

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF051420\
 Data File : BF120282.D
 Acq On : 14 May 2020 11:07
 Operator : MA/SJ
 Sample : SSTDICC060
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC060

Quant Time: May 14 11:43:35 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF051420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 14 10:46:42 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.63	152	298432	20.00	ng	0.00
21) Naphthalene-d8	7.91	136	1319983	20.00	ng	0.00
39) Acenaphthene-d10	9.66	164	722685	20.00	ng	0.00
64) Phenanthrene-d10	11.15	188	1331078	20.00	ng	0.00
76) Chrysene-d12	13.79	240	740892	20.00	ng	0.00
87) Perylene-d12	15.17	264	676425	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.23	112	1889543	95.69	ng	0.00
7) Phenol-d6	6.29	99	2584930	109.17	ng	0.00
23) Nitrobenzene-d5	7.20	82	2227429	91.50	ng	0.00
42) 2,4,6-Tribromophenol	10.46	330	701936	93.06	ng	0.00
45) 2-Fluorobiphenyl	8.99	172	3800171	83.67	ng	0.00
79) Terphenyl-d14	12.74	244	3874762	99.89	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.20	88	455447	50.742	ng	99
3) Pyridine	2.86	79	1239954	50.079	ng	100
4) n-Nitrosodimethylamine	2.84	42	496443	49.581	ng	100
6) Aniline	6.29	93	1676847	57.411	ng	96
8) 2-Chlorophenol	6.42	128	1114188	54.735	ng	97
9) Benzaldehyde	6.17	77	726257	61.114	ng	97
10) Phenol	6.30	94	1531801	58.204	ng	94
11) bis(2-Chloroethyl)ether	6.37	93	1132262	53.086	ng	99
12) 1,3-Dichlorobenzene	6.56	146	1186508	55.414	ng	100
13) 1,4-Dichlorobenzene	6.64	146	1157186	51.222	ng	99
14) 1,2-Dichlorobenzene	6.80	146	1116577	55.967	ng	99
15) Benzyl Alcohol	6.79	79	867489	55.565	ng	99
16) 2,2'-oxybis(1-Chloropropan	6.92	45	1705166	53.801	ng	96
17) 2-Methylphenol	6.90	107	917847	53.435	ng	95
18) Hexachloroethane	7.13	117	453423	56.813	ng	99
19) n-Nitroso-di-n-propylamine	7.06	70	778216	54.676	ng	99
20) 3+4-Methylphenols	7.06	107	1201978	54.452	ng	95
22) Acetophenone	7.05	105	1565868	49.002	ng	97
24) Nitrobenzene	7.22	77	1216938	48.991	ng	98
25) Isophorone	7.46	82	2349582	48.301	ng	99
26) 2-Nitrophenol	7.53	139	639461	47.914	ng	94
27) 2,4-Dimethylphenol	7.58	122	918470	47.321	ng	97
28) bis(2-Chloroethoxy)methane	7.67	93	1425314	48.121	ng	100
29) 2,4-Dichlorophenol	7.78	162	883598	45.581	ng	98
30) 1,2,4-Trichlorobenzene	7.86	180	979461	44.866	ng	99
31) Naphthalene	7.93	128	3183529	47.856	ng	98
32) Benzoic acid	7.75	122	662361	51.350	ng	98
33) 4-Chloroaniline	7.99	127	1483110	50.684	ng	98
34) Hexachlorobutadiene	8.05	225	603866	48.094	ng	97
35) Caprolactam	8.39	113	305041	51.699	ng	97
36) 4-Chloro-3-methylphenol	8.49	107	984946	47.170	ng	99
37) 2-Methylnaphthalene	8.63	142	2092829	46.038	ng	98
38) 1-Methylnaphthalene	8.72	142	2008872	45.342	ng	99
40) 1,2,4,5-Tetrachlorobenzene	8.79	216	948441	43.488	ng	100

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41) Hexachlorocyclopentadiene	8.77	237	390136	50.156	nq	99
43) 2,4,6-Trichlorophenol	8.91	196	645106	46.263	nq	99
44) 2,4,5-Trichlorophenol	8.96	196	742337	46.451	nq	97
46) 1,1'-Biphenyl	9.09	154	2437886	42.470	nq	98
47) 2-Chloronaphthalene	9.12	162	1975672	45.080	nq	99
48) 2-Nitroaniline	9.22	65	804577	51.589	nq	99
49) Acenaphthylene	9.53	152	3215821	46.968	nq	99
50) Dimethylphthalate	9.40	163	2619172	48.311	nq	99
51) 2,6-Dinitrotoluene	9.47	165	592970	49.906	nq	# 86
52) Acenaphthene	9.70	154	1980626	47.764	nq	99
53) 3-Nitroaniline	9.64	138	718715	50.883	nq	96
54) 2,4-Dinitrophenol	9.75	184	302063	50.910	nq	95
55) Dibenzofuran	9.87	168	2801177	46.422	nq	99
56) 4-Nitrophenol	9.82	139	404229	52.767	nq	98
57) 2,4-Dinitrotoluene	9.87	165	676586	46.651	nq	98
58) Fluorene	10.22	166	2139801	45.191	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.00	232	637020	50.652	nq	99
60) Diethylphthalate	10.10	149	2562771	47.395	nq	99
61) 4-Chlorophenyl-phenylether	10.21	204	1062253	44.400	nq	95
62) 4-Nitroaniline	10.26	138	694314	49.161	nq	100
63) Azobenzene	10.37	77	2485256	49.058	nq	97
65) 4,6-Dinitro-2-methylphenol	10.28	198	407302	45.700	nq	95
66) n-Nitrosodiphenylamine	10.33	169	2090458	43.697	nq	100
67) 4-Bromophenyl-phenylether	10.70	248	730067	43.713	nq	98
68) Hexachlorobenzene	10.76	284	758478	43.729	nq	97
69) Atrazine	10.87	200	620887	61.170	nq	99
70) Pentachlorophenol	10.96	266	359062	50.436	nq	98
71) Phenanthrene	11.17	178	3104789	46.738	nq	98
72) Anthracene	11.23	178	3139163	44.073	nq	99
73) Carbazole	11.39	167	2804370	39.316	nq	99
74) Di-n-butylphthalate	11.72	149	3503460	38.912	nq	99
75) Fluoranthene	12.36	202	3096009	39.533	nq	98
77) Benzidine	12.49	184	1206382	67.548	nq	96
78) Pyrene	12.59	202	3150987	51.356	nq	98
80) Butylbenzylphthalate	13.21	149	1497284	51.356	nq	99
81) Benzo(a)anthracene	13.78	228	2269857	48.438	nq	100
82) 3,3'-Dichlorobenzidine	13.74	252	841403	49.549	nq	98
83) Chrysene	13.82	228	2411832	47.319	nq	99
84) Bis(2-ethylhexyl)phthalate	13.77	149	1811980	48.934	nq	99
85) Di-n-octyl phthalate	14.39	149	3334046	48.659	nq	99
86) Indeno(1,2,3-cd)pyrene	16.52	276	2102327	44.870	nq	99
88) Benzo(b)fluoranthene	14.79	252	2199435	49.176	nq	99
89) Benzo(k)fluoranthene	14.82	252	2089638	52.591	nq	99
90) Benzo(a)pyrene	15.13	252	1965614	49.577	nq	99
91) Dibenzo(a,h)anthracene	16.53	278	1782100	49.915	nq	99
92) Benzo(g,h,i)perylene	16.92	276	1802087	50.698	nq	98

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

