	ng	90
37) 2-Methylnaphthalene	8.78	142	598910	43.80	ng	# 95
39) 1,2,4,5-Tetrachlorobenzene	8.95	216	234964	33.53	ng	# 1
40) Hexachlorocyclopentadiene	8.94	237	269958	78.50	ng	98
42) 2,4,6-Trichlorophenol	9.05	196	179498	38.85	ng	97

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

43) 2,4,5-Trichlorophenol	9.10	196	189061	47.57	nq	99
45) 1,1'-Biphenyl	9.23	154	670766	38.45	nq	98
46) 2-Chloronaphthalene	9.26	162	518120	37.83	nq	96
47) 2-Nitroaniline	9.36	65	207140	42.62	nq	# 82
48) Acenaphthylene	9.68	152	917529	42.11	nq	99
49) Dimethylphthalate	9.54	163	608550	40.55	nq	99
50) 2,6-Dinitrotoluene	9.60	165	134419	40.41	nq	# 47
51) Acenaphthene	9.85	154	557529	41.74	nq	99
52) 3-Nitroaniline	9.76	138	90642	21.44	nq	84
53) 2,4-Dinitrophenol	9.87	184	47958	32.65	nq	# 80
54) Dibenzofuran	10.02	168	800242	45.77	nq	95
55) 4-Nitrophenol	9.94	139	280281	87.22	nq	# 80
56) 2,4-Dinitrotoluene	10.00	165	181757	41.79	nq	# 39
57) Fluorene	10.35	166	548306	40.70	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.14	232	145605	47.35	nq	# 47
59) Diethylphthalate	10.24	149	627914	41.69	nq	98
60) 4-Chlorophenyl-phenylether	10.35	204	261233	41.73	nq	96
61) 4-Nitroaniline	10.38	138	151505	37.23	nq	85
62) Azobenzene	10.51	77	788709	45.91	nq	96
64) 4,6-Dinitro-2-methylphenol	10.41	198	59863	28.14	nq	# 75
65) n-Nitrosodiphenylamine	10.48	169	542076	40.27	nq	100
66) 4-Bromophenyl-phenylether	10.84	248	170799	42.62	nq	93
67) Hexachlorobenzene	10.90	284	164840	37.16	nq	95
68) Atrazine	11.00	200	186228	44.40	nq	93
69) Pentachlorophenol	11.10	266	145533	55.43	nq	97
70) Phenanthrene	11.31	178	816060	37.54	nq	99
71) Anthracene	11.37	178	919733	42.55	nq	99
72) Carbazole	11.52	167	758127	37.26	nq	99
73) Di-n-butylphthalate	11.86	149	1068253	45.14	nq	99
74) Fluoranthene	12.49	202	816764	37.06	nq	97
76) Benzidine	12.62	184	539523	42.88	nq	98
77) Pyrene	12.72	202	810971	38.42	nq	100
79) Butylbenzylphthalate	13.35	149	415034	44.08	nq	# 83
80) Benzo(a)anthracene	13.91	228	686436	42.82	nq	99
81) 3,3'-Dichlorobenzidine	13.87	252	130284	21.62	nq	100
82) Chrysene	13.94	228	672942	42.43	nq	99
83) Bis(2-ethylhexyl)phthalate	13.91	149	483694	45.00	nq	98
84) Di-n-octyl phthalate	14.51	149	758143	41.52	nq	# 100
85) Indeno(1,2,3-cd)pyrene	16.64	276	550589	45.13	nq	# 100
87) Benzo(b)fluoranthene	14.91	252	552901m	44.24	nq	
88) Benzo(k)fluoranthene	14.94	252	451014m	41.45	nq	
89) Benzo(a)pyrene	15.25	252	470021	44.42	nq	97
90) Dibenzo(a,h)anthracene	16.65	278	447540	50.65	nq	99
91) Benzo(g,h,i)perylene	17.04	276	467546	51.13	nq	96

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

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