

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF073118\
 Data File : BF107876.D
 Acq On : 1 Aug 2018 3:41
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil

8/13/2018 12:44:30 PM

Quant Time: Aug 13 12:27:14 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF073018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 30 17:48:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	134379	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	518170	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	197062	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	375276	20.00	ng	0.00
75) Chrysene-d12	14.15	240	343789	20.00	ng	0.00
86) Perylene-d12	15.69	264	255927	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.60	112	568829	71.94	ng	0.00
7) Phenol-d6	6.61	99	768692	73.77	ng	0.00
23) Nitrobenzene-d5	7.53	82	730696	81.24	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	175274	65.46	ng	0.00
44) 2-Fluorobiphenyl	9.32	172	1282307	90.68	ng	0.00
78) Terphenyl-d14	13.08	244	1228420	77.48	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.79	88	143933	39.910	ng	# 84
3) Pyridine	3.60	79	322972	36.517	ng	# 84
4) n-Nitrosodimethylamine	3.57	42	66821	16.610	ng	93
6) Aniline	6.64	93	427793	39.469	ng	# 73
8) 2-Chlorophenol	6.75	128	369218	37.175	ng	95
9) Benzaldehyde	6.52	77	196133	31.196	ng	96
10) Phenol	6.62	94	395817	31.852	ng	93
11) bis(2-Chloroethyl)ether	6.70	93	431723	39.070	ng	94
12) 1,3-Dichlorobenzene	6.91	146	395003	37.920	ng	97
13) 1,4-Dichlorobenzene	6.98	146	446202	38.480	ng	98
14) 1,2-Dichlorobenzene	7.14	146	395837	36.774	ng	99
15) Benzyl Alcohol	7.11	79	232753	35.936	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.23	45	588859	37.065	ng	74
17) 2-Methylphenol	7.22	107	275442	37.568	ng	# 78
18) Hexachloroethane	7.47	117	118353	28.396	ng	96
19) n-Nitroso-di-n-propylamine	7.37	70	248538	37.598	ng	100
20) 3+4-Methylphenols	7.38	107	328214	36.462	ng	# 58
22) Acetophenone	7.38	105	491394	39.417	ng	# 91
24) Nitrobenzene	7.55	77	374751	39.735	ng	99
25) Isophorone	7.78	82	597040	40.438	ng	99
26) 2-Nitrophenol	7.87	139	158226	33.429	ng	98
27) 2,4-Dimethylphenol	7.90	122	234851	37.060	ng	100
28) bis(2-Chloroethoxy)methane	7.99	93	395827	40.682	ng	99
29) 2,4-Dichlorophenol	8.11	162	274146	38.246	ng	98
30) 1,2,4-Trichlorobenzene	8.18	180	312956	38.271	ng	98
31) Naphthalene	8.27	128	955975	39.689	ng	99
32) Benzoic acid	8.01	122	97697m	23.577	ng	
33) 4-Chloroaniline	8.32	127	364813	35.459	ng	99
34) Hexachlorobutadiene	8.37	225	192970	38.187	ng	98
35) Caprolactam	8.68	113	74629	34.585	ng	# 79
36) 4-Chloro-3-methylphenol	8.81	107	269214	34.582	ng	97
37) 2-Methylnaphthalene	8.96	142	558920	36.941	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.12	216	286084	42.579	ng	# 97
40) Hexachlorocyclopentadiene	9.10	237	1461	7.873	ng	84

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.24	196	151315	40.738	ng	100
43) 2,4,5-Trichlorophenol	9.30	196	161614	36.142	ng	95
45) 1,1'-Biphenyl	9.42	154	768559	42.855	ng	99
46) 2-Chloronaphthalene	9.45	162	547088	40.454	ng	96
47) 2-Nitroaniline	9.56	65	180670	41.071	ng	95
48) Acenaphthylene	9.87	152	797861	38.055	ng	100
49) Dimethylphthalate	9.71	163	610509	41.252	ng	99
50) 2,6-Dinitrotoluene	9.79	165	113474	35.282	ng #	91
51) Acenaphthene	10.04	154	470843	38.660	ng	99
52) 3-Nitroaniline	9.98	138	160654	39.640	ng	96
54) Dibenzofuran	10.21	168	697217	37.038	ng	100
55) 4-Nitrophenol	10.18	139	27208m	21.916	ng	
56) 2,4-Dinitrotoluene	10.20	165	131183	31.033	ng #	91
57) Fluorene	10.55	166	519302	35.497	ng	93
58) 2,3,4,6-Tetrachlorophenol	10.34	232	141468	34.444	ng #	77
59) Diethylphthalate	10.41	149	601543	37.804	ng	98
60) 4-Chlorophenyl-phenylether	10.54	204	285162	35.781	ng	91
61) 4-Nitroaniline	10.60	138	153927	42.798	ng	91
62) Azobenzene	10.70	77	523358	35.316	ng	95
64) 4,6-Dinitro-2-methylphenol	10.62	198	2361	1.156	ng #	1
65) n-Nitrosodiphenylamine	10.67	169	485441	39.285	ng	100
66) 4-Bromophenyl-phenylether	11.04	248	163912	37.474	ng #	85
67) Hexachlorobenzene	11.11	284	180923	37.301	ng #	81
68) Atrazine	11.18	200	133225	36.428	ng	98
69) Pentachlorophenol	11.31	266	82865	34.705	ng	94
70) Phenanthrene	11.52	178	688425	38.634	ng	98
71) Anthracene	11.58	178	825332	40.156	ng	99
72) Carbazole	11.74	167	823103	40.525	ng	98
73) Di-n-butylphthalate	12.04	149	892867	40.294	ng	99
74) Fluoranthene	12.72	202	890081	42.518	ng	97
76) Benzidine	12.85	184	507213	45.631	ng	97
77) Pyrene	12.95	202	911257	36.810	ng	97
79) Butylbenzylphthalate	13.55	149	419178	38.667	ng	92
80) Benzo(a)anthracene	14.14	228	726744	37.318	ng	99
81) 3,3'-Dichlorobenzidine	14.10	252	323369	42.835	ng	99
82) Chrysene	14.18	228	790364	37.627	ng	98
83) Bis(2-ethylhexyl)phthalate	14.10	149	557708	40.922	ng	99
84) Di-n-octyl phthalate	14.72	149	970367	44.359	ng	97
85) Indeno(1,2,3-cd)pyrene	17.28	276	524225	32.811	ng #	93
87) Benzo(b)fluoranthene	15.23	252	471746	37.374	ng	98
88) Benzo(k)fluoranthene	15.26	252	648352m	41.837	ng	
89) Benzo(a)pyrene	15.62	252	477255	36.626	ng	99
90) Dibenzo(a,h)anthracene	17.29	278	439813	38.694	ng	94
91) Benzo(g,h,i)perylene	17.76	276	416621	37.881	ng #	94

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

