

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060819\
 Data File : BN006211.D
 Acq On : 09 Jun 2019 03:41
 Operator : JU/SJ
 Sample : K3017-11
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A51E4

Manual Integrations
 APPROVED

mohammad
 6/11/2019 4:33:10 PM

Quant Time: Jun 09 06:34:57 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	5326	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	20507	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	10621	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	20965m	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	16905m	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	15576m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	6179	0.20	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	13457m	0.24	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.01	152	2295	0.039	ng/ul#	91
12) Phenanthrene	17.09	178	10025	0.149	ng/ul	97
13) Anthracene	17.18	178	1750m	0.027	ng/ul	
15) Fluoranthene	19.12	202	23460m	0.316	ng/ul	
17) Pyrene	19.49	202	20492	0.230	ng/ul	99
18) Benzo(a)anthracene	21.24	228	6878m	0.092	ng/ul	
19) Chrysene	21.29	228	10884	0.155	ng/ul	96
21) Benzo(b)fluoranthene	22.82	252	14407m	0.222	ng/ul	
22) Benzo(k)fluoranthene	22.86	252	6128m	0.097	ng/ul	
23) Benzo(a)pyrene	23.39	252	8107	0.134	ng/ul#	74
24) Indeno(1,2,3-cd)pyrene	25.72	276	6607	0.101	ng/ul#	87
25) Dibenzo(a,h)anthracene	25.72	278	1571	0.030	ng/ul#	29
26) Benzo(g,h,i)perylene	26.40	276	5090	0.093	ng/ul#	85

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

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