

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061719\
 Data File : BN006349.D
 Acq On : 17 Jun 2019 15:50
 Operator : JU/SJ
 Sample : K3332-11
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A41W5

Manual Integrations
 APPROVED

mohammad
 6/20/2019 10:51:10 AM

Quant Time: Jun 18 02:30:34 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	4877	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	19623	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	10236	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	18012m	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	11925m	0.40	ng/ul	0.00
20) Perylene-d12	23.51	264	10370m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	5116	0.17	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	11049m	0.23	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.09	178	3028m	0.052	ng/ul	93
15) Fluoranthene	19.12	202	8362m	0.131	ng/ul	
17) Pyrene	19.49	202	7577m	0.121	ng/ul	
18) Benzo(a)anthracene	21.25	228	3516	0.067	ng/ul	
19) Chrysene	21.31	228	3038m	0.061	ng/ul	
21) Benzo(b)fluoranthene	22.84	252	5575m	0.129	ng/ul	
22) Benzo(k)fluoranthene	22.88	252	1337m	0.032	ng/ul	
23) Benzo(a)pyrene	23.40	252	3320m	0.082	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.75	276	2426m	0.056	ng/ul	
26) Benzo(g,h,i)perylene	26.43	276	2004m	0.055	ng/ul	

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

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