

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091619\
 Data File : BM022687.D
 Acq On : 16 Sep 2019 12:43
 Operator : HP/JU
 Sample : K4780-15MSD
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 H0BM3MSD

Manual Integrations
 APPROVED

mohammad
 9/17/2019 4:11:21 PM

Quant Time: Sep 17 01:53:45 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	9937	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	41531	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	25396	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	54398m	0.40	ng/ul	0.00
16) Chrysene-d12	21.39	240	49252m	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	53119	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	19478	0.29	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	56566m	0.38	ng/ul	0.00
Target Compounds				Ovalue		
8) Acenaphthene	14.54	153	26200	0.266	ng/ul	99
11) Pentachlorophenol	16.86	266	3228m	0.293	ng/ul	
17) Pyrene	19.63	202	75908m	0.340	ng/ul	

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

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