

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061019\
 Data File : BN006235.D
 Acq On : 10 Jun 2019 22:19
 Operator : JU/SJ
 Sample : K3017-22
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A51F5

Manual Integrations
 APPROVED

mohammad
 6/11/2019 4:29:01 PM

Quant Time: Jun 11 02:23:02 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	4293	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	16283	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	8505	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	17563m	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	13623m	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	17766	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	4908	0.20	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	10935m	0.23	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.01	152	1800	0.038	ng/ul#	86
8) Acenaphthene	14.36	153	706	0.022	ng/ul#	93
12) Phenanthrene	17.09	178	22300	0.396	ng/ul	98
13) Anthracene	17.18	178	3419m	0.064	ng/ul	
15) Fluoranthene	19.12	202	72626m	1.167	ng/ul	
17) Pyrene	19.49	202	57399	0.800	ng/ul	99
18) Benzo(a)anthracene	21.25	228	22813m	0.379	ng/ul	
19) Chrysene	21.30	228	31851	0.563	ng/ul	98
21) Benzo(b)fluoranthene	22.82	252	45387m	0.613	ng/ul	
22) Benzo(k)fluoranthene	22.86	252	17069m	0.237	ng/ul	
23) Benzo(a)pyrene	23.39	252	27187	0.394	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.72	276	23040	0.309	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.73	278	5248	0.088	ng/ul#	56
26) Benzo(g,h,i)perylene	26.41	276	20828	0.333	ng/ul	98

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

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