

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040715\
 Data File : BF078338.D
 Acq On : 8 Apr 2015 12:44
 Operator : TP/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040EC

Quant Time: Apr 08 16:06:26 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 03 14:42:25 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	47956	20.00	ng	-0.05
21) Naphthalene-d8	8.51	136	198038	20.00	ng	-0.05
38) Acenaphthene-d10	10.67	164	97542	20.00	ng	-0.06
63) Phenanthrene-d10	12.50	188	187044	20.00	ng	-0.06
75) Chrysene-d12	15.79	240	197118	20.00	ng	-0.05
86) Perylene-d12	17.57	264	206819	20.00	ng	0.02

System Monitoring Compounds						
5) 2-Fluorophenol	5.20	112	231930	79.74	ng	-0.07
7) Phenol-d6	6.53	99	298851	81.99	ng	-0.03
23) Nitrobenzene-d5	7.63	82	255299	78.14	ng	-0.05
41) 2,4,6-Tribromophenol	11.65	330	72481	89.60	ng	-0.06
44) 2-Fluorobiphenyl	9.86	172	530080	77.68	ng	-0.05
78) Terphenyl-d14	14.50	244	662994	76.00	ng	-0.06

Target Compounds				Qvalue		
2) 1,4-Dioxane	1.74	88	51671m	39.29	ng	
3) Pyridine	2.34	79	153486	44.55	ng	# 87
4) n-Nitrosodimethylamine	2.29	42	53388m	40.26	ng	
6) Aniline	6.52	93	208030	40.24	ng	92
8) 2-Chlorophenol	6.67	128	135215	39.32	ng	94
9) Benzaldehyde	6.37	77	89211	36.93	ng	94
10) Phenol	6.54	94	179284	41.05	ng	95
11) bis(2-Chloroethyl)ether	6.64	93	126183	40.17	ng	98
12) 1,3-Dichlorobenzene	6.85	146	151243	40.03	ng	# 93
13) 1,4-Dichlorobenzene	6.96	146	153139	40.27	ng	96
14) 1,2-Dichlorobenzene	7.14	146	138579	38.87	ng	95
15) Benzyl Alcohol	7.13	79	105914	41.35	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.31	45	166035	39.63	ng	96
17) 2-Methylphenol	7.30	107	104007	38.95	ng	96
18) Hexachloroethane	7.55	117	51551	40.20	ng	# 81
19) n-Nitroso-di-n-propylamine	7.47	70	101501	42.20	ng	95
20) 3+4-Methylphenols	7.49	107	143670	40.33	ng	91
22) Acetophenone	7.45	105	182447	39.54	ng	# 88
24) Nitrobenzene	7.65	77	134225	39.30	ng	# 80
25) Isophorone	7.96	82	259764	41.89	ng	100
26) 2-Nitrophenol	8.05	139	66652	40.95	ng	98
27) 2,4-Dimethylphenol	8.13	122	118541	39.17	ng	95
28) bis(2-Chloroethoxy)methane	8.25	93	152101	38.25	ng	100
29) 2,4-Dichlorophenol	8.36	162	111013	40.32	ng	99
30) 1,2,4-Trichlorobenzene	8.44	180	120447	38.15	ng	# 97
31) Naphthalene	8.53	128	392097	38.04	ng	99
32) Benzoic acid	8.27	122	67540	47.52	ng	96
33) 4-Chloroaniline	8.62	127	174132	40.71	ng	99
34) Hexachlorobutadiene	8.69	225	68926	39.76	ng	97
35) Caprolactam	9.07	113	36338	44.21	ng	88
36) 4-Chloro-3-methylphenol	9.25	107	117216	42.19	ng	85
37) 2-Methylnaphthalene	9.40	142	271556	38.80	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.61	216	120399	39.84	ng	100
40) Hexachlorocyclopentadiene	9.58	237	28237	26.67	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.76	196	82871	43.48	nq	99
43) 2,4,5-Trichlorophenol	9.80	196	83483	41.92	nq	98
45) 1,1'-Biphenyl	9.98	154	318591	39.93	nq	99
46) 2-Chloronaphthalene	10.00	162	256283	40.83	nq	97
47) 2-Nitroaniline	10.13	65	68338	44.00	nq	96
48) Acenaphthylene	10.50	152	412946	40.74	nq	99
49) Dimethylphthalate	10.38	163	299146	40.82	nq	# 97
50) 2,6-Dinitrotoluene	10.45	165	64971	43.09	nq	97
51) Acenaphthene	10.72	154	235043	41.18	nq	99
52) 3-Nitroaniline	10.65	138	74763	43.61	nq	96
53) 2,4-Dinitrophenol	10.78	184	6146	21.02	nq	94
54) Dibenzofuran	10.93	168	366264	42.02	nq	95
55) 4-Nitrophenol	10.89	139	48797	40.08	nq	# 77
56) 2,4-Dinitrotoluene	10.94	165	85701	43.89	nq	96
57) Fluorene	11.36	166	296187	41.92	nq	99
58) 2,3,4,6-Tetrachlorophenol	11.09	232	67173	45.50	nq	98
59) Diethylphthalate	11.25	149	299138	42.34	nq	99
60) 4-Chlorophenyl-phenylether	11.37	204	134777	41.46	nq	# 82
61) 4-Nitroaniline	11.40	138	82219	47.44	nq	99
62) Azobenzene	11.56	77	263704	40.89	nq	95
64) 4,6-Dinitro-2-methylphenol	11.44	198	14506	21.40	nq	89
65) n-Nitrosodiphenylamine	11.52	169	254703	39.37	nq	99
66) 4-Bromophenyl-phenylether	11.97	248	82854	41.83	nq	# 85
67) Hexachlorobenzene	12.02	284	83847	38.63	nq	# 86
68) Atrazine	12.20	200	80465	42.94	nq	98
69) Pentachlorophenol	12.28	266	42623	47.74	nq	98
70) Phenanthrene	12.53	178	427979	39.56	nq	100
71) Anthracene	12.60	178	437934	41.08	nq	99
72) Carbazole	12.81	167	402184	40.25	nq	99
73) Di-n-butylphthalate	13.28	149	509847	45.04	nq	99
74) Fluoranthene	14.00	202	474206	41.56	nq	98
76) Benzidine	14.19	184	263224	34.68	nq	99
77) Pyrene	14.27	202	493210	39.86	nq	99
79) Butylbenzylphthalate	15.13	149	230857	48.95	nq	# 81
80) Benzo(a)anthracene	15.78	228	486814	41.51	nq	99
81) 3,3'-Dichlorobenzidine	15.77	252	171979	43.78	nq	99
82) Chrysene	15.83	228	456712	41.45	nq	99
83) Bis(2-ethylhexyl)phthalate	15.86	149	344989	46.64	nq	# 95
84) Di-n-octyl phthalate	16.68	149	602385	49.60	nq	# 100
85) Indeno(1,2,3-cd)pyrene	18.89	276	552931	44.22	nq	# 85
87) Benzo(b)fluoranthene	17.12	252	466773m	36.52	nq	
88) Benzo(k)fluoranthene	17.15	252	469927	41.37	nq	# 95
89) Benzo(a)pyrene	17.51	252	447225	40.09	nq	# 94
90) Dibenzo(a,h)anthracene	18.92	278	449268	40.23	nq	99
91) Benzo(g,h,i)perylene	19.27	276	441562	39.43	nq	99

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

