

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM081820\
 Data File : BM027127.D
 Acq On : 18 Aug 2020 15:31
 Operator : JU/CG
 Sample : L1955-09
 Misc : MDL-W-02-0.07PPM
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-WATER-02

Quant Time: Aug 18 16:10:53 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM081720.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 18 11:17:38 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	930	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	2883	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.27	164	1862	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	4245	0.40	ng/ul	0.00
16) Chrysene-d12	21.21	240	4168	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	4153	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	1761	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.04	212	4597	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.45	128	398	0.046	ng/ul#	57
5) 2-Methylnaphthalene	12.08	142	267	0.045	ng/ul	99
7) Acenaphthylene	13.99	152	356	0.044	ng/ul#	74
8) Acenaphthene	14.33	153	325	0.046	ng/ul#	92
9) Fluorene	15.32	166	388	0.046	ng/ul#	94
11) Pentachlorophenol	16.68	266	62	0.044	ng/ul#	79
12) Phenanthrene	17.05	178	679	0.047	ng/ul#	79
13) Anthracene	17.14	178	545	0.042	ng/ul#	68
15) Fluoranthene	19.07	202	834	0.048	ng/ul	75
17) Pyrene	19.44	202	833	0.047	ng/ul#	82
18) Benzo(a)anthracene	21.19	228	786	0.047	ng/ul#	86
19) Chrysene	21.24	228	967	0.054	ng/ul#	86
21) Benzo(b)fluoranthene	22.74	252	912	0.051	ng/ul#	38
22) Benzo(k)fluoranthene	22.79	252	868	0.047	ng/ul#	41
23) Benzo(a)pyrene	23.30	252	860	0.053	ng/ul#	28
24) Indeno(1,2,3-cd)pyrene	25.59	276	1029	0.046	ng/ul#	73
25) Dibenzo(a,h)anthracene	25.60	278	782	0.043	ng/ul#	41
26) Benzo(g,h,i)perylene	26.25	276	855	0.045	ng/ul#	71

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

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