

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122316\
 Data File : BF091938.D
 Acq On : 23 Dec 2016 21:49
 Operator : UM/SJ
 Sample : H6226-07MS
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 IDW-01-122116MS

Quant Time: Dec 23 23:15:44 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 21 16:17:51 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	189912	20.00	ng	-0.01
21) Naphthalene-d8	8.03	136	822588	20.00	ng	-0.01
38) Acenaphthene-d10	9.78	164	389086	20.00	ng	-0.01
63) Phenanthrene-d10	11.25	188	566822	20.00	ng	-0.01
75) Chrysene-d12	13.89	240	366374	20.00	ng	-0.01
86) Perylene-d12	15.29	264	359359	20.00	ng	-0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	1515131	133.21	ng	0.05
7) Phenol-d6	6.44	99	1849875	133.22	ng	0.03
23) Nitrobenzene-d5	7.31	82	1464354	98.99	ng	-0.01
41) 2,4,6-Tribromophenol	10.58	330	488164	151.60	ng	0.00
44) 2-Fluorobiphenyl	9.10	172	2245774	124.21	ng	-0.01
78) Terphenyl-d14	12.84	244	1471340	97.40	ng	-0.01

Target Compounds						Qvalue
2) 1,4-Dioxane	2.34	88	217106	38.77	ng	98
3) Pyridine	3.02	79	987998	50.55	ng	96
4) n-Nitrosodimethylamine	2.99	42	366756	49.75	ng	100
6) Aniline	6.42	93	177401	9.84	ng	# 52
8) 2-Chlorophenol	6.54	128	598404	46.25	ng	94
9) Benzaldehyde	6.28	77	95891	9.35	ng	96
10) Phenol	6.45	94	785172	49.62	ng	# 73
11) bis(2-Chloroethyl)ether	6.49	93	777140	61.72	ng	98
12) 1,3-Dichlorobenzene	6.68	146	687008	49.74	ng	# 94
13) 1,4-Dichlorobenzene	6.76	146	693449	49.33	ng	96
14) 1,2-Dichlorobenzene	6.91	146	545759	44.74	ng	96
15) Benzyl Alcohol	6.91	79	580306	53.78	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.02	45	936903	49.97	ng	96
17) 2-Methylphenol	7.05	107	540514	51.95	ng	98
18) Hexachloroethane	7.25	117	256174	49.15	ng	96
19) n-Nitroso-di-n-propylamine	7.17	70	540001	58.45	ng	95
20) 3+4-Methylphenols	7.20	107	720156	58.03	ng	# 71
22) Acetophenone	7.16	105	1016690	50.11	ng	# 91
24) Nitrobenzene	7.33	77	766563	48.65	ng	98
25) Isophorone	7.57	82	1231586	45.13	ng	95
26) 2-Nitrophenol	7.65	139	381664	49.12	ng	# 86
27) 2,4-Dimethylphenol	7.71	122	540905	41.97	ng	95
28) bis(2-Chloroethoxy)methane	7.79	93	876907	52.98	ng	100
29) 2,4-Dichlorophenol	7.92	162	546711	50.10	ng	96
30) 1,2,4-Trichlorobenzene	7.97	180	560617	46.88	ng	95
31) Naphthalene	8.05	128	1919069	50.97	ng	99
32) Benzoic acid	7.87	122	369168	38.62	ng	99
33) 4-Chloroaniline	8.12	127	237207	13.97	ng	98
34) Hexachlorobutadiene	8.17	225	309090	48.97	ng	99
35) Caprolactam	8.49	113	143042	39.89	ng	# 61
36) 4-Chloro-3-methylphenol	8.62	107	584143	46.52	ng	100
37) 2-Methylnaphthalene	8.74	142	1157559	55.26	ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.91	216	530502	53.97	ng	98
40) Hexachlorocyclopentadiene	8.90	237	286608	66.19	ng	97

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42) 2,4,6-Trichlorophenol	9.04	196	349101	50.78	nq	99
43) 2,4,5-Trichlorophenol	9.08	196	339646	48.01	nq #	93
45) 1,1'-Biphenyl	9.21	154	1405924	52.77	nq	98
46) 2-Chloronaphthalene	9.23	162	1050009	49.77	nq	95
47) 2-Nitroaniline	9.33	65	284164	37.83	nq	95
48) Acenaphthylene	9.64	152	1523473	46.95	nq	99
49) Dimethylphthalate	9.52	163	1308972	52.60	nq	99
50) 2,6-Dinitrotoluene	9.57	165	280469	51.49	nq #	59
51) Acenaphthene	9.81	154	977667	44.44	nq	98
52) 3-Nitroaniline	9.74	138	113830	16.89	nq	98
53) 2,4-Dinitrophenol	9.86	184	230485	76.05	nq #	91
54) Dibenzofuran	9.98	168	1335210	44.95	nq	97
55) 4-Nitrophenol	9.95	139	307752m	67.24	nq	
56) 2,4-Dinitrotoluene	9.98	165	374640	54.80	nq	92
57) Fluorene	10.33	166	1027059	47.52	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.12	232	273385	48.86	nq	99
59) Diethylphthalate	10.21	149	1293551	53.52	nq	98
60) 4-Chlorophenyl-phenylether	10.33	204	514779	54.40	nq #	77
61) 4-Nitroaniline	10.36	138	186927	26.76	nq	99
62) Azobenzene	10.48	77	1291401	48.53	nq	97
64) 4,6-Dinitro-2-methylphenol	10.38	198	145829	42.55	nq #	39
65) n-Nitrosodiphenylamine	10.44	169	874722	52.01	nq	100
66) 4-Bromophenyl-phenylether	10.81	248	314954	53.42	nq	96
67) Hexachlorobenzene	10.88	284	337435	53.54	nq	97
68) Atrazine	10.97	200	130788	24.11	nq	97
69) Pentachlorophenol	11.08	266	313869	101.50	nq	99
70) Phenanthrene	11.29	178	1639585	53.35	nq	98
71) Anthracene	11.33	178	1454642	60.38	nq	99
72) Carbazole	11.50	167	1362823	Below	Cal	97
73) Di-n-butylphthalate	11.82	149	1688917	59.30	nq	99
74) Fluoranthene	12.47	202	1276777	47.57	nq	98
76) Benzidine	12.59	184	152629	12.81	nq	97
77) Pyrene	12.69	202	1266257	47.11	nq	99
79) Butylbenzylphthalate	13.32	149	676304	55.32	nq	89
80) Benzo(a)anthracene	13.88	228	1056243	51.07	nq	100
81) 3,3'-Dichlorobenzidine	13.85	252	172162	21.95	nq #	95
82) Chrysene	13.92	228	875903	42.22	nq	99
83) Bis(2-ethylhexyl)phthalate	13.88	149	811466	56.36	nq #	99
84) Di-n-octyl phthalate	14.49	149	1348893	50.58	nq	100
85) Indeno(1,2,3-cd)pyrene	16.63	276	993480	55.28	nq	98
87) Benzo(b)fluoranthene	14.90	252	1238130m	55.34	nq	
88) Benzo(k)fluoranthene	14.92	252	852746m	44.70	nq	
89) Benzo(a)pyrene	15.23	252	989693	49.84	nq	99
90) Dibenzo(a,h)anthracene	16.64	278	832154	50.98	nq	96
91) Benzo(g,h,i)perylene	17.03	276	809559	47.70	nq	99

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

