

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060619\
 Data File : BN006157.D
 Acq On : 07 Jun 2019 04:39
 Operator : JU/SJ
 Sample : K3016-07
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A51C6

Manual Integrations
 APPROVED

mohammad
 6/11/2019 5:08:01 PM

Quant Time: Jun 07 06:44:37 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 : Fri Jun 07 01:31:14 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	5929	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	21699	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	10915	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	22335m	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	19953m	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	18033m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	5831	0.18	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	12917m	0.22	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.01	152	4368	0.072	ng/ul#	93
12) Phenanthrene	17.09	178	13415	0.187	ng/ul	98
13) Anthracene	17.18	178	3217m	0.047	ng/ul	
15) Fluoranthene	19.12	202	33954m	0.429	ng/ul	
17) Pyrene	19.49	202	30256	0.288	ng/ul	96
18) Benzo(a)anthracene	21.25	228	12704m	0.144	ng/ul	
19) Chrysene	21.30	228	17190	0.207	ng/ul	95
21) Benzo(b)fluoranthene	22.83	252	20573m	0.274	ng/ul	
22) Benzo(k)fluoranthene	22.87	252	7983m	0.109	ng/ul	
23) Benzo(a)pyrene	23.40	252	13210	0.189	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.72	276	10068	0.133	ng/ul	98
25) Dibenzo(a,h)anthracene	25.73	278	2901	0.048	ng/ul#	33
26) Benzo(g,h,i)perylene	26.41	276	9461	0.149	ng/ul#	90

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

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