

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM081820\
 Data File : BM027129.D
 Acq On : 18 Aug 2020 16:44
 Operator : JU/CG
 Sample : L1955-10
 Misc : MDL-W-03-0.07PPM
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-WATER-03

Quant Time: Aug 18 17:34:32 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	848	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	2753	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.27	164	1455	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	4085	0.40	ng/ul	0.00
16) Chrysene-d12	21.20	240	4030	0.40	ng/ul	0.00
20) Perylene-d12	23.39	264	4056	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	1735	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.04	212	4470	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.46	128	412	0.050	ng/ul#	59
5) 2-Methylnaphthalene	12.08	142	276	0.049	ng/ul	98
7) Acenaphthylene	13.99	152	304	0.048	ng/ul#	76
8) Acenaphthene	14.33	153	280	0.051	ng/ul	93
9) Fluorene	15.32	166	400	0.061	ng/ul#	91
11) Pentachlorophenol	16.69	266	47	0.035	ng/ul#	89
12) Phenanthrene	17.05	178	703	0.051	ng/ul#	80
13) Anthracene	17.15	178	532	0.042	ng/ul#	72
15) Fluoranthene	19.07	202	853	0.051	ng/ul	79
17) Pyrene	19.44	202	884	0.051	ng/ul#	81
18) Benzo(a)anthracene	21.19	228	865	0.054	ng/ul	93
19) Chrysene	21.24	228	1074	0.062	ng/ul	93
21) Benzo(b)fluoranthene	22.74	252	994	0.057	ng/ul#	47
22) Benzo(k)fluoranthene	22.79	252	1208	0.067	ng/ul#	48
23) Benzo(a)pyrene	23.30	252	909	0.058	ng/ul#	36
24) Indeno(1,2,3-cd)pyrene	25.59	276	1173	0.054	ng/ul#	90
25) Dibenzo(a,h)anthracene	25.61	278	960	0.054	ng/ul#	44
26) Benzo(g,h,i)perylene	26.26	276	1022	0.055	ng/ul#	71

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

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