

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061319\
 Data File : BN006320.D
 Acq On : 13 Jun 2019 20:34
 Operator : JU/SJ
 Sample : K3213-14
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DB8K5

Manual Integrations
 APPROVED

mohammad
 6/14/2019 8:49:35 AM

Quant Time: Jun 14 02:00:17 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 : Thu Jun 13 11:20:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	4131	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	15397	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	8109	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	15055m	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	11087m	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	9885m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	6588	0.28	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	15480m	0.38	ng/ul	0.00
Target Compounds				Ovalue		
19) Chrysene	21.30	228	1536	0.033	ng/ul#	82
21) Benzo(b)fluoranthene	22.83	252	1579m	0.038	ng/ul	
22) Benzo(k)fluoranthene	22.87	252	1443m	0.036	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.74	276	1661m	0.040	ng/ul	
25) Dibenzo(a,h)anthracene	25.75	278	1224m	0.037	ng/ul	
26) Benzo(g,h,i)perylene	26.42	276	1389m	0.040	ng/ul	

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

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