

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060819\
 Data File : BN006204.D
 Acq On : 08 Jun 2019 23:41
 Operator : JU/SJ
 Sample : K3017-02
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A51G3

Manual Integrations
 APPROVED

mohammad
 6/11/2019 4:32:02 PM

Quant Time: Jun 09 06:01:54 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	7.65	152	4712	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	17730	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	9127	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	18943m	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	15653m	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	14772m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	4707	0.17	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	10436m	0.21	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.01	152	4184	0.082	ng/ul#	94
9) Fluorene	15.35	166	832	0.020	ng/ul#	91
12) Phenanthrene	17.09	178	11456	0.188	ng/ul	97
13) Anthracene	17.18	178	3240m	0.056	ng/ul	
15) Fluoranthene	19.12	202	30317m	0.452	ng/ul	
17) Pyrene	19.49	202	29764	0.361	ng/ul	99
18) Benzo(a)anthracene	21.24	228	12051	0.174	ng/ul#	87
19) Chrysene	21.30	228	19267	0.296	ng/ul#	90
21) Benzo(b)fluoranthene	22.82	252	23687m	0.385	ng/ul	
22) Benzo(k)fluoranthene	22.86	252	9519m	0.159	ng/ul	
23) Benzo(a)pyrene	23.39	252	11938	0.208	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.72	276	9957	0.160	ng/ul#	87
25) Dibenzo(a,h)anthracene	25.72	278	2389	0.048	ng/ul#	31
26) Benzo(g,h,i)perylene	26.40	276	8611	0.166	ng/ul#	87

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

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