

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101118\
 Data File : BF109788.D
 Acq On : 11 Oct 2018 15:52
 Operator : JU/SJ
 Sample : J5309-15MSD
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S-23MSD

Quant Time: Oct 11 16:38:05 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF100818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 08 16:02:11 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	83373	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	334495	20.00	ng	0.00
38) Acenaphthene-d10	9.72	164	158673	20.00	ng	0.00
63) Phenanthrene-d10	11.20	188	283815	20.00	ng	0.00
75) Chrysene-d12	13.83	240	181176	20.00	ng	0.00
86) Perylene-d12	15.23	264	213507	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.32	112	541593	115.31	ng	0.00
7) Phenol-d6	6.35	99	712131	113.49	ng	0.00
23) Nitrobenzene-d5	7.26	82	437894	72.16	ng	0.00
41) 2,4,6-Tribromophenol	10.52	330	176572	102.13	ng	0.00
44) 2-Fluorobiphenyl	9.05	172	827602	71.83	ng	0.00
78) Terphenyl-d14	12.78	244	676141	72.24	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.40	88	70887	28.756	ng	93
3) Pyridine	3.14	79	134052	26.359	ng	98
4) n-Nitrosodimethylamine	3.06	42	67493	36.768	ng	93
6) Aniline	6.37	93	103206	14.120	ng	# 41
8) 2-Chlorophenol	6.49	128	213851	37.020	ng	99
9) Benzaldehyde	6.25	77	95678	23.696	ng	98
10) Phenol	6.36	94	238397	33.134	ng	# 76
11) bis(2-Chloroethyl)ether	6.43	93	177467	29.751	ng	99
12) 1,3-Dichlorobenzene	6.63	146	217130	32.864	ng	98
13) 1,4-Dichlorobenzene	6.72	146	240336	33.370	ng	99
14) 1,2-Dichlorobenzene	6.86	146	213182	33.705	ng	98
15) Benzyl Alcohol	6.85	79	171140	36.791	ng	100
16) 2,2'-oxybis(1-Chloropropan	6.97	45	266410	33.682	ng	98
17) 2-Methylphenol	6.97	107	166996	36.557	ng	99
18) Hexachloroethane	7.20	117	86340	33.916	ng	96
19) n-Nitroso-di-n-propylamine	7.12	70	152045	34.554	ng	98
20) 3+4-Methylphenols	7.12	107	203931	36.229	ng	93
22) Acetophenone	7.11	105	309671	38.284	ng	# 98
24) Nitrobenzene	7.28	77	220237	34.880	ng	99
25) Isophorone	7.52	82	419197	41.602	ng	100
26) 2-Nitrophenol	7.60	139	103316	34.981	ng	95
27) 2,4-Dimethylphenol	7.65	122	184907	43.355	ng	99
28) bis(2-Chloroethoxy)methane	7.73	93	262081	38.452	ng	98
29) 2,4-Dichlorophenol	7.85	162	178021	37.213	ng	98
30) 1,2,4-Trichlorobenzene	7.92	180	184779	34.892	ng	98
31) Naphthalene	8.00	128	616117	39.828	ng	100
32) Benzoic acid	7.65	122	184907	60.639	ng	# 12
33) 4-Chloroaniline	8.06	127	54760	8.038	ng	97
34) Hexachlorobutadiene	8.12	225	113756	34.548	ng	97
35) Caprolactam	8.42	113	43546	30.480	ng	90
36) 4-Chloro-3-methylphenol	8.55	107	191934	38.008	ng	100
37) 2-Methylnaphthalene	8.69	142	396692	39.148	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.86	216	191215	35.402	ng	99
40) Hexachlorocyclopentadiene	8.85	237	143508	62.419	ng	99

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42) 2,4,6-Trichlorophenol	8.97	196	115100	35.647	nq	99
43) 2,4,5-Trichlorophenol	9.02	196	135310	36.447	nq	98
45) 1,1'-Biphenyl	9.15	154	495417	34.950	nq	100
46) 2-Chloronaphthalene	9.17	162	380059	35.788	nq	99
47) 2-Nitroaniline	9.27	65	122395	38.063	nq	99
48) Acenaphthylene	9.59	152	603364	38.971	nq	99
49) Dimethylphthalate	9.45	163	515098	44.263	nq	99
50) 2,6-Dinitrotoluene	9.52	165	93546	37.084	nq	99
51) Acenaphthene	9.76	154	336793	36.696	nq	99
52) 3-Nitroaniline	9.69	138	42458	15.015	nq	99
53) 2,4-Dinitrophenol	9.82	184	7256	8.915	nq	# 40
54) Dibenzofuran	9.93	168	500180	36.357	nq	98
55) 4-Nitrophenol	9.87	139	77632	50.157	nq	89
56) 2,4-Dinitrotoluene	9.92	165	115395	36.656	nq	96
57) Fluorene	10.27	166	401653	39.095	nq	100
58) 2,3,4,6-Tetrachlorophenol	10.06	232	101497	33.632	nq	97
59) Diethylphthalate	10.15	149	432218	37.485	nq	100
60) 4-Chlorophenyl-phenylether	10.26	204	192657	35.545	nq	99
61) 4-Nitroaniline	10.30	138	86079	32.871	nq	98
62) Azobenzene	10.42	77	434972	37.149	nq	99
64) 4,6-Dinitro-2-methylphenol	10.33	198	25066	18.145	nq	91
65) n-Nitrosodiphenylamine	10.39	169	367347	38.180	nq	99
66) 4-Bromophenyl-phenylether	10.75	248	120815	37.048	nq	97
67) Hexachlorobenzene	10.82	284	128013	36.251	nq	96
68) Atrazine	10.91	200	110114	36.672	nq	99
69) Pentachlorophenol	11.02	266	85158	42.856	nq	99
70) Phenanthrene	11.23	178	535271	37.687	nq	99
71) Anthracene	11.28	178	596290	40.613	nq	99
72) Carbazole	11.44	167	485691	34.108	nq	99
73) Di-n-butylphthalate	11.76	149	599197	36.280	nq	99
74) Fluoranthene	12.41	202	514352	34.665	nq	98
76) Benzidine	12.54	184	113225	16.823	nq	97
77) Pyrene	12.63	202	497885	37.508	nq	98
79) Butylbenzylphthalate	13.25	149	210336	36.548	nq	98
80) Benzo(a)anthracene	13.82	228	379154	37.923	nq	100
81) 3,3'-Dichlorobenzidine	13.79	252	81386	20.216	nq	99
82) Chrysene	13.86	228	400019	38.091	nq	99
83) Bis(2-ethylhexyl)phthalate	13.82	149	279253	39.772	nq	98
84) Di-n-octyl phthalate	14.42	149	514732	46.379	nq	100
85) Indeno(1,2,3-cd)pyrene	16.59	276	515559	68.588	nq	100
87) Benzo(b)fluoranthene	14.83	252	386058	32.720	nq	99
88) Benzo(k)fluoranthene	14.86	252	449378	37.907	nq	98
89) Benzo(a)pyrene	15.17	252	420671	39.289	nq	97
90) Dibenzo(a,h)anthracene	16.59	278	434612	49.424	nq	98
91) Benzo(g,h,i)perylene	16.99	276	430511	49.027	nq	99

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

