

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061019\
 Data File : BN006233.D
 Acq On : 10 Jun 2019 21:10
 Operator : JU/SJ
 Sample : K3017-20
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A51F1

Manual Integrations
 APPROVED

mohammad
 6/11/2019 4:28:49 PM

Quant Time: Jun 11 02:11:32 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	4004	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	15188	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	6823	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	13107m	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	12317m	0.40	ng/ul	0.00
20) Perylene-d12	23.50	264	19440	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	4811	0.21	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	9858m	0.28	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.01	152	6210	0.164	ng/ul	97
8) Acenaphthene	14.36	153	1277	0.049	ng/ul	96
9) Fluorene	15.35	166	2055	0.067	ng/ul#	98
12) Phenanthrene	17.09	178	73437m	1.745	ng/ul	
13) Anthracene	17.18	178	10083m	0.251	ng/ul	
15) Fluoranthene	19.12	202	237259m	5.107	ng/ul	
17) Pyrene	19.49	202	186858	2.880	ng/ul	100
18) Benzo(a)anthracene	21.25	228	89181m	1.637	ng/ul	
19) Chrysene	21.30	228	117216	2.292	ng/ul	98
21) Benzo(b)fluoranthene	22.83	252	167568m	2.067	ng/ul	
22) Benzo(k)fluoranthene	22.87	252	55724m	0.706	ng/ul	
23) Benzo(a)pyrene	23.40	252	99217	1.314	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	25.72	276	82117	1.005	ng/ul#	87
25) Dibenzo(a,h)anthracene	25.73	278	18109	0.277	ng/ul#	90
26) Benzo(g,h,i)perylene	26.41	276	71697	1.048	ng/ul	96

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

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