

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM081920\
 Data File : BM027137.D
 Acq On : 19 Aug 2020 11:02
 Operator : JU/CG
 Sample : L1955-12
 Misc : MDL-W-05-0.07PPM
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-WATER-05

Quant Time: Aug 19 11:53:36 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	745	0.40	ng/ul	0.00
2) Naphthalene-d8	10.40	136	2316	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	1536	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	3596	0.40	ng/ul	0.00
16) Chrysene-d12	21.20	240	3493	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	3440	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	1425	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.04	212	3978	0.39	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.45	128	387	0.056	ng/ul#	55
5) 2-Methylnaphthalene	12.08	142	262	0.055	ng/ul	93
7) Acenaphthylene	13.98	152	384	0.058	ng/ul#	72
8) Acenaphthene	14.32	153	356	0.061	ng/ul	94
9) Fluorene	15.32	166	432	0.062	ng/ul#	94
11) Pentachlorophenol	16.69	266	56	0.047	ng/ul	96
12) Phenanthrene	17.05	178	719	0.059	ng/ul#	81
13) Anthracene	17.15	178	631	0.057	ng/ul#	70
15) Fluoranthene	19.07	202	886	0.061	ng/ul	80
17) Pyrene	19.44	202	913	0.061	ng/ul#	85
18) Benzo(a)anthracene	21.19	228	856	0.062	ng/ul#	87
19) Chrysene	21.24	228	1079	0.072	ng/ul#	92
21) Benzo(b)fluoranthene	22.75	252	999	0.068	ng/ul#	55
22) Benzo(k)fluoranthene	22.79	252	1112	0.072	ng/ul#	50
23) Benzo(a)pyrene	23.30	252	890	0.067	ng/ul#	41
24) Indeno(1,2,3-cd)pyrene	25.59	276	1141	0.062	ng/ul#	90
25) Dibenzo(a,h)anthracene	25.60	278	953	0.064	ng/ul#	47
26) Benzo(g,h,i)perylene	26.25	276	978	0.061	ng/ul#	69

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

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