

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061819\
 Data File : BN006375.D
 Acq On : 18 Jun 2019 13:30
 Operator : JU/SJ
 Sample : K3335-06
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4220

Manual Integrations
 APPROVED

mohammad
 6/21/2019 8:38:55 AM

Quant Time: Jun 18 16:41:23 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 : Tue Jun 18 15:57:40 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	4287	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	17242	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	9104	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	16934m	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	10260m	0.40	ng/ul	0.00
20) Perylene-d12	23.50	264	8809m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	6084	0.23	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	11794m	0.26	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.01	152	1069	0.021	ng/ul#	82
12) Phenanthrene	17.09	178	17860	0.329	ng/ul	99
13) Anthracene	17.18	178	2724m	0.052	ng/ul	
15) Fluoranthene	19.12	202	48409m	0.806	ng/ul	
17) Pyrene	19.48	202	39807	0.737	ng/ul	98
18) Benzo(a)anthracene	21.25	228	14807m	0.326	ng/ul	
19) Chrysene	21.30	228	17126	0.402	ng/ul	99
21) Benzo(b)fluoranthene	22.83	252	26731m	0.728	ng/ul	
22) Benzo(k)fluoranthene	22.87	252	11150m	0.312	ng/ul	
23) Benzo(a)pyrene	23.40	252	16243	0.475	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.74	276	11443	0.309	ng/ul#	86
25) Dibenzo(a,h)anthracene	25.74	278	2631	0.089	ng/ul#	84
26) Benzo(g,h,i)perylene	26.42	276	9252	0.298	ng/ul	99

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

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