

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN021420\
 Data File : BN009726.D
 Acq On : 14 Feb 2020 15:25
 Operator : CG/JU
 Sample : K5695-13
 Misc : MDL-WATER-06
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MDL-WATER-06

Manual Integrations
 APPROVED

mohammad
 2/17/2020 3:50:13 PM

Quant Time: Feb 17 01:55:08 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Feb 17 00:47:25 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.42	152	108811	20.00	ng/ul	0.00
18) Naphthalene-d8	10.16	136	447014	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.06	164	286057	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.81	188	613799	20.00	ng/ul	0.00
78) Chrysene-d12	21.03	240	599942	20.00	ng/ul	0.00
86) Perylene-d12	23.14	264	595728	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.06	96	11579	4.38	ng/uL	0.00
5) Phenol-d5	6.62	99	262971	28.41	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.78	67	190107	29.76	ng/ul	0.00
9) 2-Chlorophenol-d4	6.96	132	208232	29.88	ng/ul	0.00
13) 4-Methylphenol-d8	8.13	113	207491	28.18	ng/ul	0.00
19) Nitrobenzene-d5	8.56	128	100414	30.70	ng/ul	0.00
22) 2-Nitrophenol-d4	9.28	143	109433	30.59	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.80	165	196539	28.58	ng/ul	0.00
29) 4-Chloroaniline-d4	10.32	131	278540	36.05	ng/ul	0.00
44) Dimethylphthalate-d6	13.48	166	653688	30.59	ng/ul	0.00
47) Acenaphthylene-d8	13.75	160	784133	31.16	ng/ul	0.00
52) 4-Nitrophenol-d4	14.30	143	93512	23.96	ng/ul	0.00
58) Fluorene-d10	15.06	176	561912	30.46	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.22	200	91422	23.98	ng/ul	0.00
71) Anthracene-d10	16.91	188	859459	30.55	ng/ul	0.00
79) Pyrene-d10	19.22	212	928437	29.61	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.00	264	906767	30.48	ng/ul	0.00

Target Compounds

					Qvalue
2) 1,4-Dioxane	3.09	88	2713	0.89 ng/uL	99
4) Benzaldehyde	6.59	77	18999	3.09 ng/ul	97
6) Phenol	6.65	94	26704	2.69 ng/ul	96
8) Bis(2-Chloroethyl)ether	6.87	93	22175	2.84 ng/ul	99
10) 2-Chlorophenol	6.99	128	19542	2.66 ng/ul	95
11) 2-Methylphenol	7.88	108	16965	2.34 ng/ul	94
12) 2,2'-oxybis(1-Chloropropan	7.95	45	56591m	2.96 ng/ul	
14) Acetophenone	8.24	105	31331	2.61 ng/ul	99
15) N-Nitroso-di-n-propylamine	8.23	70	17239	2.75 ng/ul	99
16) 4-Methylphenol	8.19	108	20943	2.63 ng/ul	94
17) Hexachloroethane	8.46	117	9936	2.98 ng/ul	88
20) Nitrobenzene	8.60	77	27830	2.89 ng/ul	98
21) Isophorone	9.13	82	42683	2.52 ng/ul	96
23) 2-Nitrophenol	9.30	139	10939	2.73 ng/ul#	79
24) 2,4-Dimethylphenol	9.38	107	23557	2.71 ng/ul	93
25) Bis(2-Chloroethoxy)methane	9.62	93	28437	2.74 ng/ul	97
27) 2,4-Dichlorophenol	9.83	162	18893	2.72 ng/ul#	74
28) Naphthalene	10.22	128	69485	2.88 ng/ul	98
30) 4-Chloroaniline	10.35	127	22745	2.87 ng/ul	95
31) Hexachlorobutadiene	10.50	225	13119	2.83 ng/ul#	89
32) Caprolactam	11.18	113	3895m	1.67 ng/ul	
33) 4-Chloro-3-methylphenol	11.49	107	18042	2.27 ng/ul	94
34) 2-Methylnaphthalene	11.85	142	46758	2.79 ng/ul	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
35) 1-Methylnaphthalene	12.06	142	48640	2.90	ng/uL	98
37) 1,2,4,5-Tetrachlorobenzene	12.23	216	24695	2.91	ng/ul	93
38) Hexachlorocyclopentadiene	12.19	237	3117	0.85	ng/ul#	74
39) 2,4,6-Trichlorophenol	12.49	196	10975	2.04	ng/ul	99
40) 2,4,5-Trichlorophenol	12.56	196	14311m	2.48	ng/ul	
41) 1,1'-Biphenyl	12.88	154	62383	2.84	ng/ul	96
42) 2-Chloronaphthalene	12.92	162	47368	2.71	ng/ul	95
43) 2-Nitroaniline	13.15	65	14041	2.20	ng/ul	89
45) Dimethylphthalate	13.53	163	65265	2.93	ng/ul	99
46) 2,6-Dinitrotoluene	13.66	165	9675	2.22	ng/ul#	82
48) Acenaphthylene	13.77	152	77577	2.85	ng/ul	98
49) 3-Nitroaniline	14.00	138	8833	2.19	ng/ul#	90
50) Acenaphthene	14.12	153	53397	2.86	ng/ul	99
51) 2,4-Dinitrophenol	14.23	184	2145	0.86	ng/ul#	89
53) 4-Nitrophenol	14.31	109	9138	2.60	ng/ul	91
54) Dibenzofuran	14.46	168	75369	2.88	ng/ul	95
55) 2,4-Dinitrotoluene	14.46	165	15877	2.45	ng/ul#	99
56) 2,3,4,6-Tetrachlorophenol	14.70	232	11802	2.38	ng/ul#	92
57) Diethylphthalate	14.92	149	61445	2.68	ng/ul	97
59) Fluorene	15.12	166	62105	2.82	ng/ul	96
60) 4-Chlorophenyl-phenylether	15.12	204	31265	2.88	ng/ul	96
61) 4-Nitroaniline	15.17	138	7384	1.50	ng/ul	90
64) 4,6-Dinitro-2-methylphenol	15.23	198	10002	2.44	ng/ul#	75
65) N-Nitrosodiphenylamine	15.34	169	49318	2.71	ng/ul	94
66) 4-Bromophenyl-phenylether	16.02	248	16864	2.75	ng/ul	98
67) Hexachlorobenzene	16.13	284	18647	2.74	ng/ul	97
68) Atrazine	16.31	200	16020	2.32	ng/ul	91
69) Pentachlorophenol	16.49	266	6049	1.54	ng/ul#	78
70) Phenanthrene	16.86	178	100162	2.86	ng/ul	98
72) Anthracene	16.95	178	102444	2.86	ng/ul	99
73) 1,2,3,4-Tetrachlorobenzene	12.85	216	27331	3.05	ng/uL	95
74) Pentachlorobenzene	14.39	250	23735	2.74	ng/uL	95
75) Carbazole	17.23	167	72674	2.30	ng/ul	98
76) Di-n-butylphthalate	17.81	149	86413	2.22	ng/ul	99
77) Fluoranthene	18.89	202	104533	2.55	ng/ul	99
80) Pyrene	19.25	202	119517	2.76	ng/ul	99
81) Butylbenzylphthalate	20.18	149	35107	1.97	ng/ul	97
82) 3,3'-Dichlorobenzidine	20.96	252	29127	2.23	ng/ul#	92
83) Benzo(a)anthracene	21.02	228	107542	2.61	ng/ul	98
84) Bis(2-ethylhexyl)phthalate	20.97	149	54998	2.03	ng/ul	98
85) Chrysene	21.07	228	111800	2.75	ng/ul	98
87) Di-n-octyl phthalate	21.82	149	88852	1.97	ng/ul	100
88) Benzo(b)fluoranthene	22.52	252	105579m	2.73	ng/ul	
89) Benzo(k)fluoranthene	22.56	252	100381	2.74	ng/ul	99
91) Benzo(a)pyrene	23.05	252	102908	2.96	ng/ul	99
92) Indeno(1,2,3-cd)pyrene	25.20	276	108852	2.64	ng/ul	96
93) Dibenzo(a,h)anthracene	25.20	278	94035	2.67	ng/ul	97
94) Benzo(g,h,i)perylene	25.83	276	93275	2.70	ng/ul	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
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(#) = qualifier out of range (m) = manual integration (+) = signals summed

