

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN062419\
 Data File : BN006534.D
 Acq On : 24 Jun 2019 17:02
 Operator : HP/JU
 Sample : K3362-13
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A41X9

Manual Integrations
 APPROVED

mohammad
 6/25/2019 10:20:03 PM

Quant Time: Jun 25 01:51:40 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	7.63	152	4943	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.41	136	18550	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.28	164	9247	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	17780m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.27	240	9871m	0.40	ng/ul	0.00
20) Perylene-d12	23.51	264	9616m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	7234	0.25	ng/ul	-0.02
14) Fluoranthene-d10	19.08	212	13599m	0.28	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.08	178	2436m	0.043	ng/ul	
15) Fluoranthene	19.11	202	4293m	0.068	ng/ul	
17) Pyrene	19.48	202	3723	0.072	ng/ul#	90
18) Benzo(a)anthracene	21.25	228	1630	0.037	ng/ul#	81
19) Chrysene	21.30	228	1457	0.036	ng/ul#	82
21) Benzo(b)fluoranthene	22.84	252	2368	0.059	ng/ul#	37
23) Benzo(a)pyrene	23.41	252	1220m	0.033	ng/ul	

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

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