

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061819\
 Data File : BN006386.D
 Acq On : 18 Jun 2019 20:43
 Operator : JU/SJ
 Sample : K3335-05
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A41Z7

Manual Integrations
 APPROVED

mohammad
 6/21/2019 8:39:25 AM

Quant Time: Jun 19 02:10:25 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 : Mon Jun 17 11:13:07 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	3573	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	13559	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	6586	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	12413	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	9197	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	12710	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	2846	0.14	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	5676	0.17	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.09	178	2329	0.058	ng/ul#	91
15) Fluoranthene	19.12	202	6587	0.150	ng/ul	95
17) Pyrene	19.48	202	5201m	0.107	ng/ul	
18) Benzo(a)anthracene	21.25	228	2515	0.062	ng/ul#	91
19) Chrysene	21.30	228	2698	0.071	ng/ul	93
21) Benzo(b)fluoranthene	22.83	252	4069m	0.077	ng/ul	
22) Benzo(k)fluoranthene	22.87	252	1473m	0.029	ng/ul	
23) Benzo(a)pyrene	23.39	252	2198	0.045	ng/ul#	70
24) Indeno(1,2,3-cd)pyrene	25.73	276	1804	0.034	ng/ul#	85
26) Benzo(g,h,i)perylene	26.42	276	1151	0.026	ng/ul#	61

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

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