

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091019\
 Data File : BM022656.D
 Acq On : 11 Sep 2019 14:28
 Operator : HP/JU
 Sample : K4706-08
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BF574

Manual Integrations
 APPROVED

mohammad
 9/16/2019 8:54:57 AM

Quant Time: Sep 12 01:16:53 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	17724	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	77502	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.48	164	45590	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	102184m	0.40	ng/ul	0.00
16) Chrysene-d12	21.39	240	87905m	0.40	ng/ul	0.00
20) Perylene-d12	23.68	264	82255	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.23	152	34031	0.27	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	92983m	0.33	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.69	128	477905	2.164	ng/ul	100
5) 2-Methylnaphthalene	12.30	142	82565	0.545	ng/ul	100
7) Acenaphthylene	14.20	152	7248	0.034	ng/ul#	73
8) Acenaphthene	14.54	153	240480	1.361	ng/ul	98
9) Fluorene	15.52	166	349506	1.727	ng/ul	99
12) Phenanthrene	17.26	178	1088004m	3.342	ng/ul	
13) Anthracene	17.35	178	62238m	0.223	ng/ul	
15) Fluoranthene	19.27	202	209883m	0.597	ng/ul	
17) Pyrene	19.63	202	126352m	0.317	ng/ul	
18) Benzo(a)anthracene	21.38	228	18530	0.055	ng/ul	95
19) Chrysene	21.43	228	11872m	0.034	ng/ul	
21) Benzo(b)fluoranthene	22.99	252	11522m	0.032	ng/ul	

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

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