

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM091719\
 Data File : BM022704.D
 Acq On : 17 Sep 2019 13:30
 Operator : HP/JU
 Sample : K4780-04
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 H0BN8

Manual Integrations
 APPROVED

mohammad
 9/19/2019 4:44:34 PM

Quant Time: Sep 17 18:20:15 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	9665	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	41077	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	25303	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	51699m	0.40	ng/ul	0.00
16) Chrysene-d12	21.39	240	49983m	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	55416	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	18871	0.28	ng/ul	0.00
14) Fluoranthene-d10	19.23	212	51765m	0.36	ng/ul	0.00
Target Compounds				Qvalue		
17) Pyrene	19.62	202	6839	0.030	ng/ul#	77
18) Benzo(a)anthracene	21.37	228	4141	0.022	ng/ul#	91
19) Chrysene	21.42	228	5699	0.028	ng/ul#	93
21) Benzo(b)fluoranthene	22.98	252	10021m	0.042	ng/ul	
23) Benzo(a)pyrene	23.57	252	5403	0.026	ng/ul#	1
26) Benzo(g,h,i)perylene	26.68	276	4295	0.021	ng/ul#	72

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

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