

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082020\
 Data File : BM027155.D
 Acq On : 20 Aug 2020 13:46
 Operator : JU/CG
 Sample : L1955-08
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-WATER-01

Manual Integrations
 APPROVED

mohammad
 8/21/2020 12:49:28 PM

Quant Time: Aug 20 14:18:00 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 : Thu Aug 20 06:55:50 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	483	0.40	ng/ul	0.00
2) Naphthalene-d8	10.40	136	1400	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	971	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	2772	0.40	ng/ul	0.00
16) Chrysene-d12	21.20	240	2699	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	2789	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	870	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.04	212	3048	0.39	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.45	128	235	0.056	ng/ul#	33
5) 2-Methylnaphthalene	12.07	142	152	0.053	ng/ul	98
7) Acenaphthylene	13.98	152	239	0.057	ng/ul#	58
8) Acenaphthene	14.33	153	213	0.058	ng/ul	95
9) Fluorene	15.32	166	249	0.057	ng/ul#	89
11) Pentachlorophenol	16.69	266	54m	0.059	ng/ul	
12) Phenanthrene	17.05	178	548	0.058	ng/ul#	73
13) Anthracene	17.15	178	457	0.054	ng/ul#	63
15) Fluoranthene	19.07	202	678	0.060	ng/ul	74
17) Pyrene	19.43	202	697	0.060	ng/ul#	79
18) Benzo(a)anthracene	21.19	228	620	0.058	ng/ul#	88
19) Chrysene	21.24	228	787	0.068	ng/ul#	90
21) Benzo(b)fluoranthene	22.75	252	715	0.060	ng/ul#	43
22) Benzo(k)fluoranthene	22.79	252	769	0.062	ng/ul#	44
23) Benzo(a)pyrene	23.30	252	651	0.060	ng/ul#	32
24) Indeno(1,2,3-cd)pyrene	25.58	276	858	0.057	ng/ul#	82
25) Dibenzo(a,h)anthracene	25.61	278	722	0.059	ng/ul#	27
26) Benzo(g,h,i)perylene	26.25	276	761	0.059	ng/ul#	68

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

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