

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030620\
 Data File : BF119275.D
 Acq On : 7 Mar 2020 4:00
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Mar 09 04:06:51 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 06 23:38:46 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	105409	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	382007	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	199469	20.00	ng	0.00
64) Phenanthrene-d10	11.25	188	340179	20.00	ng	0.00
76) Chrysene-d12	13.89	240	278561	20.00	ng	0.00
87) Perylene-d12	15.31	264	257539	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.35	112	535798	84.95	ng	0.00
7) Phenol-d6	6.38	99	626822	81.71	ng	0.00
23) Nitrobenzene-d5	7.29	82	585332	80.90	ng	0.00
42) 2,4,6-Tribromophenol	10.55	330	193165	74.03	ng	0.00
45) 2-Fluorobiphenyl	9.08	172	1103611	81.51	ng	0.00
79) Terphenyl-d14	12.83	244	1229361	75.98	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.34	88	105207	37.463	ng	99
3) Pyridine	3.07	79	291523	38.011	ng	98
4) n-Nitrosodimethylamine	3.02	42	94273	40.577	ng	92
6) Aniline	6.39	93	368076	38.886	ng	# 77
8) 2-Chlorophenol	6.52	128	279145	41.009	ng	98
9) Benzaldehyde	6.27	77	181721	35.457	ng	98
10) Phenol	6.40	94	339897	40.982	ng	95
11) bis(2-Chloroethyl)ether	6.46	93	248743	39.197	ng	98
12) 1,3-Dichlorobenzene	6.67	146	329741	40.417	ng	100
13) 1,4-Dichlorobenzene	6.75	146	335356	41.077	ng	97
14) 1,2-Dichlorobenzene	6.90	146	310281	41.672	ng	98
15) Benzyl Alcohol	6.88	79	228350	41.188	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.00	45	209543	38.118	ng	# 48
17) 2-Methylphenol	7.00	107	221232	40.087	ng	94
18) Hexachloroethane	7.23	117	95675	32.756	ng	96
19) n-Nitroso-di-n-propylamine	7.15	70	186934	41.820	ng	96
20) 3+4-Methylphenols	7.15	107	292620	43.279	ng	90
22) Acetophenone	7.14	105	368406	41.681	ng	# 94
24) Nitrobenzene	7.31	77	287087	40.126	ng	95
25) Isophorone	7.55	82	488041	38.841	ng	100
26) 2-Nitrophenol	7.63	139	134527	35.487	ng	97
27) 2,4-Dimethylphenol	7.67	122	199632	38.802	ng	98
28) bis(2-Chloroethoxy)methane	7.76	93	315776	38.610	ng	99
29) 2,4-Dichlorophenol	7.88	162	236742	39.087	ng	98
30) 1,2,4-Trichlorobenzene	7.95	180	277067	38.582	ng	99
31) Naphthalene	8.03	128	784183	39.582	ng	98
32) Benzoic acid	7.83	122	106815	31.173	ng	99
33) 4-Chloroaniline	8.09	127	335621	39.387	ng	97
34) Hexachlorobutadiene	8.15	225	172701	38.608	ng	98
35) Caprolactam	8.46	113	75088	40.262	ng	92
36) 4-Chloro-3-methylphenol	8.58	107	247316	40.347	ng	99
37) 2-Methylnaphthalene	8.72	142	523511	39.266	ng	99
38) 1-Methylnaphthalene	8.82	142	492671	39.292	ng	98
40) 1,2,4,5-Tetrachlorobenzene	8.89	216	269788	40.116	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.87	237	1327	9.873	ng	75
43) 2,4,6-Trichlorophenol	9.01	196	171567	40.398	ng	98
44) 2,4,5-Trichlorophenol	9.06	196	172582	38.725	ng	98
46) 1,1'-Biphenyl	9.19	154	662791	41.113	ng	96
47) 2-Chloronaphthalene	9.21	162	527215	40.582	ng	98
48) 2-Nitroaniline	9.32	65	161278	44.509	ng	98
49) Acenaphthylene	9.62	152	752323	40.096	ng	98
50) Dimethylphthalate	9.49	163	623294	40.864	ng	99
51) 2,6-Dinitrotoluene	9.55	165	128153	37.817	ng #	76
52) Acenaphthene	9.79	154	455015	40.217	ng	99
53) 3-Nitroaniline	9.73	138	158890	42.293	ng	97
54) 2,4-Dinitrophenol	9.87	184	5109	10.287	ng	92
55) Dibenzofuran	9.96	168	670470	39.703	ng	100
56) 4-Nitrophenol	9.92	139	80685	35.746	ng	98
57) 2,4-Dinitrotoluene	9.96	165	159528	36.318	ng #	86
58) Fluorene	10.31	166	535347	40.797	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.10	232	139662	37.140	ng	99
60) Diethylphthalate	10.18	149	592902	41.248	ng	99
61) 4-Chlorophenyl-phenylether	10.30	204	283519	40.819	ng	96
62) 4-Nitroaniline	10.34	138	149745	38.959	ng	95
63) Azobenzene	10.46	77	500166	40.032	ng	95
65) 4,6-Dinitro-2-methylphenol	10.39	198	14679	7.131	ng	97
66) n-Nitrosodiphenylamine	10.42	169	476744	42.801	ng	99
67) 4-Bromophenyl-phenylether	10.79	248	180277	40.543	ng	96
68) Hexachlorobenzene	10.86	284	204130	39.817	ng	99
69) Atrazine	10.95	200	133082	34.196	ng	99
70) Pentachlorophenol	11.07	266	143974	69.716	ng	96
71) Phenanthrene	11.27	178	758505	40.886	ng	97
72) Anthracene	11.32	178	758479	40.919	ng	98
73) Carbazole	11.48	167	683568	41.395	ng	98
74) Di-n-butylphthalate	11.80	149	863859	43.197	ng	98
75) Fluoranthene	12.46	202	791661	40.202	ng	99
77) Benzidine	12.58	184	569005	50.226	ng	99
78) Pyrene	12.69	202	795982	35.586	ng	98
80) Butylbenzylphthalate	13.30	149	353724	37.574	ng	100
81) Benzo(a)anthracene	13.87	228	767589	40.442	ng	99
82) 3,3'-Dichlorobenzidine	13.83	252	316248	42.910	ng	99
83) Chrysene	13.92	228	741928	39.230	ng	98
84) Bis(2-ethylhexyl)phthalate	13.86	149	481552	40.489	ng	100
85) Di-n-octyl phthalate	14.46	149	805269	42.495	ng	99
86) Indeno(1,2,3-cd)pyrene	16.72	276	559392	30.548	ng	99
88) Benzo(b)fluoranthene	14.91	252	735913	43.351	ng	98
89) Benzo(k)fluoranthene	14.93	252	659499	41.922	ng	99
90) Benzo(a)pyrene	15.26	252	608764	39.734	ng	99
91) Dibenzo(a,h)anthracene	16.72	278	472124	35.022	ng	99
92) Benzo(g,h,i)perylene	17.14	276	420427	31.565	ng	99

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

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