

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061719\
 Data File : BN006359.D
 Acq On : 17 Jun 2019 21:35
 Operator : JU/SJ
 Sample : K3335-02
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4225

Manual Integrations
 APPROVED

mohammad
 6/21/2019 8:22:41 AM

Quant Time: Jun 18 04:11:16 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	4016	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	15881	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	8252	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	15176m	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	10355m	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	9267m	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	3740	0.15	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	8112m	0.20	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.09	178	7637	0.157	ng/ul	97
13) Anthracene	17.18	178	1536m	0.033	ng/ul	
15) Fluoranthene	19.12	202	14254m	0.265	ng/ul	
17) Pyrene	19.48	202	12206m	0.224	ng/ul	
18) Benzo(a)anthracene	21.25	228	5718	0.125	ng/ul	95
19) Chrysene	21.30	228	5417	0.126	ng/ul	96
21) Benzo(b)fluoranthene	22.83	252	7693m	0.199	ng/ul	
22) Benzo(k)fluoranthene	22.87	252	2900m	0.077	ng/ul	
23) Benzo(a)pyrene	23.39	252	4777m	0.133	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.73	276	3137m	0.081	ng/ul	
26) Benzo(g,h,i)perylene	26.42	276	2804m	0.086	ng/ul	

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

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