

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092217\
 Data File : BF098718.D
 Acq On : 23 Sep 2017 1:31
 Operator : SJ/JU
 Sample : I5340-01MS
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LT-TS-001MS

Quant Time: Sep 23 04:51:24 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 22 10:38:37 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	169089	20.00	ng	0.00
21) Naphthalene-d8	8.21	136	687491	20.00	ng	0.00
38) Acenaphthene-d10	9.97	164	281787	20.00	ng	0.00
63) Phenanthrene-d10	11.46	188	394489	20.00	ng	0.00
75) Chrysene-d12	14.09	240	335779	20.00	ng	0.00
86) Perylene-d12	15.59	264	297167	20.00	ng	0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.55	112	1062883	95.45	ng	0.01
7) Phenol-d6	6.56	99	1306811	96.92	ng	0.01
23) Nitrobenzene-d5	7.49	82	780293	76.99	ng	0.00
41) 2,4,6-Tribromophenol	10.76	330	236213	91.66	ng	0.00
44) 2-Fluorobiphenyl	9.29	172	1226662	69.27	ng	0.00
78) Terphenyl-d14	13.03	244	870856	55.69	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.80	88	149317	28.389	ng	98
3) Pyridine	3.57	79	367192	26.269	ng	99
4) n-Nitrosodimethylamine	3.52	42	201940	32.970	ng	95
6) Aniline	6.59	93	375092	20.999	ng	99
8) 2-Chlorophenol	6.72	128	400325	33.097	ng	98
9) Benzaldehyde	6.48	77	118493	15.667	ng	99
10) Phenol	6.57	94	513188	34.483	ng	95
11) bis(2-Chloroethyl)ether	6.66	93	392511	34.235	ng	98
12) 1,3-Dichlorobenzene	6.87	146	399659	31.281	ng	98
13) 1,4-Dichlorobenzene	6.95	146	404153	31.254	ng	99
14) 1,2-Dichlorobenzene	7.10	146	383240	31.779	ng	97
15) Benzyl Alcohol	7.07	79	344287	35.385	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.20	45	591967	32.400	ng	98
17) 2-Methylphenol	7.17	107	328240	33.303	ng	98
18) Hexachloroethane	7.44	117	126223	28.354	ng	98
19) n-Nitroso-di-n-propylamine	7.34	70	276050	33.047	ng	98
20) 3+4-Methylphenols	7.33	107	413728	32.557	ng	90
22) Acetophenone	7.34	105	546954	32.161	ng	# 96
24) Nitrobenzene	7.51	77	414931	35.446	ng	98
25) Isophorone	7.75	82	749208	33.894	ng	99
26) 2-Nitrophenol	7.83	139	194426	38.750	ng	93
27) 2,4-Dimethylphenol	7.86	122	333540	30.725	ng	97
28) bis(2-Chloroethoxy)methane	7.96	93	485574	35.303	ng	99
29) 2,4-Dichlorophenol	8.06	162	318533	34.111	ng	98
30) 1,2,4-Trichlorobenzene	8.15	180	308749	32.781	ng	99
31) Naphthalene	8.23	128	1081583	33.147	ng	99
32) Benzoic acid	7.94	122	116313	21.569	ng	99
33) 4-Chloroaniline	8.28	127	304918	22.477	ng	98
34) Hexachlorobutadiene	8.35	225	156609	30.864	ng	99
35) Caprolactam	8.64	113	100165	34.399	ng	98
36) 4-Chloro-3-methylphenol	8.75	107	341121	32.003	ng	99
37) 2-Methylnaphthalene	8.93	142	713793	34.069	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.09	216	271091	32.529	ng	99
40) Hexachlorocyclopentadiene	9.07	237	75132	20.746	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.20	196	202506	36.056	nq	98
43) 2,4,5-Trichlorophenol	9.24	196	206584	35.872	nq	95
45) 1,1'-Biphenyl	9.39	154	813502	33.603	nq	99
46) 2-Chloronaphthalene	9.42	162	621743	34.249	nq	97
47) 2-Nitroaniline	9.50	65	203677	39.959	nq	95
48) Acenaphthylene	9.83	152	920660	32.973	nq	100
49) Dimethylphthalate	9.69	163	936059	44.296	nq	100
50) 2,6-Dinitrotoluene	9.74	165	157911	37.869	nq	95
51) Acenaphthene	10.00	154	555651	32.061	nq	99
52) 3-Nitroaniline	9.92	138	157984	31.782	nq	100
53) 2,4-Dinitrophenol	10.02	184	26850	24.836	nq #	80
54) Dibenzofuran	10.17	168	802091	33.893	nq	99
55) 4-Nitrophenol	10.07	139	271886	62.855	nq	91
56) 2,4-Dinitrotoluene	10.15	165	195801	34.953	nq	91
57) Fluorene	10.52	166	593655	31.210	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.29	232	147858	36.087	nq	99
59) Diethylphthalate	10.39	149	699839	33.795	nq	98
60) 4-Chlorophenyl-phenylether	10.50	204	269002	32.282	nq	98
61) 4-Nitroaniline	10.53	138	157860	33.112	nq	97
62) Azobenzene	10.67	77	630630	32.561	nq	96
64) 4,6-Dinitro-2-methylphenol	10.55	198	23785	18.785	nq	90
65) n-Nitrosodiphenylamine	10.62	169	521442	38.401	nq	100
66) 4-Bromophenyl-phenylether	11.00	248	153114	38.440	nq	93
67) Hexachlorobenzene	11.06	284	153641	33.850	nq	95
68) Atrazine	11.14	200	157131	38.469	nq	99
69) Pentachlorophenol	11.25	266	178795	68.920	nq	98
70) Phenanthrene	11.48	178	764087	35.705	nq	98
71) Anthracene	11.53	178	754163	34.562	nq	100
72) Carbazole	11.68	167	715612	33.563	nq	99
73) Di-n-butylphthalate	12.01	149	891093	35.982	nq	99
74) Fluoranthene	12.67	202	743908	34.881	nq	97
77) Pyrene	12.90	202	774796	31.080	nq	99
79) Butylbenzylphthalate	13.51	149	343891	33.345	nq	95
80) Benzo(a)anthracene	14.09	228	695908	34.407	nq	100
81) 3,3'-Dichlorobenzidine	14.04	252	160224	22.199	nq	96
82) Chrysene	14.12	228	645962	33.072	nq	99
83) Bis(2-ethylhexyl)phthalate	14.06	149	496918	34.834	nq	99
84) Di-n-octyl phthalate	14.68	149	884208	41.271	nq	98
85) Indeno(1,2,3-cd)pyrene	17.10	276	474914	31.344	nq	99
87) Benzo(b)fluoranthene	15.14	252	634923	36.841	nq	98
88) Benzo(k)fluoranthene	15.17	252	546254	31.718	nq	99
89) Benzo(a)pyrene	15.52	252	530675	33.906	nq	98
90) Dibenzo(a,h)anthracene	17.11	278	399633	30.191	nq	100
91) Benzo(g,h,i)perylene	17.56	276	378175	28.421	nq	98

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

