

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM091719\
 Data File : BM022702.D
 Acq On : 17 Sep 2019 12:17
 Operator : HP/JU
 Sample : K4780-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 H0BM9

Manual Integrations
 APPROVED

mohammad
 9/19/2019 4:44:22 PM

Quant Time: Sep 17 18:12:37 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM091019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 10 15:06:45 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	9887	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	41451	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	25314	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	53769m	0.40	ng/ul	-0.01
16) Chrysene-d12	21.39	240	48369m	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	49253	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	18410	0.27	ng/ul	0.00
14) Fluoranthene-d10	19.23	212	44444m	0.30	ng/ul	0.00
Target Compounds				Qvalue		
17) Pyrene	19.62	202	5897	0.027	ng/ul#	75
19) Chrysene	21.42	228	4970	0.026	ng/ul#	92
21) Benzo(b)fluoranthene	22.98	252	9092m	0.042	ng/ul	
23) Benzo(a)pyrene	23.57	252	4637	0.025	ng/ul#	1

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

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