

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM081920\
 Data File : BM027135.D
 Acq On : 19 Aug 2020 09:50
 Operator : JU/CG
 Sample : L1955-11
 Misc : MDL-W-04-0.07PPM
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-WATER-04

Quant Time: Aug 19 11:17:27 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 : Wed Aug 19 11:15:05 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	900	0.40	ng/ul	0.00
2) Naphthalene-d8	10.40	136	2844	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	1886	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	4286	0.40	ng/ul	0.00
16) Chrysene-d12	21.21	240	4193	0.40	ng/ul	0.00
20) Perylene-d12	23.39	264	4176	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	1268	0.26	ng/ul	0.00
14) Fluoranthene-d10	19.04	212	4429	0.36	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.45	128	386	0.045	ng/ul#	59
5) 2-Methylnaphthalene	12.07	142	198	0.034	ng/ul	92
7) Acenaphthylene	13.99	152	384	0.047	ng/ul#	75
8) Acenaphthene	14.33	153	354	0.050	ng/ul	95
9) Fluorene	15.32	166	414	0.049	ng/ul#	96
11) Pentachlorophenol	16.69	266	59	0.042	ng/ul#	71
12) Phenanthrene	17.05	178	725	0.050	ng/ul#	79
13) Anthracene	17.15	178	589	0.045	ng/ul#	69
15) Fluoranthene	19.07	202	924	0.053	ng/ul	79
17) Pyrene	19.44	202	967	0.054	ng/ul#	84
18) Benzo(a)anthracene	21.19	228	1048	0.063	ng/ul	91
19) Chrysene	21.24	228	1388	0.077	ng/ul	91
21) Benzo(b)fluoranthene	22.75	252	1278	0.071	ng/ul#	59
22) Benzo(k)fluoranthene	22.79	252	1348	0.072	ng/ul#	63
23) Benzo(a)pyrene	23.30	252	1187	0.073	ng/ul#	54
24) Indeno(1,2,3-cd)pyrene	25.58	276	1460	0.065	ng/ul#	92
25) Dibenzo(a,h)anthracene	25.60	278	1137	0.062	ng/ul#	54
26) Benzo(g,h,i)perylene	26.25	276	1280	0.066	ng/ul#	83

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

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