

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060820\
 Data File : BN010808.D
 Acq On : 08 Jun 2020 15:01
 Operator : CG/JU
 Sample : L1955-14
 Misc : MDL-WATER-07
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MDL-WATER-07

Quant Time: Jun 08 15:31:14 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN060320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 08 10:24:50 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	1777	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	5571	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	2893	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	5582	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	4603	0.40	ng/ul	0.00
20) Perylene-d12	23.29	264	4343	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	2476	0.19	ng/ul	0.01
14) Fluoranthene-d10	18.97	212	5022	0.21	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.37	128	1032	0.063	ng/ul#	82
5) 2-Methylnaphthalene	12.00	142	602	0.058	ng/ul	98
7) Acenaphthylene	13.90	152	804	0.062	ng/ul#	82
8) Acenaphthene	14.25	153	670	0.062	ng/ul	95
9) Fluorene	15.25	166	702	0.059	ng/ul#	94
11) Pentachlorophenol	16.62	266	77	0.059	ng/ul#	91
12) Phenanthrene	16.98	178	1129	0.062	ng/ul#	86
13) Anthracene	17.08	178	861	0.057	ng/ul#	78
15) Fluoranthene	19.00	202	1267	0.062	ng/ul	90
17) Pyrene	19.37	202	1260	0.063	ng/ul#	87
18) Benzo(a)anthracene	21.13	228	963	0.059	ng/ul#	89
19) Chrysene	21.18	228	1318	0.063	ng/ul	94
21) Benzo(b)fluoranthene	22.66	252	1161	0.060	ng/ul#	48
22) Benzo(k)fluoranthene	22.70	252	1344	0.066	ng/ul#	51
23) Benzo(a)pyrene	23.21	252	1118	0.068	ng/ul#	34
24) Indeno(1,2,3-cd)pyrene	25.42	276	1370	0.061	ng/ul#	54
25) Dibenzo(a,h)anthracene	25.44	278	1129	0.062	ng/ul#	57
26) Benzo(g,h,i)perylene	26.06	276	1309	0.063	ng/ul#	80

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

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