

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080819\
 Data File : BF116060.D
 Acq On : 8 Aug 2019 11:13
 Operator : HP/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Aug 08 11:38:24 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 31 15:31:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	157976	20.00	ng	-0.01
21) Naphthalene-d8	8.12	136	649848	20.00	ng	-0.01
39) Acenaphthene-d10	9.87	164	345262	20.00	ng	-0.02
64) Phenanthrene-d10	11.36	188	602622	20.00	ng	-0.02
76) Chrysene-d12	14.00	240	431615	20.00	ng	-0.02
87) Perylene-d12	15.47	264	407659	20.00	ng	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.46	112	743931	78.83	ng	0.00
7) Phenol-d6	6.48	99	954798	80.19	ng	0.00
23) Nitrobenzene-d5	7.41	82	883221	78.45	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	266219	79.53	ng	-0.02
45) 2-Fluorobiphenyl	9.20	172	1446814	78.49	ng	-0.02
79) Terphenyl-d14	12.94	244	1628309	79.49	ng	-0.02

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.60	88	173590	36.413	ng	99
3) Pyridine	3.36	79	491086	36.876	ng	97
4) n-Nitrosodimethylamine	3.31	42	199569	37.974	ng	96
6) Aniline	6.50	93	684068	40.119	ng	94
8) 2-Chlorophenol	6.63	128	423573	40.591	ng	96
9) Benzaldehyde	6.39	77	308216	37.519	ng	97
10) Phenol	6.49	94	554562	39.553	ng	90
11) bis(2-Chloroethyl)ether	6.57	93	419604	40.706	ng	100
12) 1,3-Dichlorobenzene	6.78	146	481280	39.908	ng	98
13) 1,4-Dichlorobenzene	6.86	146	486312	40.274	ng	100
14) 1,2-Dichlorobenzene	7.02	146	450791	40.544	ng	99
15) Benzyl Alcohol	6.99	79	359998	39.843	ng	100
16) 2,2'-oxybis(1-Chloropropan	7.12	45	566263	40.274	ng	93
17) 2-Methylphenol	7.10	107	361713	40.937	ng	98
18) Hexachloroethane	7.35	117	180047	40.499	ng	97
19) n-Nitroso-di-n-propylamine	7.26	70	293473	39.777	ng	97
20) 3+4-Methylphenols	7.25	107	423860	40.537	ng	# 85
22) Acetophenone	7.25	105	566153	39.724	ng	97
24) Nitrobenzene	7.43	77	476719	39.217	ng	96
25) Isophorone	7.66	82	864275	40.148	ng	100
26) 2-Nitrophenol	7.74	139	234253	40.881	ng	99
27) 2,4-Dimethylphenol	7.78	122	342454	39.724	ng	100
28) bis(2-Chloroethoxy)methane	7.87	93	541962	40.618	ng	99
29) 2,4-Dichlorophenol	7.99	162	371600	40.824	ng	97
30) 1,2,4-Trichlorobenzene	8.06	180	420486	40.525	ng	97
31) Naphthalene	8.15	128	1229123	40.193	ng	100
32) Benzoic acid	7.93	122	236815	38.963	ng	98
33) 4-Chloroaniline	8.19	127	560210	41.000	ng	99
34) Hexachlorobutadiene	8.26	225	241005	40.325	ng	99
35) Caprolactam	8.58	113	117747	39.996	ng	94
36) 4-Chloro-3-methylphenol	8.68	107	380175	40.147	ng	98
37) 2-Methylnaphthalene	8.83	142	797811	39.613	ng	100
38) 1-Methylnaphthalene	8.93	142	753399	39.542	ng	100
40) 1,2,4,5-Tetrachlorobenzene	9.00	216	375146	40.643	ng	99

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080819\
 Data File : BF116060.D
 Acq On : 8 Aug 2019 11:13
 Operator : HP/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Aug 08 11:38:24 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 31 15:31:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.99	237	127303	34.370	ng	98
43) 2,4,6-Trichlorophenol	9.12	196	250400	39.413	ng	99
44) 2,4,5-Trichlorophenol	9.16	196	255591	38.918	ng	98
46) 1,1'-Biphenyl	9.30	154	966979	39.670	ng	99
47) 2-Chloronaphthalene	9.33	162	807912	39.823	ng	99
48) 2-Nitroaniline	9.42	65	266246	39.704	ng	91
49) Acenaphthylene	9.74	152	1169488	39.513	ng	100
50) Dimethylphthalate	9.60	163	955577	40.648	ng	100
51) 2,6-Dinitrotoluene	9.66	165	207357	40.588	ng	91
52) Acenaphthene	9.91	154	710304	39.119	ng	99
53) 3-Nitroaniline	9.84	138	252463	39.994	ng	100
54) 2,4-Dinitrophenol	9.95	184	84759	37.247	ng	95
55) Dibenzofuran	10.08	168	1026844	38.887	ng	96
56) 4-Nitrophenol	10.00	139	134495	33.259	ng	98
57) 2,4-Dinitrotoluene	10.07	165	264172	38.838	ng	95
58) Fluorene	10.43	166	805187	39.633	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.20	232	215025	38.917	ng	99
60) Diethylphthalate	10.30	149	887707	40.446	ng	99
61) 4-Chlorophenyl-phenylether	10.41	204	414594	40.294	ng	96
62) 4-Nitroaniline	10.45	138	259077	39.713	ng	99
63) Azobenzene	10.57	77	904959	40.087	ng	97
65) 4,6-Dinitro-2-methylphenol	10.49	198	134537	42.128	ng	96
66) n-Nitrosodiphenylamine	10.53	169	715369	39.440	ng	99
67) 4-Bromophenyl-phenylether	10.90	248	262391	40.674	ng	92
68) Hexachlorobenzene	10.98	284	301362	40.399	ng	97
69) Atrazine	11.06	200	237629	39.823	ng	99
70) Pentachlorophenol	11.17	266	120094	33.160	ng	97
71) Phenanthrene	11.39	178	1199238	38.927	ng	99
72) Anthracene	11.44	178	1226500	39.338	ng	99
73) Carbazole	11.59	167	1135001	40.125	ng	99
74) Di-n-butylphthalate	11.92	149	1368565	41.475	ng	100
75) Fluoranthene	12.57	202	1264414	39.204	ng	98
77) Benzidine	12.69	184	579042	39.710	ng	99
78) Pyrene	12.80	202	1301040	39.988	ng	99
80) Butylbenzylphthalate	13.41	149	585593	42.549	ng	95
81) Benzo(a)anthracene	13.99	228	1138829	41.281	ng	100
82) 3,3'-Dichlorobenzidine	13.95	252	415064	43.306	ng	# 99
83) Chrysene	14.03	228	1086829	40.579	ng	99
84) Bis(2-ethylhexyl)phthalate	13.97	149	798807	44.664	ng	100
85) Di-n-octyl phthalate	14.59	149	1243739	43.782	ng	100
86) Indeno(1,2,3-cd)pyrene	16.95	276	808854	35.957	ng	99
88) Benzo(b)fluoranthene	15.04	252	1011683	42.920	ng	# 97
89) Benzo(k)fluoranthene	15.07	252	880594	38.356	ng	100
90) Benzo(a)pyrene	15.41	252	850210	40.350	ng	99
91) Dibenzo(a,h)anthracene	16.95	278	669784	36.722	ng	99
92) Benzo(g,h,i)perylene	17.39	276	639615	35.708	ng	99

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

