

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061719\
 Data File : BN006364.D
 Acq On : 18 Jun 2019 00:28
 Operator : JU/SJ
 Sample : K3335-21
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A41X1

Manual Integrations
 APPROVED

mohammad
 6/21/2019 8:22:55 AM

Quant Time: Jun 18 04:36:17 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	4641	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	18679	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	9773	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	17946m	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	11010m	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	16980	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	4073	0.14	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	7128m	0.15	ng/ul	0.00
Target Compounds				Ovalue		
9) Fluorene	15.35	166	935	0.021	ng/ul#	96
12) Phenanthrene	17.09	178	18270	0.317	ng/ul	98
13) Anthracene	17.18	178	4713m	0.086	ng/ul	
15) Fluoranthene	19.11	202	27985m	0.440	ng/ul	
17) Pyrene	19.48	202	21911	0.378	ng/ul	99
18) Benzo(a)anthracene	21.24	228	9521m	0.196	ng/ul	
19) Chrysene	21.30	228	9847	0.215	ng/ul	99
21) Benzo(b)fluoranthene	22.82	252	12076m	0.171	ng/ul	
22) Benzo(k)fluoranthene	22.86	252	3631m	0.053	ng/ul	
23) Benzo(a)pyrene	23.39	252	7310	0.111	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.73	276	6288	0.088	ng/ul#	91
25) Dibenzo(a,h)anthracene	25.73	278	2316	0.041	ng/ul#	85
26) Benzo(g,h,i)perylene	26.41	276	3030	0.051	ng/ul	92

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

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