

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070518\
 Data File : BF107144.D
 Acq On : 6 Jul 2018 1:27
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Jul 06 04:14:25 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 29 14:38:29 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	49330	20.00	ng	-0.01
21) Naphthalene-d8	8.23	136	195434	20.00	ng	-0.02
38) Acenaphthene-d10	9.99	164	90893	20.00	ng	-0.02
63) Phenanthrene-d10	11.48	188	149221	20.00	ng	-0.02
75) Chrysene-d12	14.14	240	145596	20.00	ng	-0.03
86) Perylene-d12	15.67	264	116945	20.00	ng	-0.05

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.58	112	235054	80.69	ng	-0.02
7) Phenol-d6	6.60	99	284270	78.14	ng	-0.02
23) Nitrobenzene-d5	7.52	82	261041	74.23	ng	-0.02
41) 2,4,6-Tribromophenol	10.79	330	72074	68.42	ng	-0.01
44) 2-Fluorobiphenyl	9.31	172	482531	90.88	ng	-0.02
78) Terphenyl-d14	13.07	244	461643	76.80	ng	-0.02

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.74	88	53873	35.970	ng	97
3) Pyridine	3.54	79	147277	36.258	ng	96
4) n-Nitrosodimethylamine	3.51	42	58798	34.082	ng	84
6) Aniline	6.61	93	182283	36.937	ng	# 74
8) 2-Chlorophenol	6.74	128	126522	39.597	ng	97
9) Benzaldehyde	6.51	77	87604	34.060	ng	94
10) Phenol	6.61	94	156555	37.707	ng	75
11) bis(2-Chloroethyl)ether	6.68	93	129024	37.643	ng	98
12) 1,3-Dichlorobenzene	6.89	146	149051	39.828	ng	97
13) 1,4-Dichlorobenzene	6.97	146	149984	39.641	ng	98
14) 1,2-Dichlorobenzene	7.12	146	140398	40.490	ng	97
15) Benzyl Alcohol	7.10	79	106511	36.461	ng	96
16) 2,2'-oxybis(1-Chloropropan	7.22	45	155848	34.655	ng	# 46
17) 2-Methylphenol	7.21	107	110216	40.307	ng	# 93
18) Hexachloroethane	7.45	117	42809	32.175	ng	# 89
19) n-Nitroso-di-n-propylamine	7.35	70	91211	38.035	ng	94
20) 3+4-Methylphenols	7.37	107	132475	44.831	ng	90
22) Acetophenone	7.35	105	176271	40.315	ng	# 98
24) Nitrobenzene	7.54	77	132965	37.034	ng	99
25) Isophorone	7.77	82	225697	36.421	ng	97
26) 2-Nitrophenol	7.85	139	60835	35.893	ng	99
27) 2,4-Dimethylphenol	7.88	122	101467	38.732	ng	99
28) bis(2-Chloroethoxy)methane	7.97	93	155686	37.418	ng	99
29) 2,4-Dichlorophenol	8.10	162	108812	39.673	ng	92
30) 1,2,4-Trichlorobenzene	8.17	180	125129	41.414	ng	96
31) Naphthalene	8.25	128	358523	39.238	ng	100
32) Benzoic acid	8.01	122	45846	26.917	ng	99
33) 4-Chloroaniline	8.31	127	145963	38.072	ng	99
34) Hexachlorobutadiene	8.36	225	84148	42.742	ng	99
35) Caprolactam	8.68	113	31768	35.533	ng	98
36) 4-Chloro-3-methylphenol	8.80	107	106948	37.046	ng	96
37) 2-Methylnaphthalene	8.94	142	249427	41.185	ng	97
39) 1,2,4,5-Tetrachlorobenzene	9.11	216	131443	42.941	ng	# 95
42) 2,4,6-Trichlorophenol	9.23	196	74921	36.822	ng	93

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070518\
 Data File : BF107144.D
 Acq On : 6 Jul 2018 1:27
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Jul 06 04:14:25 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 29 14:38:29 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)

43) 2,4,5-Trichlorophenol	9.28	196	75690	35.650	nq	# 87
45) 1,1'-Biphenyl	9.41	154	313325	43.255	nq	97
46) 2-Chloronaphthalene	9.44	162	217004	38.058	nq	97
47) 2-Nitroaniline	9.54	65	68257	35.599	nq	96
48) Acenaphthylene	9.85	152	334920	37.205	nq	99
49) Dimethylphthalate	9.70	163	263196	38.608	nq	98
50) 2,6-Dinitrotoluene	9.78	165	55322	38.324	nq	# 85
51) Acenaphthene	10.02	154	210310	37.100	nq	98
52) 3-Nitroaniline	9.95	138	60397	36.359	nq	97
53) 2,4-Dinitrophenol	10.08	184	1655	7.410	nq	# 19
54) Dibenzofuran	10.20	168	304361	39.211	nq	96
55) 4-Nitrophenol	10.14	139	21353	18.416	nq	96
56) 2,4-Dinitrotoluene	10.18	165	64028	33.981	nq	# 96
57) Fluorene	10.54	166	236622	38.415	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.32	232	53226	31.537	nq	# 81
59) Diethylphthalate	10.40	149	247985	37.690	nq	# 93
60) 4-Chlorophenyl-phenylether	10.52	204	130018	40.342	nq	# 86
61) 4-Nitroaniline	10.57	138	56302	34.152	nq	97
62) Azobenzene	10.69	77	208014	32.104	nq	91
64) 4,6-Dinitro-2-methylphenol	10.60	198	9213	9.988	nq	# 74
65) n-Nitrosodiphenylamine	10.65	169	214116	42.660	nq	100
66) 4-Bromophenyl-phenylether	11.02	248	80036	44.942	nq	# 80
67) Hexachlorobenzene	11.09	284	78060	41.879	nq	# 86
68) Atrazine	11.17	200	60218	37.740	nq	95
69) Pentachlorophenol	11.30	266	19887	21.383	nq	98
70) Phenanthrene	11.51	178	299888	41.667	nq	99
71) Anthracene	11.56	178	300000	40.837	nq	100
72) Carbazole	11.72	167	266480	38.744	nq	97
73) Di-n-butylphthalate	12.02	149	328376	40.526	nq	99
74) Fluoranthene	12.70	202	320593	40.414	nq	97
76) Benzidine	12.82	184	143483	34.603	nq	99
77) Pyrene	12.94	202	330036	34.375	nq	100
79) Butylbenzylphthalate	13.53	149	125884	31.890	nq	# 97
80) Benzo(a)anthracene	14.12	228	343899	38.755	nq	99
81) 3,3'-Dichlorobenzidine	14.08	252	127548	40.897	nq	# 95
82) Chrysene	14.17	228	321736	39.411	nq	99
83) Bis(2-ethylhexyl)phthalate	14.08	149	162604	32.220	nq	# 96
84) Di-n-octyl phthalate	14.71	149	327016	39.752	nq	95
85) Indeno(1,2,3-cd)pyrene	17.25	276	224253	32.369	nq	# 90
87) Benzo(b)fluoranthene	15.21	252	310001	42.673	nq	# 95
88) Benzo(k)fluoranthene	15.25	252	282888	40.891	nq	# 95
89) Benzo(a)pyrene	15.61	252	251293	38.912	nq	# 96
90) Dibenzo(a,h)anthracene	17.26	278	194850	35.601	nq	# 95
91) Benzo(g,h,i)perylene	17.73	276	179955	33.639	nq	# 90

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

