

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110918\
 Data File : BN003477.D
 Acq On : 09 Nov 2018 18:11
 Operator : MJ/SJ
 Sample : J5860-05MS
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 EP2-SB-104(15-17)MS

Manual Integrations
 APPROVED

Sohil
 11/12/2018 5:07:48 PM

Quant Time: Nov 12 03:49:23 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-BN102418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 07 08:25:32 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	104285	20.00	ng	0.00
21) Naphthalene-d8	10.65	136	448200	20.00	ng	0.00
39) Acenaphthene-d10	14.49	164	251640	20.00	ng	0.00
64) Phenanthrene-d10	17.23	188	513544	20.00	ng	0.00
76) Chrysene-d12	21.41	240	431252	20.00	ng	0.00
87) Perylene-d12	23.72	264	430231	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.42	112	691303	115.46	ng	0.00
7) Phenol-d6	7.01	99	917619	107.78	ng	0.00
23) Nitrobenzene-d5	9.02	82	522909	66.91	ng	0.00
42) 2,4,6-Tribromophenol	15.97	330	401633	118.53	ng	0.00
45) 2-Fluorobiphenyl	13.10	172	1382351	74.51	ng	0.00
79) Terphenyl-d14	19.85	244	1637680	81.21	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.33	88	64227	22.593	ng	92
3) Pyridine	3.73	79	187170	28.812	ng	92
4) n-Nitrosodimethylamine	3.65	42	74389	31.102	ng	91
6) Aniline	7.18	93	263346	24.771	ng	97
8) 2-Chlorophenol	7.41	128	259930	35.247	ng	95
9) Benzaldehyde	7.00	77	126035	25.490	ng	96
10) Phenol	7.03	94	373417	41.563	ng	95
11) bis(2-Chloroethyl)ether	7.28	93	222155	30.273	ng	93
12) 1,3-Dichlorobenzene	7.73	146	262260	31.986	ng	97
13) 1,4-Dichlorobenzene	7.89	146	268693	31.826	ng	98
14) 1,2-Dichlorobenzene	8.20	146	260192	31.755	ng	98
15) Benzyl Alcohol	8.09	79	200782	32.726	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.38	45	288230	26.790	ng	97
17) 2-Methylphenol	8.28	107	213115	34.708	ng	98
18) Hexachloroethane	8.92	117	90915	30.739	ng	94
19) n-Nitroso-di-n-propylamine	8.66	70	169601	27.625	ng	95
20) 3+4-Methylphenols	8.61	107	286935	32.897	ng	97
22) Acetophenone	8.68	105	357742	32.929	ng	# 97
24) Nitrobenzene	9.06	77	249533	31.289	ng	93
25) Isophorone	9.58	82	473750	34.558	ng	96
26) 2-Nitrophenol	9.76	139	138113	32.903	ng	93
27) 2,4-Dimethylphenol	9.81	122	238396	40.393	ng	99
28) bis(2-Chloroethoxy)methane	10.06	93	309289	33.489	ng	98
29) 2,4-Dichlorophenol	10.29	162	251162	37.292	ng	99
30) 1,2,4-Trichlorobenzene	10.50	180	260705	34.050	ng	99
31) Naphthalene	10.70	128	787345	35.948	ng	99
32) Benzoic acid	9.88	122	29038m	10.157	ng	
33) 4-Chloroaniline	10.81	127	141785	14.650	ng	96
34) Hexachlorobutadiene	10.96	225	153849	33.672	ng	98
35) Caprolactam	11.59	113	74139m	28.113	ng	
36) 4-Chloro-3-methylphenol	11.92	107	250556	33.356	ng	93
37) 2-Methylnaphthalene	12.30	142	583637	37.288	ng	100
38) 1-Methylnaphthalene	12.53	142	553826	36.228	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.67	216	304274	36.551	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.64	237	378511	102.086	ng	100
43) 2,4,6-Trichlorophenol	12.91	196	200699	38.159	ng	99
44) 2,4,5-Trichlorophenol	12.98	196	218548	39.837	ng	98
46) 1,1'-Biphenyl	13.32	154	734803	33.270	ng	99
47) 2-Chloronaphthalene	13.36	162	566457	34.771	ng	99
48) 2-Nitroaniline	13.57	65	141314	29.522	ng	87
49) Acenaphthylene	14.21	152	912801	36.606	ng	99
50) Dimethylphthalate	13.95	163	873066	42.910	ng	100
51) 2,6-Dinitrotoluene	14.07	165	156146	33.812	ng	94
52) Acenaphthene	14.55	154	525396	34.775	ng	98
53) 3-Nitroaniline	14.40	138	109331	22.147	ng	94
54) 2,4-Dinitrophenol	14.60	184	96473	39.840	ng	# 89
55) Dibenzofuran	14.89	168	843031	34.297	ng	100
56) 4-Nitrophenol	14.70	139	346066	96.088	ng	94
57) 2,4-Dinitrotoluene	14.86	165	207344	33.114	ng	# 76
58) Fluorene	15.53	166	664131	35.014	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.10	232	192665	39.993	ng	97
60) Diethylphthalate	15.30	149	675882	32.735	ng	100
61) 4-Chlorophenyl-phenylether	15.53	204	348889	33.108	ng	98
62) 4-Nitroaniline	15.56	138	149731	31.046	ng	98
63) Azobenzene	15.82	77	579891	30.390	ng	94
65) 4,6-Dinitro-2-methylphenol	15.61	198	102264	28.539	ng	95
66) n-Nitrosodiphenylamine	15.75	169	580965	36.137	ng	100
67) 4-Bromophenyl-phenylether	16.42	248	222316	38.438	ng	94
68) Hexachlorobenzene	16.53	284	252816	38.370	ng	95
69) Atrazine	16.69	200	203051	35.639	ng	98
70) Pentachlorophenol	16.87	266	277646	83.141	ng	99
71) Phenanthrene	17.27	178	980639	36.322	ng	98
72) Anthracene	17.36	178	998648	37.069	ng	99
73) Carbazole	17.63	167	863677	30.957	ng	99
74) Di-n-butylphthalate	18.18	149	1085016	33.340	ng	100
75) Fluoranthene	19.28	202	1026466	32.271	ng	99
77) Benzidine	19.47	184	433110	34.306	ng	99
78) Pyrene	19.65	202	1016494	40.105	ng	99
80) Butylbenzylphthalate	20.54	149	422934	35.936	ng	93
81) Benzo(a)anthracene	21.39	228	897091	35.720	ng	98
82) 3,3'-Dichlorobenzidine	21.33	252	291678	28.065	ng	98
83) Chrysene	21.45	228	845215	35.010	ng	99
84) Bis(2-ethylhexyl)phthalate	21.30	149	619409	35.489	ng	98
85) Di-n-octyl phthalate	22.20	149	950697	31.743	ng	98
86) Indeno(1,2,3-cd)pyrene	26.08	276	1057125	38.207	ng	99
88) Benzo(b)fluoranthene	23.02	252	917880	38.748	ng	98
89) Benzo(k)fluoranthene	23.07	252	851107	36.634	ng	98
90) Benzo(a)pyrene	23.62	252	856792	37.951	ng	98
91) Dibenzo(a,h)anthracene	26.09	278	887710	39.726	ng	99
92) Benzo(g,h,i)perylene	26.81	276	878613	40.956	ng	98

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

