

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120518\
 Data File : BF111110.D
 Acq On : 5 Dec 2018 15:45
 Operator : JU/SJ
 Sample : SSTDICC060
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC060

Manual Integrations
 APPROVED

mohammad
 12/6/2018 12:11:38 PM

Quant Time: Dec 05 17:28:14 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 05 15:25:21 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	564706	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	2310417	20.00	ng	0.00
39) Acenaphthene-d10	9.85	164	1071203	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	1775368	20.00	ng	0.00
76) Chrysene-d12	13.96	240	984050	20.00	ng	0.00
87) Perylene-d12	15.39	264	923119	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.42	112	3893438	114.33	ng	0.00
7) Phenol-d6	6.44	99	5080282	114.48	ng	0.00
23) Nitrobenzene-d5	7.37	82	4588323	113.66	ng	0.00
42) 2,4,6-Tribromophenol	10.63	330	960979	90.33	ng	0.00
45) 2-Fluorobiphenyl	9.17	172	6448318	87.21	ng	0.00
79) Terphenyl-d14	12.91	244	5430135	107.42	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.55	88	1208054	74.048	ng	# 82
3) Pyridine	3.28	79	2726561	76.322	ng	# 73
4) n-Nitrosodimethylamine	3.22	42	1365968	83.561	ng	# 85
6) Aniline	6.47	93	3493372	63.679	ng	# 54
8) 2-Chlorophenol	6.59	128	2289557	56.733	ng	91
9) Benzaldehyde	6.35	77	1313987	55.472	ng	98
10) Phenol	6.45	94	2845525m	56.937	ng	
11) bis(2-Chloroethyl)ether	6.55	93	2340594	62.165	ng	88
12) 1,3-Dichlorobenzene	6.75	146	2438033	54.991	ng	97
13) 1,4-Dichlorobenzene	6.83	146	2459055	53.936	ng	95
14) 1,2-Dichlorobenzene	6.98	146	2246192	52.346	ng	97
15) Benzyl Alcohol	6.95	79	2073802	60.809	ng	93
16) 2,2'-oxybis(1-Chloropropan	7.09	45	4498494	75.075	ng	66
17) 2-Methylphenol	7.06	107	1966335	61.280	ng	# 83
18) Hexachloroethane	7.32	117	955265	56.960	ng	# 88
19) n-Nitroso-di-n-propylamine	7.23	70	1789023	61.889	ng	# 86
20) 3+4-Methylphenols	7.22	107	2240064	52.517	ng	# 85
22) Acetophenone	7.22	105	3060014	56.504	ng	# 83
24) Nitrobenzene	7.39	77	2429135	57.862	ng	93
25) Isophorone	7.63	82	4115793	62.152	ng	96
26) 2-Nitrophenol	7.70	139	1169019	57.316	ng	97
27) 2,4-Dimethylphenol	7.75	122	1888424	61.309	ng	96
28) bis(2-Chloroethoxy)methane	7.84	93	2741659	61.799	ng	98
29) 2,4-Dichlorophenol	7.95	162	1908277	59.205	ng	98
30) 1,2,4-Trichlorobenzene	8.03	180	1882784	52.123	ng	95
31) Naphthalene	8.11	128	5378932	49.666	ng	# 90
32) Benzoic acid	7.88	122	1333243m	58.219	ng	
33) 4-Chloroaniline	8.16	127	2630361	57.062	ng	95
34) Hexachlorobutadiene	8.24	225	1020768	49.615	ng	100
35) Caprolactam	8.55	113	706677	64.220	ng	# 67
36) 4-Chloro-3-methylphenol	8.64	107	2058649	57.484	ng	99
37) 2-Methylnaphthalene	8.80	142	3709908	51.862	ng	97
38) 1-Methylnaphthalene	8.90	142	3601868	51.765	ng	# 100
40) 1,2,4,5-Tetrachlorobenzene	8.97	216	1567964	46.531	ng	# 96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.96	237	894270	263.738	nq	99
43) 2,4,6-Trichlorophenol	9.08	196	1183832	58.822	nq	99
44) 2,4,5-Trichlorophenol	9.12	196	1226239	57.865	nq	95
46) 1,1'-Biphenyl	9.27	154	4524431	45.813	nq	87
47) 2-Chloronaphthalene	9.29	162	3674196	51.311	nq	92
48) 2-Nitroaniline	9.38	65	1340789	59.864	nq	99
49) Acenaphthylene	9.71	152	5181812	50.891	nq #	90
50) Dimethylphthalate	9.57	163	3984074	50.712	nq	97
51) 2,6-Dinitrotoluene	9.63	165	987714	56.474	nq	97
52) Acenaphthene	9.88	154	3399912	52.086	nq	96
53) 3-Nitroaniline	9.79	138	1230681	60.374	nq	97
54) 2,4-Dinitrophenol	9.89	184	291894	52.503	nq	91
55) Dibenzofuran	10.05	168	4681314	49.159	nq	91
56) 4-Nitrophenol	9.95	139	921149	90.607	nq	93
57) 2,4-Dinitrotoluene	10.03	165	1260577	55.553	nq	96
58) Fluorene	10.39	166	3333530	44.993	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.16	232	890835m	51.223	nq	
60) Diethylphthalate	10.27	149	4058919	50.101	nq	96
61) 4-Chlorophenyl-phenylether	10.39	204	1660193	42.841	nq	94
62) 4-Nitroaniline	10.41	138	1121384	59.084	nq	91
63) Azobenzene	10.55	77	4267909	54.048	nq	88
65) 4,6-Dinitro-2-methylphenol	10.43	198	485643	57.828	nq #	56
66) n-Nitrosodiphenylamine	10.50	169	3432219	58.073	nq	90
67) 4-Bromophenyl-phenylether	10.87	248	1068477	55.261	nq #	90
68) Hexachlorobenzene	10.94	284	1099411	53.383	nq #	82
69) Atrazine	11.03	200	582012	38.266	nq	98
70) Pentachlorophenol	11.13	266	734398	89.630	nq	100
71) Phenanthrene	11.35	178	4629231m	49.232	nq	
72) Anthracene	11.40	178	4643070	48.487	nq #	92
73) Carbazole	11.55	167	4880153	52.506	nq #	90
74) Di-n-butylphthalate	11.89	149	5354957	49.277	nq #	93
75) Fluoranthene	12.53	202	4528777	46.862	nq	91
77) Benzidine	12.65	184	1176309	46.734	nq	100
78) Pyrene	12.76	202	4585492	63.800	nq	92
80) Butylbenzylphthalate	13.39	149	2352227	72.490	nq	96
81) Benzo(a)anthracene	13.94	228	3262369	56.154	nq	96
82) 3,3'-Dichlorobenzidine	13.91	252	1193306	59.903	nq	100
83) Chrysene	13.99	228	3213422	56.187	nq	96
84) Bis(2-ethylhexyl)phthalate	13.95	149	2658726	61.491	nq	93
85) Di-n-octyl phthalate	14.56	149	4098268	60.046	nq #	83
86) Indeno(1,2,3-cd)pyrene	16.80	276	2963846m	72.197	nq	
88) Benzo(b)fluoranthene	14.97	252	3084317	56.875	nq #	95
89) Benzo(k)fluoranthene	15.00	252	2579873	46.219	nq #	93
90) Benzo(a)pyrene	15.33	252	2746697	54.413	nq #	96
91) Dibenzo(a,h)anthracene	16.82	278	2465608	59.200	nq #	94
92) Benzo(g,h,i)perylene	17.23	276	2505943	62.403	nq #	92

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

