

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091019\
 Data File : BM022630.D
 Acq On : 10 Sep 2019 21:25
 Operator : HP/JU
 Sample : K4708-08
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 CB5Q5

Manual Integrations
 APPROVED

mohammad
 9/11/2019 8:21:51 AM

Quant Time: Sep 10 23:58:16 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	7.85	152	17014	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	70885	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.48	164	37667	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	80936m	0.40	ng/ul	0.00
16) Chrysene-d12	21.39	240	62443m	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	65171	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.23	152	28993	0.25	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	55240m	0.25	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.25	178	10594m	0.041	ng/ul	
15) Fluoranthene	19.27	202	22571m	0.081	ng/ul	
17) Pyrene	19.63	202	19841	0.070	ng/ul#	95
18) Benzo(a)anthracene	21.37	228	9044	0.038	ng/ul#	91
19) Chrysene	21.42	228	11950	0.048	ng/ul#	93
21) Benzo(b)fluoranthene	22.98	252	15611m	0.055	ng/ul	
23) Benzo(a)pyrene	23.57	252	9366	0.038	ng/ul#	1
26) Benzo(g,h,i)perylene	26.69	276	7752	0.032	ng/ul#	66

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

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