

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061719\
 Data File : BN006365.D
 Acq On : 18 Jun 2019 01:02
 Operator : JU/SJ
 Sample : K3335-23
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A41Y1

Manual Integrations
 APPROVED

mohammad
 6/21/2019 8:22:59 AM

Quant Time: Jun 18 04:42:36 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 : Mon Jun 17 11:13:07 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	3945	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	15869	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	8753	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	16492m	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	9614m	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	8148m	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	3862	0.16	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	7901m	0.18	ng/ul	0.00
Target Compounds						
12) Phenanthrene	17.09	178	5741	0.108	ng/ul	96
13) Anthracene	17.18	178	1781m	0.035	ng/ul	
15) Fluoranthene	19.11	202	14302m	0.245	ng/ul	
17) Pyrene	19.48	202	13837m	0.273	ng/ul	
18) Benzo(a)anthracene	21.25	228	5165	0.121	ng/ul	96
19) Chrysene	21.30	228	4934	0.124	ng/ul	96
21) Benzo(b)fluoranthene	22.83	252	5028m	0.148	ng/ul	
22) Benzo(k)fluoranthene	22.86	252	1773m	0.054	ng/ul	
23) Benzo(a)pyrene	23.39	252	3853m	0.122	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.74	276	2198m	0.064	ng/ul	
26) Benzo(g,h,i)perylene	26.42	276	2011m	0.070	ng/ul	

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

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