

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091019\
 Data File : BM022631.D
 Acq On : 10 Sep 2019 22:01
 Operator : HP/JU
 Sample : K4708-09
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 CB5Q6

Manual Integrations
 APPROVED

mohammad
 9/11/2019 8:21:54 AM

Quant Time: Sep 11 00:10:43 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	15991	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	66848	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.48	164	35723	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	74691m	0.40	ng/ul	0.00
16) Chrysene-d12	21.39	240	57274m	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	53384m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.23	152	37590	0.34	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	69621m	0.34	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.69	128	7786	0.041	ng/ul#	92
5) 2-Methylnaphthalene	12.31	142	3243	0.025	ng/ul	99
12) Phenanthrene	17.25	178	16473m	0.069	ng/ul	
15) Fluoranthene	19.27	202	33682m	0.131	ng/ul	
17) Pyrene	19.63	202	29073m	0.112	ng/ul	
18) Benzo(a)anthracene	21.37	228	13921m	0.064	ng/ul	
19) Chrysene	21.42	228	17946	0.078	ng/ul	94
21) Benzo(b)fluoranthene	22.98	252	22707m	0.098	ng/ul	
22) Benzo(k)fluoranthene	23.02	252	9492m	0.040	ng/ul	
23) Benzo(a)pyrene	23.57	252	14079m	0.069	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.98	276	11523m	0.047	ng/ul	
26) Benzo(g,h,i)perylene	26.69	276	12149m	0.062	ng/ul	

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

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