

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011515\
 Data File : BF076875.D
 Acq On : 15 Jan 2015 9:52
 Operator : IZ / TP
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Jan 15 13:03:57 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF011415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 14 23:27:58 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.02	152	33160	20.00	ng	0.00
21) Naphthalene-d8	10.93	136	136047	20.00	ng	0.00
38) Acenaphthene-d10	15.07	164	71219	20.00	ng	-0.01
63) Phenanthrene-d10	17.82	188	123261	20.00	ng	0.00
75) Chrysene-d12	21.31	240	126475	20.00	ng	0.00
86) Perylene-d12	22.97	264	114195	20.00	ng	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.09	112	169428	84.01	ng	0.00
7) Phenol-d6	7.42	99	214791	84.42	ng	0.00
23) Nitrobenzene-d5	9.32	82	211091	85.96	ng	0.00
41) 2,4,6-Tribromophenol	16.72	330	50664	87.64	ng	0.00
44) 2-Fluorobiphenyl	13.59	172	415052	88.69	ng	0.00
78) Terphenyl-d14	20.04	244	462986	84.28	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	1.32	88	35476	41.08	ng	98
3) Pyridine	1.80	79	96039	42.98	ng	92
4) n-Nitrosodimethylamine	1.76	42	40101	36.23	ng	89
6) Aniline	7.30	93	141990	40.33	ng	96
8) 2-Chlorophenol	7.56	128	98792	42.49	ng	98
9) Benzaldehyde	7.03	77	51492	34.58	ng	95
10) Phenol	7.44	94	112261	41.55	ng	94
11) bis(2-Chloroethyl)ether	7.56	93	88181	41.71	ng	91
12) 1,3-Dichlorobenzene	7.86	146	110682	41.84	ng	99
13) 1,4-Dichlorobenzene	8.06	146	115408	41.14	ng	99
14) 1,2-Dichlorobenzene	8.38	146	107644	41.65	ng	97
15) Benzyl Alcohol	8.45	79	85237	41.40	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.80	45	113777	40.04	ng	85
17) 2-Methylphenol	8.79	107	76330	40.17	ng	98
18) Hexachloroethane	9.12	117	40763	41.78	ng	94
19) n-Nitroso-di-n-propylamine	9.08	70	71175	39.96	ng	94
20) 3+4-Methylphenols	9.18	107	102289	40.66	ng	94
22) Acetophenone	9.01	105	142679	42.82	ng	96
24) Nitrobenzene	9.36	77	107662	43.04	ng	98
25) Isophorone	9.96	82	187867	42.01	ng	97
26) 2-Nitrophenol	10.09	139	51682	40.14	ng	# 84
27) 2,4-Dimethylphenol	10.37	122	81531	42.15	ng	97
28) bis(2-Chloroethoxy)methane	10.57	93	106717	41.98	ng	97
29) 2,4-Dichlorophenol	10.70	162	77254	42.85	ng	96
30) 1,2,4-Trichlorobenzene	10.84	180	87185	43.54	ng	94
31) Naphthalene	10.97	128	294806	42.09	ng	99
32) Benzoic acid	10.72	122	44851m	41.85	ng	
33) 4-Chloroaniline	11.21	127	119692	42.29	ng	98
34) Hexachlorobutadiene	11.34	225	51267	43.15	ng	97
35) Caprolactam	12.06	113	20642	38.89	ng	# 82
36) 4-Chloro-3-methylphenol	12.52	107	87177	42.23	ng	99
37) 2-Methylnaphthalene	12.63	142	204168	42.68	ng	95
39) 1,2,4,5-Tetrachlorobenzene	13.03	216	76556	43.48	ng	98
40) Hexachlorocyclopentadiene	13.02	237	37927	41.12	ng	99

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011515\
 Data File : BF076875.D
 Acq On : 15 Jan 2015 9:52
 Operator : IZ / TP
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Jan 15 13:03:57 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF011415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 14 23:27:58 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.39	196	56462	44.01	nq	92
43) 2,4,5-Trichlorophenol	13.47	196	54534	41.68	nq	94
45) 1,1'-Biphenyl	13.79	154	232787	44.09	nq	98
46) 2-Chloronaphthalene	13.76	162	188090	43.29	nq	96
47) 2-Nitroaniline	14.11	65	57529	43.65	nq	91
48) Acenaphthylene	14.72	152	314919	42.97	nq	100
49) Dimethylphthalate	14.65	163	216582	42.88	nq	97
50) 2,6-Dinitrotoluene	14.76	165	44764	44.36	nq	94
51) Acenaphthene	15.15	154	177169	42.61	nq	95
52) 3-Nitroaniline	15.10	138	54805	41.37	nq	86
53) 2,4-Dinitrophenol	15.39	184	14040	32.02	nq	97
54) Dibenzofuran	15.57	168	263408	43.49	nq	99
55) 4-Nitrophenol	15.68	139	32770	41.39	nq	98
56) 2,4-Dinitrotoluene	15.68	165	61906	40.67	nq	96
57) Fluorene	16.28	166	227924	44.41	nq	98
58) 2,3,4,6-Tetrachlorophenol	15.89	232	41692	43.76	nq	# 95
59) Diethylphthalate	16.27	149	230935	45.24	nq	98
60) 4-Chlorophenyl-phenylether	16.36	204	91364	43.52	nq	95
61) 4-Nitroaniline	16.40	138	54330	42.54	nq	95
62) Azobenzene	16.64	77	240234	45.17	nq	98
64) 4,6-Dinitro-2-methylphenol	16.47	198	31337	39.17	nq	78
65) n-Nitrosodiphenylamine	16.60	169	181871	41.38	nq	96
66) 4-Bromophenyl-phenylether	17.19	248	52611	42.13	nq	96
67) Hexachlorobenzene	17.21	284	55899	42.48	nq	92
68) Atrazine	17.56	200	60764	45.56	nq	99
69) Pentachlorophenol	17.57	266	21973	37.94	nq	96
70) Phenanthrene	17.85	178	325718	42.37	nq	99
71) Anthracene	17.92	178	315312	42.07	nq	99
72) Carbazole	18.21	167	315377	43.38	nq	99
73) Di-n-butylphthalate	18.79	149	381392	45.32	nq	98
74) Fluoranthene	19.49	202	343676	43.63	nq	98
76) Benzidine	19.72	184	154452	38.16	nq	99
77) Pyrene	19.77	202	354853	40.94	nq	99
79) Butylbenzylphthalate	20.70	149	173011	44.07	nq	98
80) Benzo(a)anthracene	21.29	228	324099	40.84	nq	100
81) 3,3'-Dichlorobenzidine	21.31	252	105105	41.68	nq	# 98
82) Chrysene	21.34	228	307700	42.52	nq	100
83) Bis(2-ethylhexyl)phthalate	21.43	149	255471	47.04	nq	# 95
84) Di-n-octyl phthalate	22.21	149	453943m	48.16	nq	
85) Indeno(1,2,3-cd)pyrene	24.11	276	341329m	44.51	nq	
87) Benzo(b)fluoranthene	22.55	252	361359	46.98	nq	98
88) Benzo(k)fluoranthene	22.59	252	256086m	38.11	nq	
89) Benzo(a)pyrene	22.91	252	293973	43.33	nq	# 96
90) Dibenzo(a,h)anthracene	24.13	278	272983	43.85	nq	# 95
91) Benzo(g,h,i)perylene	24.40	276	273629	44.22	nq	# 94

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

