

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061019\
 Data File : BN006218.D
 Acq On : 10 Jun 2019 12:32
 Operator : JU/SJ
 Sample : K3016-10
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A51C9

Manual Integrations
 APPROVED

mohammad
 6/11/2019 4:27:46 PM

Quant Time: Jun 11 00:41:41 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	4112	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	15630	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	8229	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	15995m	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	12584m	0.40	ng/ul	0.00
20) Perylene-d12	23.50	264	11615m	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	5971	0.25	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	13134m	0.31	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.09	178	3372m	0.066	ng/ul	
15) Fluoranthene	19.12	202	7267m	0.128	ng/ul	
17) Pyrene	19.49	202	6687m	0.101	ng/ul	
18) Benzo(a)anthracene	21.25	228	3135	0.056	ng/ul#	83
19) Chrysene	21.31	228	3535	0.068	ng/ul#	91
21) Benzo(b)fluoranthene	22.84	252	4845m	0.100	ng/ul	
22) Benzo(k)fluoranthene	22.87	252	2281m	0.048	ng/ul	
23) Benzo(a)pyrene	23.40	252	2977m	0.066	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.73	276	2227m	0.046	ng/ul	
26) Benzo(g,h,i)perylene	26.42	276	1582m	0.039	ng/ul	

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

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