

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061019\
 Data File : BN006223.D
 Acq On : 10 Jun 2019 15:24
 Operator : JU/SJ
 Sample : K3016-09
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A51C8

Manual Integrations
 APPROVED

mohammad
 6/11/2019 4:28:03 PM

Quant Time: Jun 11 01:02:12 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN053019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Jun 09 05:10:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	4819	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	18161	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	9401	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	19036m	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	14911m	0.40	ng/ul	0.00
20) Perylene-d12	23.50	264	13724m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	5827	0.21	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	14667m	0.29	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.09	178	2616m	0.043	ng/ul	
15) Fluoranthene	19.12	202	6466m	0.096	ng/ul	
17) Pyrene	19.49	202	6284m	0.080	ng/ul	
18) Benzo(a)anthracene	21.25	228	2294	0.035	ng/ul#	86
19) Chrysene	21.30	228	3670	0.059	ng/ul#	92
21) Benzo(b)fluoranthene	22.83	252	4593m	0.080	ng/ul	
22) Benzo(k)fluoranthene	22.87	252	2367m	0.042	ng/ul	
23) Benzo(a)pyrene	23.40	252	2928m	0.055	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.73	276	2190m	0.038	ng/ul	
26) Benzo(g,h,i)perylene	26.41	276	2256m	0.047	ng/ul	

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

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