

Data Path : Z:\HPCHEM1\BNA F\DATA\BF101017\
 Data File : BF099313.D
 Acq On : 10 Oct 2017 21:22
 Operator : SJ/JU
 Sample : I5745-03MS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HR-06-100617-AMS

Quant Time: Oct 11 00:52:23 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 05 17:20:48 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	86539	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	340976	20.00	ng	0.00
38) Acenaphthene-d10	9.90	164	137723	20.00	ng	0.00
63) Phenanthrene-d10	11.39	188	203504	20.00	ng	0.00
75) Chrysene-d12	14.03	240	172847	20.00	ng	0.00
86) Perylene-d12	15.50	264	153644	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.50	112	650326	119.63	ng	0.02
7) Phenol-d6	6.50	99	713617	106.41	ng	0.00
23) Nitrobenzene-d5	7.43	82	480127	90.30	ng	0.00
41) 2,4,6-Tribromophenol	10.70	330	143049	126.93	ng	0.00
44) 2-Fluorobiphenyl	9.22	172	884809	107.25	ng	0.00
78) Terphenyl-d14	12.97	244	633401	83.34	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.80	88	96552	37.181	ng	# 81
3) Pyridine	3.55	79	224823	32.004	ng	91
4) n-Nitrosodimethylamine	3.48	42	110101	36.960	ng	84
6) Aniline	6.53	93	168374	18.603	ng	99
8) 2-Chlorophenol	6.66	128	276173	44.860	ng	94
9) Benzaldehyde	6.42	77	80693	20.743	ng	91
10) Phenol	6.52	94	291836	38.911	ng	97
11) bis(2-Chloroethyl)ether	6.60	93	258388	44.316	ng	97
12) 1,3-Dichlorobenzene	6.81	146	287990	44.668	ng	97
13) 1,4-Dichlorobenzene	6.89	146	291717	44.471	ng	97
14) 1,2-Dichlorobenzene	7.04	146	274268	45.250	ng	97
15) Benzyl Alcohol	7.01	79	211839	43.059	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.14	45	358890	39.507	ng	93
17) 2-Methylphenol	7.12	107	221451	43.963	ng	97
18) Hexachloroethane	7.37	117	91721	38.900	ng	98
19) n-Nitroso-di-n-propylamine	7.28	70	170889	39.912	ng	94
20) 3+4-Methylphenols	7.27	107	258930	40.633	ng	# 70
22) Acetophenone	7.28	105	365700	44.267	ng	# 97
24) Nitrobenzene	7.45	77	268655	44.543	ng	97
25) Isophorone	7.69	82	492139	44.769	ng	100
26) 2-Nitrophenol	7.77	139	144872	49.950	ng	# 85
27) 2,4-Dimethylphenol	7.80	122	235313	42.961	ng	96
28) bis(2-Chloroethoxy)methane	7.89	93	317908	46.406	ng	99
29) 2,4-Dichlorophenol	8.01	162	224306	47.697	ng	95
30) 1,2,4-Trichlorobenzene	8.09	180	230088	48.919	ng	99
31) Naphthalene	8.17	128	784533	48.989	ng	99
32) Benzoic acid	7.93	122	99120	22.030	ng	92
33) 4-Chloroaniline	8.22	127	53393	7.984	ng	97
34) Hexachlorobutadiene	8.28	225	115499	47.675	ng	99
35) Caprolactam	8.60	113	56836	37.677	ng	89
36) 4-Chloro-3-methylphenol	8.70	107	223588	42.300	ng	96
37) 2-Methylnaphthalene	8.86	142	516674	49.629	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.03	216	204985	49.908	ng	99
40) Hexachlorocyclopentadiene	9.01	237	31397	16.727	ng	99

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42) 2,4,6-Trichlorophenol	9.14	196	140814	48.394	ng	100
43) 2,4,5-Trichlorophenol	9.19	196	149338	51.413	ng	93
45) 1,1'-Biphenyl	9.32	154	585900	49.500	ng	100
46) 2-Chloronaphthalene	9.35	162	454851	51.181	ng	97
47) 2-Nitroaniline	9.45	65	126521	46.494	ng	91
48) Acenaphthylene	9.77	152	667036	49.030	ng	99
49) Dimethylphthalate	9.62	163	537089	51.670	ng	99
50) 2,6-Dinitrotoluene	9.69	165	113005	49.937	ng	# 80
51) Acenaphthene	9.94	154	386236	44.818	ng	99
52) 3-Nitroaniline	9.86	138	63758	24.163	ng	96
53) 2,4-Dinitrophenol	9.97	184	23301	24.164	ng	# 86
54) Dibenzofuran	10.11	168	576509	50.243	ng	98
55) 4-Nitrophenol	10.03	139	143978	64.531	ng	85
56) 2,4-Dinitrotoluene	10.10	165	135742	46.219	ng	# 83
57) Fluorene	10.45	166	434509	47.114	ng	100
58) 2,3,4,6-Tetrachlorophenol	10.23	232	97859	48.771	ng	95
59) Diethylphthalate	10.32	149	485805	47.830	ng	97
60) 4-Chlorophenyl-phenylether	10.44	204	205441	49.590	ng	96
61) 4-Nitroaniline	10.47	138	105817	41.568	ng	89
62) Azobenzene	10.60	77	408830	43.776	ng	90
64) 4,6-Dinitro-2-methylphenol	10.50	198	19577	15.811	ng	87
65) n-Nitrosodiphenylamine	10.56	169	388515	55.108	ng	98
66) 4-Bromophenyl-phenylether	10.93	248	110595	55.520	ng	# 89
67) Hexachlorobenzene	11.00	284	105519	49.597	ng	98
68) Atrazine	11.09	200	113405	53.465	ng	98
69) Pentachlorophenol	11.20	266	105583	79.741	ng	99
70) Phenanthrene	11.42	178	564000	51.776	ng	98
71) Anthracene	11.47	178	572307	51.348	ng	99
72) Carbazole	11.62	167	533622	48.891	ng	99
73) Di-n-butylphthalate	11.94	149	639111	48.648	ng	98
74) Fluoranthene	12.60	202	568313	51.522	ng	98
76) Benzidine	12.73	184	298215	46.647	ng	97
77) Pyrene	12.83	202	588756	44.670	ng	99
79) Butylbenzylphthalate	13.44	149	266322	42.994	ng	95
80) Benzo(a)anthracene	14.02	228	531069	50.741	ng	99
81) 3,3'-Dichlorobenzidine	13.98	252	159105	41.468	ng	99
82) Chrysene	14.06	228	504221	50.602	ng	99
83) Bis(2-ethylhexyl)phthalate	13.99	149	367049	43.034	ng	99
84) Di-n-octyl phthalate	14.60	149	679500	49.475	ng	# 95
85) Indeno(1,2,3-cd)pyrene	17.00	276	400689	50.269	ng	96
87) Benzo(b)fluoranthene	15.07	252	509265	53.910	ng	97
88) Benzo(k)fluoranthene	15.10	252	438341	46.979	ng	98
89) Benzo(a)pyrene	15.44	252	445676	51.396	ng	97
90) Dibenzo(a,h)anthracene	17.01	278	343037	49.432	ng	96
91) Benzo(g,h,i)perylene	17.45	276	328197	47.844	ng	96

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

