

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060619\
 Data File : BN006158.D
 Acq On : 07 Jun 2019 05:14
 Operator : JU/SJ
 Sample : K3016-21
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A51D8

Manual Integrations
 APPROVED

mohammad
 6/11/2019 5:08:04 PM

Quant Time: Jun 07 06:50:12 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	4986	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	18817	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	9620	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	20137m	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	17710m	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	15925m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	4645	0.16	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	10367m	0.19	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.48	128	13410	0.258	ng/ul	99
5) 2-Methylnaphthalene	12.10	142	3841	0.106	ng/ul	99
7) Acenaphthylene	14.02	152	8340	0.156	ng/ul#	87
8) Acenaphthene	14.36	153	10579	0.285	ng/ul	100
9) Fluorene	15.36	166	13278	0.306	ng/ul	99
12) Phenanthrene	17.09	178	172271m	2.665	ng/ul	
13) Anthracene	17.18	178	47407m	0.768	ng/ul	
15) Fluoranthene	19.12	202	261550m	3.664	ng/ul	
17) Pyrene	19.48	202	215947	2.315	ng/ul	98
18) Benzo(a)anthracene	21.25	228	118145m	1.508	ng/ul	
19) Chrysene	21.30	228	116097	1.579	ng/ul	98
21) Benzo(b)fluoranthene	22.83	252	137614m	2.072	ng/ul	
22) Benzo(k)fluoranthene	22.87	252	58309m	0.902	ng/ul	
23) Benzo(a)pyrene	23.40	252	93837	1.517	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	25.73	276	65696	0.981	ng/ul#	88
25) Dibenzo(a,h)anthracene	25.73	278	17271	0.323	ng/ul#	91
26) Benzo(a,h,i)perylene	26.41	276	47382	0.845	ng/ul	97

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

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