

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102219\
 Data File : BN008374.D
 Acq On : 24 Oct 2019 05:52
 Operator : JU
 Sample : K5235-13
 Misc :
 ALS Vial : 66 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPL1

Manual Integrations
 APPROVED

mohammad
 10/24/2019 8:07:34 AM

Quant Time: Oct 24 06:43:11 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 23 14:07:55 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	1669	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.34	136	8255	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.21	164	5169	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.96	188	11307m	0.40	ng/ul	-0.01
16) Chrysene-d12	21.17	240	10042	0.40	ng/ul	-0.01
20) Perylene-d12	23.33	264	9163	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.94	152	3012	0.22	ng/ul	-0.02
14) Fluoranthene-d10	19.00	212	8832	0.23	ng/ul	-0.01
Target Compounds				Ovalue		
12) Phenanthrene	17.00	178	1099	0.033	ng/ul#	85
15) Fluoranthene	19.03	202	3127	0.072	ng/ul	91
17) Pyrene	19.39	202	2892	0.072	ng/ul#	85
18) Benzo(a)anthracene	21.15	228	1383	0.038	ng/ul#	87
19) Chrysene	21.20	228	1548	0.035	ng/ul#	90
21) Benzo(b)fluoranthene	22.70	252	1793m	0.060	ng/ul	
23) Benzo(a)pyrene	23.24	252	1124	0.036	ng/ul#	5

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

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