

Data Path : Z:\HPCHEM1\BNA F\DATA\BF080717\
 Data File : BF097451.D
 Acq On : 8 Aug 2017 2:17
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Aug 08 04:52:19 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 07 16:02:51 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.41	152	88403	20.00	ng	0.00
21) Naphthalene-d8	7.70	136	366177	20.00	ng	0.00
38) Acenaphthene-d10	9.44	164	144207	20.00	ng	0.00
63) Phenanthrene-d10	10.91	188	211713	20.00	ng	0.00
75) Chrysene-d12	13.54	240	159934	20.00	ng	0.00
86) Perylene-d12	14.87	264	127708	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.00	112	439991	75.03	ng	0.00
7) Phenol-d6	6.09	99	575833	86.01	ng	0.00
23) Nitrobenzene-d5	6.99	82	510767	81.83	ng	0.00
41) 2,4,6-Tribromophenol	10.23	330	85321	74.88	ng	0.00
44) 2-Fluorobiphenyl	8.77	172	741049	87.21	ng	0.00
78) Terphenyl-d14	12.49	244	595853	75.96	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.86	88	90182	36.851	ng	# 73
3) Pyridine	2.46	79	274332	36.609	ng	# 59
4) n-Nitrosodimethylamine	2.42	42	151803	36.566	ng	81
6) Aniline	6.07	93	373944	44.387	ng	97
8) 2-Chlorophenol	6.20	128	269462	42.973	ng	94
9) Benzaldehyde	5.96	77	204464	51.598	ng	98
10) Phenol	6.10	94	348719	43.914	ng	97
11) bis(2-Chloroethyl)ether	6.16	93	257233	40.957	ng	88
12) 1,3-Dichlorobenzene	6.35	146	279285	41.329	ng	99
13) 1,4-Dichlorobenzene	6.43	146	284795	42.527	ng	95
14) 1,2-Dichlorobenzene	6.58	146	246783	42.205	ng	98
15) Benzyl Alcohol	6.58	79	196258	46.302	ng	99
16) 2,2'-oxybis(1-Chloropropan	6.70	45	514382	40.833	ng	60
17) 2-Methylphenol	6.70	107	201146	41.563	ng	# 87
18) Hexachloroethane	6.92	117	91441	37.323	ng	# 79
19) n-Nitroso-di-n-propylamine	6.85	70	185458	42.086	ng	# 83
20) 3+4-Methylphenols	6.86	107	282062	44.625	ng	# 62
22) Acetophenone	6.84	105	361541	41.114	ng	96
24) Nitrobenzene	7.01	77	277514	42.689	ng	91
25) Isophorone	7.25	82	471186	38.546	ng	97
26) 2-Nitrophenol	7.32	139	136294	38.752	ng	# 84
27) 2,4-Dimethylphenol	7.39	122	233769	40.293	ng	96
28) bis(2-Chloroethoxy)methane	7.47	93	318794	39.963	ng	99
29) 2,4-Dichlorophenol	7.58	162	212644	42.545	ng	98
30) 1,2,4-Trichlorobenzene	7.65	180	187955	39.453	ng	97
31) Naphthalene	7.72	128	706000	40.901	ng	100
32) Benzoic acid	7.56	122	144516	33.358	ng	96
33) 4-Chloroaniline	7.78	127	295707	42.102	ng	98
34) Hexachlorobutadiene	7.84	225	100494	39.790	ng	97
35) Caprolactam	8.16	113	66070	41.225	ng	# 43
36) 4-Chloro-3-methylphenol	8.29	107	222832	40.866	ng	90
37) 2-Methylnaphthalene	8.41	142	431829	39.512	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.58	216	171284	44.515	ng	99
40) Hexachlorocyclopentadiene	8.56	237	3437	8.935	ng	100

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42) 2,4,6-Trichlorophenol	8.70	196	114803	41.172	nq	99
43) 2,4,5-Trichlorophenol	8.76	196	110454	39.576	nq	95
45) 1,1'-Biphenyl	8.87	154	534637	43.405	nq	98
46) 2-Chloronaphthalene	8.89	162	383063	42.262	nq	# 92
47) 2-Nitroaniline	9.00	65	149127	45.334	nq	96
48) Acenaphthylene	9.30	152	571621	41.086	nq	99
49) Dimethylphthalate	9.18	163	472746	43.170	nq	99
50) 2,6-Dinitrotoluene	9.25	165	95153	40.968	nq	# 68
51) Acenaphthene	9.47	154	343175	41.876	nq	99
52) 3-Nitroaniline	9.42	138	124704	43.256	nq	91
53) 2,4-Dinitrophenol	9.56	184	7069	6.694	nq	# 71
54) Dibenzofuran	9.65	168	467282	42.808	nq	96
55) 4-Nitrophenol	9.65	139	309529	180.544	nq	# 34
56) 2,4-Dinitrotoluene	9.66	165	119233	44.656	nq	# 78
57) Fluorene	9.99	166	349591	42.611	nq	100
58) 2,3,4,6-Tetrachlorophenol	9.78	232	72207	37.470	nq	97
59) Diethylphthalate	9.88	149	438238	43.620	nq	99
60) 4-Chlorophenyl-phenylether	9.98	204	146923	43.368	nq	96
61) 4-Nitroaniline	10.03	138	109674	40.037	nq	92
62) Azobenzene	10.14	77	385113	41.320	nq	93
64) 4,6-Dinitro-2-methylphenol	10.07	198	15653	11.898	nq	89
65) n-Nitrosodiphenylamine	10.10	169	330337	44.844	nq	99
66) 4-Bromophenyl-phenylether	10.46	248	91810	43.454	nq	96
67) Hexachlorobenzene	10.53	284	96912	40.903	nq	# 86
68) Atrazine	10.64	200	79894	37.823	nq	95
69) Pentachlorophenol	10.74	266	55976	57.100	nq	98
70) Phenanthrene	10.93	178	468608	41.310	nq	98
71) Anthracene	10.99	178	483978	41.656	nq	98
72) Carbazole	11.15	167	475587	41.622	nq	99
73) Di-n-butylphthalate	11.49	149	559543	39.616	nq	99
74) Fluoranthene	12.12	202	460518	41.440	nq	98
76) Benzidine	12.25	184	222840	37.551	nq	# 93
77) Pyrene	12.34	202	467749	35.724	nq	99
79) Butylbenzylphthalate	12.97	149	245655	36.884	nq	# 82
80) Benzo(a)anthracene	13.53	228	411425	39.757	nq	98
81) 3,3'-Dichlorobenzidine	13.50	252	167170	42.837	nq	# 96
82) Chrysene	13.56	228	392513	38.844	nq	99
83) Bis(2-ethylhexyl)phthalate	13.54	149	311266	35.794	nq	100
84) Di-n-octyl phthalate	14.14	149	592467	37.126	nq	# 92
85) Indeno(1,2,3-cd)pyrene	16.07	276	249031	28.741	nq	99
87) Benzo(b)fluoranthene	14.53	252	329363	42.904	nq	97
88) Benzo(k)fluoranthene	14.55	252	310464	42.440	nq	97
89) Benzo(a)pyrene	14.83	252	283073	40.289	nq	99
90) Dibenzo(a,h)anthracene	16.07	278	208453	36.179	nq	97
91) Benzo(g,h,i)perylene	16.42	276	210572	34.851	nq	98

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

