

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010320\
 Data File : BF118463.D
 Acq On : 3 Jan 2020 18:01
 Operator : JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Jan 04 00:04:47 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF122419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 24 15:58:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	77575	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	311677	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	160846	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	311256	20.00	ng	0.00
76) Chrysene-d12	14.05	240	231415	20.00	ng	0.00
87) Perylene-d12	15.53	264	179109	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.46	112	336186	73.98	ng	0.00
7) Phenol-d6	6.49	99	437963	77.69	ng	0.00
23) Nitrobenzene-d5	7.42	82	437870	80.35	ng	0.00
42) 2,4,6-Tribromophenol	10.69	330	154212	77.64	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	880380	83.55	ng	0.00
79) Terphenyl-d14	12.98	244	1044051	74.85	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.58	88	71788	39.736	ng	97
3) Pyridine	3.34	79	129321	26.867	ng	# 92
4) n-Nitrosodimethylamine	3.30	42	71684	36.021	ng	# 74
6) Aniline	6.52	93	273986	38.273	ng	96
8) 2-Chlorophenol	6.64	128	203768	39.589	ng	99
9) Benzaldehyde	6.40	77	113634	34.380	ng	98
10) Phenol	6.50	94	245863	38.801	ng	83
11) bis(2-Chloroethyl)ether	6.59	93	173882	38.371	ng	97
12) 1,3-Dichlorobenzene	6.79	146	230919	38.955	ng	99
13) 1,4-Dichlorobenzene	6.87	146	237450	39.437	ng	99
14) 1,2-Dichlorobenzene	7.02	146	223074	39.242	ng	99
15) Benzyl Alcohol	7.00	79	171946	40.058	ng	96
16) 2,2'-oxybis(1-Chloropropan	7.12	45	197167	40.774	ng	69
17) 2-Methylphenol	7.11	107	162813	39.439	ng	94
18) Hexachloroethane	7.36	117	89051	40.299	ng	96
19) n-Nitroso-di-n-propylamine	7.26	70	143442	41.302	ng	92
20) 3+4-Methylphenols	7.27	107	211682	40.064	ng	96
22) Acetophenone	7.26	105	298124	39.027	ng	# 94
24) Nitrobenzene	7.44	77	213240	39.064	ng	99
25) Isophorone	7.67	82	368909	38.844	ng	97
26) 2-Nitrophenol	7.76	139	112909	39.050	ng	96
27) 2,4-Dimethylphenol	7.79	122	150179	38.069	ng	99
28) bis(2-Chloroethoxy)methane	7.88	93	240317	38.692	ng	98
29) 2,4-Dichlorophenol	8.00	162	174130	38.350	ng	97
30) 1,2,4-Trichlorobenzene	8.08	180	202018	38.194	ng	96
31) Naphthalene	8.16	128	651412	41.121	ng	99
32) Benzoic acid	7.94	122	103718	33.129	ng	97
33) 4-Chloroaniline	8.21	127	278199	40.965	ng	99
34) Hexachlorobutadiene	8.27	225	132958	42.229	ng	100
35) Caprolactam	8.59	113	56040	39.532	ng	93
36) 4-Chloro-3-methylphenol	8.70	107	197428	40.498	ng	98
37) 2-Methylnaphthalene	8.85	142	424718	39.159	ng	98
38) 1-Methylnaphthalene	8.95	142	399533	39.099	ng	100
40) 1,2,4,5-Tetrachlorobenzene	9.02	216	199041	41.186	ng	99

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41) Hexachlorocyclopentadiene	9.00	237	53220	32.236	nq	98
43) 2,4,6-Trichlorophenol	9.14	196	120315	37.265	nq	98
44) 2,4,5-Trichlorophenol	9.19	196	122460	36.871	nq	98
46) 1,1'-Biphenyl	9.32	154	520417	40.907	nq	98
47) 2-Chloronaphthalene	9.35	162	410930	40.133	nq	98
48) 2-Nitroaniline	9.45	65	124307	42.982	nq	95
49) Acenaphthylene	9.76	152	631004	41.553	nq	100
50) Dimethylphthalate	9.62	163	491060	40.672	nq	99
51) 2,6-Dinitrotoluene	9.69	165	106393	40.873	nq	93
52) Acenaphthene	9.93	154	377564	40.819	nq	99
53) 3-Nitroaniline	9.86	138	122833	40.205	nq	89
54) 2,4-Dinitrophenol	9.98	184	42583	35.527	nq	98
55) Dibenzofuran	10.10	168	571686	41.829	nq	95
56) 4-Nitrophenol	10.04	139	83540	44.293	nq	98
57) 2,4-Dinitrotoluene	10.10	165	145545	41.845	nq	# 77
58) Fluorene	10.45	166	453656	41.118	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.23	232	108229	38.352	nq	97
60) Diethylphthalate	10.32	149	484299	40.546	nq	98
61) 4-Chlorophenyl-phenylether	10.44	204	227541	41.313	nq	96
62) 4-Nitroaniline	10.48	138	114399	36.044	nq	89
63) Azobenzene	10.60	77	437741	42.785	nq	94
65) 4,6-Dinitro-2-methylphenol	10.52	198	65807	39.192	nq	92
66) n-Nitrosodiphenylamine	10.56	169	405057	40.604	nq	99
67) 4-Bromophenyl-phenylether	10.93	248	140323	38.523	nq	94
68) Hexachlorobenzene	11.00	284	162049	37.794	nq	98
69) Atrazine	11.09	200	115585	37.542	nq	98
70) Pentachlorophenol	11.20	266	63910	34.418	nq	95
71) Phenanthrene	11.42	178	675416	39.805	nq	99
72) Anthracene	11.47	178	682238	40.169	nq	98
73) Carbazole	11.63	167	588393	39.043	nq	98
74) Di-n-butylphthalate	11.95	149	803919	42.070	nq	100
75) Fluoranthene	12.61	202	702132	41.002	nq	98
77) Benzidine	12.73	184	252458	39.814	nq	98
78) Pyrene	12.85	202	692139	36.626	nq	100
80) Butylbenzylphthalate	13.45	149	321424	37.776	nq	99
81) Benzo(a)anthracene	14.04	228	629591	38.868	nq	99
82) 3,3'-Dichlorobenzidine	13.99	252	207378	37.759	nq	98
83) Chrysene	14.07	228	620677	40.017	nq	99
84) Bis(2-ethylhexyl)phthalate	14.01	149	446259	39.886	nq	100
85) Di-n-octyl phthalate	14.62	149	673625	40.930	nq	99
86) Indeno(1,2,3-cd)pyrene	17.05	276	442950	33.517	nq	99
88) Benzo(b)fluoranthene	15.10	252	491075	41.304	nq	99
89) Benzo(k)fluoranthene	15.13	252	444845	38.968	nq	99
90) Benzo(a)pyrene	15.47	252	423167	40.404	nq	99
91) Dibenzo(a,h)anthracene	17.06	278	368608	40.557	nq	99
92) Benzo(g,h,i)perylene	17.50	276	332540	37.876	nq	99

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

