

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060819\
 Data File : BN006207.D
 Acq On : 09 Jun 2019 01:24
 Operator : JU/SJ
 Sample : K3017-13
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A51E6

Manual Integrations
 APPROVED

mohammad
 6/11/2019 4:33:16 PM

Quant Time: Jun 10 02:37:15 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	4983	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	19181	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	9976	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	21240m	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	17137m	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	16843m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	5011	0.17	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	11224m	0.20	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.09	178	5620	0.082	ng/ul	94
15) Fluoranthene	19.12	202	20746m	0.276	ng/ul	
17) Pyrene	19.49	202	17088	0.189	ng/ul	97
18) Benzo(a)anthracene	21.24	228	5577	0.074	ng/ul#	87
19) Chrysene	21.30	228	10308	0.145	ng/ul#	92
21) Benzo(b)fluoranthene	22.82	252	14498m	0.206	ng/ul	
22) Benzo(k)fluoranthene	22.86	252	5932m	0.087	ng/ul	
23) Benzo(a)pyrene	23.39	252	7894	0.121	ng/ul#	65
24) Indeno(1,2,3-cd)pyrene	25.72	276	6776	0.096	ng/ul	93
26) Benzo(g,h,i)perylene	26.40	276	6225	0.105	ng/ul#	81

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

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