

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030920\
 Data File : BM025466.D
 Acq On : 10 Mar 2020 23:18
 Operator : CG/JU
 Sample : L1617-12
 Misc :
 ALS Vial : 60 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Mar 11 00:16:51 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 10 23:48:01 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	3948	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	16949	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	12075	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	25590	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	18010	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	19510	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	7766	0.25	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	16669	0.24	ng/ul	0.00
Target Compounds				Ovalue		
8) Acenaphthene	14.35	153	2711	0.063	ng/ul	98
9) Fluorene	15.34	166	8901	0.169	ng/ul	99
12) Phenanthrene	17.07	178	72786	0.898	ng/ul	100
13) Anthracene	17.16	178	4520	0.064	ng/ul	97
15) Fluoranthene	19.09	202	39176	0.446	ng/ul	99
17) Pyrene	19.45	202	27367	0.339	ng/ul	99
18) Benzo(a)anthracene	21.20	228	3340	0.052	ng/ul	96
19) Chrysene	21.25	228	3303	0.046	ng/ul	98
21) Benzo(b)fluoranthene	22.76	252	2346m	0.030	ng/ul	

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

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