

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012720\
 Data File : BF118714.D
 Acq On : 27 Jan 2020 12:12
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Jan 27 14:29:57 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 27 14:24:10 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	306460	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	1112322	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	621513	20.00	ng	0.00
64) Phenanthrene-d10	11.38	188	1150246	20.00	ng	0.00
76) Chrysene-d12	14.02	240	880108	20.00	ng	0.00
87) Perylene-d12	15.48	264	879539	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.48	112	1372529	83.18	ng	0.00
7) Phenol-d6	6.49	99	1715568	79.99	ng	0.00
23) Nitrobenzene-d5	7.42	82	1542549	77.57	ng	0.00
42) 2,4,6-Tribromophenol	10.69	330	619122	86.13	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	2898633	76.66	ng	0.00
79) Terphenyl-d14	12.96	244	3535694	87.62	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.62	88	308809	38.639	ng	100
3) Pyridine	3.38	79	861091	38.227	ng	100
4) n-Nitrosodimethylamine	3.33	42	314084	32.523	ng	100
6) Aniline	6.52	93	1112462	39.393	ng	100
8) 2-Chlorophenol	6.65	128	763729	41.236	ng	100
9) Benzaldehyde	6.41	77	498376	38.210	ng	100
10) Phenol	6.50	94	947819	38.780	ng	100
11) bis(2-Chloroethyl)ether	6.59	93	713962	39.372	ng	100
12) 1,3-Dichlorobenzene	6.80	146	904412	41.721	ng	100
13) 1,4-Dichlorobenzene	6.88	146	905438	41.609	ng	100
14) 1,2-Dichlorobenzene	7.03	146	834043	40.772	ng	100
15) Benzyl Alcohol	7.00	79	656252	38.586	ng	100
16) 2,2'-oxybis(1-Chloropropan	7.13	45	886746	35.107	ng	100
17) 2-Methylphenol	7.11	107	623615	40.423	ng	100
18) Hexachloroethane	7.37	117	316688	40.135	ng	100
19) n-Nitroso-di-n-propylamine	7.27	70	509692	35.197	ng	100
20) 3+4-Methylphenols	7.26	107	754296	40.046	ng	100
22) Acetophenone	7.27	105	1024641	38.886	ng	100
24) Nitrobenzene	7.44	77	776806	38.149	ng	100
25) Isophorone	7.68	82	1372886	37.850	ng	100
26) 2-Nitrophenol	7.76	139	397585	45.787	ng	100
27) 2,4-Dimethylphenol	7.79	122	564475	37.646	ng	100
28) bis(2-Chloroethoxy)methane	7.89	93	908776	38.884	ng	100
29) 2,4-Dichlorophenol	8.00	162	679187	40.975	ng	100
30) 1,2,4-Trichlorobenzene	8.09	180	784142	40.800	ng	100
31) Naphthalene	8.17	128	2139224	42.574	ng	100
32) Benzoic acid	7.92	122	440322	39.128	ng	100
33) 4-Chloroaniline	8.21	127	949703	40.893	ng	100
34) Hexachlorobutadiene	8.29	225	485725	41.500	ng	100
35) Caprolactam	8.58	113	213073	39.280	ng	100
36) 4-Chloro-3-methylphenol	8.69	107	668576	40.845	ng	100
37) 2-Methylnaphthalene	8.86	142	1474535	41.574	ng	100
38) 1-Methylnaphthalene	8.96	142	1387252	41.170	ng	100
40) 1,2,4,5-Tetrachlorobenzene	9.02	216	797065	40.987	ng	100

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012720\
 Data File : BF118714.D
 Acq On : 27 Jan 2020 12:12
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Jan 27 14:29:57 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 27 14:24:10 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.02	237	326016	33.228	ng	100
43) 2,4,6-Trichlorophenol	9.13	196	526554	40.804	ng	100
44) 2,4,5-Trichlorophenol	9.17	196	535960	40.667	ng	100
46) 1,1'-Biphenyl	9.32	154	1816489	42.407	ng	100
47) 2-Chloronaphthalene	9.35	162	1506046	41.714	ng	100
48) 2-Nitroaniline	9.43	65	423606	38.168	ng	100
49) Acenaphthylene	9.76	152	2162699	42.580	ng	100
50) Dimethylphthalate	9.62	163	1736762	42.891	ng	100
51) 2,6-Dinitrotoluene	9.67	165	396513	43.730	ng	100
52) Acenaphthene	9.93	154	1417755	41.744	ng	100
53) 3-Nitroaniline	9.85	138	442651	42.753	ng	100
54) 2,4-Dinitrophenol	9.95	184	171027	47.246	ng	100
55) Dibenzofuran	10.10	168	2008852	42.681	ng	100
56) 4-Nitrophenol	10.00	139	329338	42.136	ng	100
57) 2,4-Dinitrotoluene	10.08	165	526283	46.041	ng	100
58) Fluorene	10.45	166	1513859	40.971	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.22	232	473277	40.828	ng	100
60) Diethylphthalate	10.32	149	1681629	42.833	ng	100
61) 4-Chlorophenyl-phenylether	10.44	204	821550	40.881	ng	100
62) 4-Nitroaniline	10.46	138	446237	42.017	ng	100
63) Azobenzene	10.60	77	1459580	38.388	ng	100
65) 4,6-Dinitro-2-methylphenol	10.49	198	252929	50.007	ng	100
66) n-Nitrosodiphenylamine	10.56	169	1418978	40.906	ng	100
67) 4-Bromophenyl-phenylether	10.93	248	573227	40.337	ng	100
68) Hexachlorobenzene	11.00	284	663738	41.197	ng	100
69) Atrazine	11.08	200	482299	40.114	ng	100
70) Pentachlorophenol	11.19	266	358893	40.065	ng	100
71) Phenanthrene	11.40	178	2346200	42.279	ng	100
72) Anthracene	11.46	178	2316258	42.183	ng	100
73) Carbazole	11.61	167	2118589	42.064	ng	100
74) Di-n-butylphthalate	11.94	149	2525793	45.138	ng	100
75) Fluoranthene	12.59	202	2519519	43.047	ng	100
77) Benzidine	12.71	184	1113535	47.353	ng	100
78) Pyrene	12.82	202	2550229	42.681	ng	100
80) Butylbenzylphthalate	13.44	149	1082174	44.254	ng	100
81) Benzo(a)anthracene	14.01	228	2206950	41.685	ng	100
82) 3,3'-Dichlorobenzidine	13.97	252	902384	47.833	ng	100
83) Chrysene	14.05	228	2179599	42.956	ng	100
84) Bis(2-ethylhexyl)phthalate	14.00	149	1382057	46.260	ng	100
85) Di-n-octyl phthalate	14.61	149	2232211	50.477	ng	100
86) Indeno(1,2,3-cd)pyrene	16.95	276	2082309	47.230	ng	100
88) Benzo(b)fluoranthene	15.06	252	2183087	39.488	ng	100
89) Benzo(k)fluoranthene	15.09	252	2068883	40.392	ng	100
90) Benzo(a)pyrene	15.42	252	1950533	40.899	ng	100
91) Dibenzo(a,h)anthracene	16.97	278	1712480	40.594	ng	100
92) Benzo(g,h,i)perylene	17.39	276	1642314	40.096	ng	100

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

