

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060819\
 Data File : BN006210.D
 Acq On : 09 Jun 2019 03:07
 Operator : JU/SJ
 Sample : K3017-10
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A51E3

Manual Integrations
 APPROVED

mohammad
 6/11/2019 4:33:05 PM

Quant Time: Jun 10 02:40:01 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	7.64	152	5702	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	22318	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	11993	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	22701m	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	18285m	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	14991m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	12364	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	25034m	0.41	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.01	152	1608	0.024	ng/ul#	82
12) Phenanthrene	17.09	178	7021	0.096	ng/ul	96
15) Fluoranthene	19.12	202	15303m	0.190	ng/ul	
17) Pyrene	19.49	202	15512	0.161	ng/ul	97
18) Benzo(a)anthracene	21.24	228	4969m	0.061	ng/ul	
19) Chrysene	21.29	228	7601m	0.100	ng/ul	
21) Benzo(b)fluoranthene	22.82	252	10118m	0.162	ng/ul	
22) Benzo(k)fluoranthene	22.86	252	3573m	0.059	ng/ul	
23) Benzo(a)pyrene	23.39	252	5918	0.102	ng/ul#	59
24) Indeno(1,2,3-cd)pyrene	25.72	276	4761	0.076	ng/ul#	83
26) Benzo(g,h,i)perylene	26.41	276	4742	0.090	ng/ul#	81

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

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