

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050815\
 Data File : BF079037.D
 Acq On : 8 May 2015 14:48
 Operator : TP/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: May 08 22:58:02 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 08 22:44:53 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.32	152	70783	20.00	ng	0.00
21) Naphthalene-d8	8.90	136	292270	20.00	ng	0.00
38) Acenaphthene-d10	11.07	164	135025	20.00	ng	0.00
63) Phenanthrene-d10	12.93	188	251333	20.00	ng	0.00
75) Chrysene-d12	16.25	240	240923	20.00	ng	0.00
86) Perylene-d12	18.08	264	245225	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.62	112	340891	82.90	ng	0.00
7) Phenol-d6	6.89	99	449460	82.69	ng	0.00
23) Nitrobenzene-d5	8.01	82	415207	81.65	ng	0.00
41) 2,4,6-Tribromophenol	12.06	330	118967	81.44	ng	0.00
44) 2-Fluorobiphenyl	10.25	172	818877	84.49	ng	0.00
78) Terphenyl-d14	14.93	244	859585	85.42	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.27	88	73548	41.14	ng	# 19
3) Pyridine	3.03	79	208322	39.81	ng	99
4) n-Nitrosodimethylamine	2.95	42	84348	37.62	ng	# 85
6) Aniline	6.91	93	315828	41.17	ng	96
8) 2-Chlorophenol	7.05	128	191887	41.14	ng	92
9) Benzaldehyde	6.78	77	142773	38.99	ng	99
10) Phenol	6.90	94	265404	42.09	ng	88
11) bis(2-Chloroethyl)ether	7.02	93	182382	39.46	ng	87
12) 1,3-Dichlorobenzene	7.24	146	208669	39.81	ng	96
13) 1,4-Dichlorobenzene	7.35	146	215815	40.00	ng	98
14) 1,2-Dichlorobenzene	7.53	146	203793	40.58	ng	99
15) Benzyl Alcohol	7.51	79	168184	41.68	ng	93
16) 2,2'-oxybis(1-Chloropropan	7.68	45	276079	39.04	ng	100
17) 2-Methylphenol	7.64	107	150363	40.06	ng	96
18) Hexachloroethane	7.94	117	71878	38.68	ng	# 81
19) n-Nitroso-di-n-propylamine	7.84	70	150032	40.87	ng	# 86
20) 3+4-Methylphenols	7.84	107	204070	40.81	ng	# 56
22) Acetophenone	7.83	105	260939	39.52	ng	# 96
24) Nitrobenzene	8.03	77	195956	38.87	ng	97
25) Isophorone	8.34	82	382300	40.55	ng	# 90
26) 2-Nitrophenol	8.43	139	97517	46.74	ng	95
27) 2,4-Dimethylphenol	8.49	122	171945	41.18	ng	99
28) bis(2-Chloroethoxy)methane	8.62	93	242996	41.81	ng	99
29) 2,4-Dichlorophenol	8.73	162	160499	43.23	ng	99
30) 1,2,4-Trichlorobenzene	8.83	180	164671	40.38	ng	97
31) Naphthalene	8.92	128	560825	40.27	ng	99
32) Benzoic acid	8.60	122	108835	42.46	ng	93
33) 4-Chloroaniline	9.00	127	241278	40.95	ng	# 92
34) Hexachlorobutadiene	9.08	225	88760	37.79	ng	99
35) Caprolactam	9.43	113	49606	41.32	ng	96
36) 4-Chloro-3-methylphenol	9.61	107	166499	42.70	ng	85
37) 2-Methylnaphthalene	9.78	142	376058	38.97	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.99	216	153891	40.47	ng	# 100
40) Hexachlorocyclopentadiene	9.98	237	62631	32.75	ng	95

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42) 2,4,6-Trichlorophenol	10.14	196	112344	42.97	nq	94
43) 2,4,5-Trichlorophenol	10.18	196	116417	44.04	nq	97
45) 1,1'-Biphenyl	10.37	154	437988	40.80	nq	99
46) 2-Chloronaphthalene	10.39	162	341750	40.82	nq	99
47) 2-Nitroaniline	10.51	65	103446	42.64	nq	89
48) Acenaphthylene	10.90	152	577192	41.99	nq	100
49) Dimethylphthalate	10.75	163	398658	40.23	nq	99
50) 2,6-Dinitrotoluene	10.83	165	81104	41.47	nq	# 63
51) Acenaphthene	11.12	154	332254	40.53	nq	99
52) 3-Nitroaniline	11.03	138	101986	41.75	nq	# 81
53) 2,4-Dinitrophenol	11.17	184	23377	42.06	nq	92
54) Dibenzofuran	11.34	168	492592	41.60	nq	96
55) 4-Nitrophenol	11.25	139	79198	40.96	nq	89
56) 2,4-Dinitrotoluene	11.33	165	107543	41.60	nq	85
57) Fluorene	11.76	166	397991	40.95	nq	100
58) 2,3,4,6-Tetrachlorophenol	11.49	232	85233	39.46	nq	# 100
59) Diethylphthalate	11.63	149	394438	40.18	nq	99
60) 4-Chlorophenyl-phenylether	11.77	204	167838	38.92	nq	92
61) 4-Nitroaniline	11.78	138	103607	42.39	nq	91
62) Azobenzene	11.97	77	388816	40.19	nq	95
64) 4,6-Dinitro-2-methylphenol	11.83	198	43294	43.11	nq	87
65) n-Nitrosodiphenylamine	11.92	169	340605	40.69	nq	99
66) 4-Bromophenyl-phenylether	12.38	248	105648	40.33	nq	91
67) Hexachlorobenzene	12.43	284	118498	39.58	nq	97
68) Atrazine	12.58	200	96440	39.62	nq	98
69) Pentachlorophenol	12.69	266	47771	31.78	nq	97
70) Phenanthrene	12.96	178	563385	40.07	nq	99
71) Anthracene	13.02	178	563687	40.87	nq	100
72) Carbazole	13.22	167	541276	41.17	nq	99
73) Di-n-butylphthalate	13.67	149	634130	40.22	nq	99
74) Fluoranthene	14.43	202	584777	39.91	nq	97
76) Benzidine	14.62	184	350547	53.99	nq	98
77) Pyrene	14.72	202	608966	43.86	nq	99
79) Butylbenzylphthalate	15.54	149	268305	44.21	nq	95
80) Benzo(a)anthracene	16.23	228	547827	41.52	nq	99
81) 3,3'-Dichlorobenzidine	16.21	252	205124	43.65	nq	# 97
82) Chrysene	16.29	228	511962	40.35	nq	98
83) Bis(2-ethylhexyl)phthalate	16.29	149	376158	40.92	nq	# 99
84) Di-n-octyl phthalate	17.12	149	657551	40.62	nq	# 100
85) Indeno(1,2,3-cd)pyrene	19.58	276	679472	42.03	nq	# 100
87) Benzo(b)fluoranthene	17.60	252	608412	45.33	nq	99
88) Benzo(k)fluoranthene	17.63	252	468696	36.07	nq	98
89) Benzo(a)pyrene	18.01	252	522545	42.28	nq	99
90) Dibenzo(a,h)anthracene	19.61	278	537667	40.93	nq	99
91) Benzo(g,h,i)perylene	20.02	276	552042	41.93	nq	98

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

