

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061719\
 Data File : BN006361.D
 Acq On : 17 Jun 2019 22:44
 Operator : JU/SJ
 Sample : K3335-10
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4237

Manual Integrations
 APPROVED

mohammad
 6/21/2019 8:22:47 AM

Quant Time: Jun 18 04:19:56 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	4302	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	17311	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	9246	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	16876m	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	10472m	0.40	ng/ul	-0.01
20) Perylene-d12	23.49	264	15899	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	4703	0.18	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	8661m	0.19	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.01	152	2698	0.052	ng/ul#	92
9) Fluorene	15.35	166	1432	0.034	ng/ul#	96
12) Phenanthrene	17.09	178	23472m	0.433	ng/ul	
13) Anthracene	17.18	178	4792m	0.093	ng/ul	
15) Fluoranthene	19.11	202	40363m	0.675	ng/ul	
17) Pyrene	19.48	202	38531	0.699	ng/ul	98
18) Benzo(a)anthracene	21.24	228	18092m	0.391	ng/ul	
19) Chrysene	21.29	228	15050	0.346	ng/ul	98
21) Benzo(b)fluoranthene	22.82	252	20484m	0.309	ng/ul	
22) Benzo(k)fluoranthene	22.86	252	8108m	0.126	ng/ul	
23) Benzo(a)pyrene	23.38	252	13907	0.225	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.72	276	8838	0.132	ng/ul#	86
25) Dibenzo(a,h)anthracene	25.73	278	2547	0.048	ng/ul#	76
26) Benzo(g,h,i)perylene	26.40	276	7017	0.125	ng/ul	98

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

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