

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012020\
 Data File : BF118590.D
 Acq On : 20 Jan 2020 14:38
 Operator : CG/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ICBF012020

Quant Time: Jan 20 15:38:58 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 20 14:45:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.88	152	442109	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	1613664	20.00	ng	0.00
39) Acenaphthene-d10	9.92	164	925583	20.00	ng	0.00
64) Phenanthrene-d10	11.41	188	1695035	20.00	ng	0.00
76) Chrysene-d12	14.04	240	1396587	20.00	ng	0.00
87) Perylene-d12	15.52	264	1245301	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.50	112	1992190	83.69	ng	0.00
7) Phenol-d6	6.50	99	2596632	83.93	ng	0.00
23) Nitrobenzene-d5	7.45	82	2456278	85.14	ng	0.00
42) 2,4,6-Tribromophenol	10.71	330	918805	85.83	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	4282630	76.05	ng	0.00
79) Terphenyl-d14	12.99	244	5073501	79.23	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.66	88	479254	41.567	ng	98
3) Pyridine	3.42	79	1332835	41.014	ng	99
4) n-Nitrosodimethylamine	3.36	42	575842	41.332	ng	95
6) Aniline	6.55	93	1677072	41.165	ng	99
8) 2-Chlorophenol	6.67	128	1113159	41.662	ng	99
9) Benzaldehyde	6.43	77	798554	42.440	ng	99
10) Phenol	6.52	94	1494488	42.385	ng	99
11) bis(2-Chloroethyl)ether	6.62	93	1094985	41.857	ng	99
12) 1,3-Dichlorobenzene	6.83	146	1306450	41.775	ng	100
13) 1,4-Dichlorobenzene	6.90	146	1327580	42.289	ng	99
14) 1,2-Dichlorobenzene	7.06	146	1266159	42.905	ng	99
15) Benzyl Alcohol	7.02	79	1037930	42.303	ng	100
16) 2,2'-oxybis(1-Chloropropan	7.16	45	1512184	41.500	ng	99
17) 2-Methylphenol	7.13	107	923363	41.488	ng	99
18) Hexachloroethane	7.40	117	479379	42.113	ng	99
19) n-Nitroso-di-n-propylamine	7.30	70	872486	41.763	ng	99
20) 3+4-Methylphenols	7.28	107	1164612	42.860	ng	# 83
22) Acetophenone	7.29	105	1625769	42.530	ng	98
24) Nitrobenzene	7.46	77	1250730	42.341	ng	98
25) Isophorone	7.70	82	2156934	40.991	ng	100
26) 2-Nitrophenol	7.78	139	543997	43.184	ng	99
27) 2,4-Dimethylphenol	7.82	122	872324	40.102	ng	99
28) bis(2-Chloroethoxy)methane	7.92	93	1404938	41.437	ng	99
29) 2,4-Dichlorophenol	8.02	162	1001716	41.657	ng	99
30) 1,2,4-Trichlorobenzene	8.11	180	1159582	41.590	ng	99
31) Naphthalene	8.19	128	3121688	42.825	ng	100
32) Benzoic acid	7.93	122	708397	43.392	ng	99
33) 4-Chloroaniline	8.23	127	1415666	42.018	ng	99
34) Hexachlorobutadiene	8.31	225	708404	41.722	ng	98
35) Caprolactam	8.60	113	312820	39.752	ng	96
36) 4-Chloro-3-methylphenol	8.71	107	993526	41.840	ng	99
37) 2-Methylnaphthalene	8.88	142	2192713	42.616	ng	99
38) 1-Methylnaphthalene	8.98	142	2093570	42.828	ng	100
40) 1,2,4,5-Tetrachlorobenzene	9.05	216	1215229	41.961	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.04	237	592805	40.571	nq	99
43) 2,4,6-Trichlorophenol	9.15	196	786814	40.942	nq	99
44) 2,4,5-Trichlorophenol	9.19	196	803318	40.929	nq	98
46) 1,1'-Biphenyl	9.35	154	2693956	42.231	nq	100
47) 2-Chloronaphthalene	9.37	162	2261989	42.070	nq	99
48) 2-Nitroaniline	9.46	65	701598	42.449	nq	98
49) Acenaphthylene	9.79	152	3192007	42.200	nq	100
50) Dimethylphthalate	9.65	163	2528459	41.929	nq	100
51) 2,6-Dinitrotoluene	9.70	165	576800	42.715	nq	97
52) Acenaphthene	9.96	154	2148165	42.471	nq	100
53) 3-Nitroaniline	9.87	138	657125	42.617	nq	99
54) 2,4-Dinitrophenol	9.97	184	234095	43.424	nq	# 78
55) Dibenzofuran	10.13	168	2968707	42.354	nq	100
56) 4-Nitrophenol	10.02	139	497310	42.724	nq	97
57) 2,4-Dinitrotoluene	10.10	165	743118	43.654	nq	99
58) Fluorene	10.47	166	2359027	42.871	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.24	232	718298	41.608	nq	99
60) Diethylphthalate	10.34	149	2501964	42.792	nq	99
61) 4-Chlorophenyl-phenylether	10.46	204	1272552	42.521	nq	99
62) 4-Nitroaniline	10.48	138	669236	42.313	nq	98
63) Azobenzene	10.62	77	2403669	42.449	nq	98
65) 4,6-Dinitro-2-methylphenol	10.51	198	324874	43.587	nq	94
66) n-Nitrosodiphenylamine	10.58	169	2153567	42.129	nq	100
67) 4-Bromophenyl-phenylether	10.95	248	873347	41.704	nq	# 93
68) Hexachlorobenzene	11.02	284	987944	41.611	nq	98
69) Atrazine	11.10	200	743255	41.950	nq	99
70) Pentachlorophenol	11.21	266	560008	42.424	nq	98
71) Phenanthrene	11.43	178	3472928	42.469	nq	99
72) Anthracene	11.49	178	3445978	42.587	nq	100
73) Carbazole	11.63	167	3155383	42.514	nq	99
74) Di-n-butylphthalate	11.97	149	3547160	43.017	nq	98
75) Fluoranthene	12.62	202	3708302	42.995	nq	99
77) Benzidine	12.73	184	1490747	39.950	nq	# 95
78) Pyrene	12.85	202	3756056	39.614	nq	100
80) Butylbenzylphthalate	13.47	149	1599840	41.228	nq	97
81) Benzo(a)anthracene	14.03	228	3488676	41.526	nq	99
82) 3,3'-Dichlorobenzidine	14.00	252	1280779	42.784	nq	98
83) Chrysene	14.07	228	3319081	41.222	nq	98
84) Bis(2-ethylhexyl)phthalate	14.03	149	2035374	42.933	nq	99
85) Di-n-octyl phthalate	14.64	149	3230152	46.031	nq	100
86) Indeno(1,2,3-cd)pyrene	16.99	276	2774334	39.655	nq	100
88) Benzo(b)fluoranthene	15.09	252	3405252	43.503	nq	98
89) Benzo(k)fluoranthene	15.12	252	2883753	39.765	nq	99
90) Benzo(a)pyrene	15.46	252	2805022	41.541	nq	100
91) Dibenzo(a,h)anthracene	17.01	278	2301289	38.529	nq	99
92) Benzo(g,h,i)perylene	17.44	276	2142384	36.942	nq	97

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

