

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101819\
 Data File : BN008258.D
 Acq On : 18 Oct 2019 21:34
 Operator : JU
 Sample : K5177-16
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPK2

Manual Integrations
 APPROVED

mohammad
 10/22/2019 1:41:05 AM

Quant Time: Oct 19 00:58:10 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 : Fri Oct 18 23:47:17 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	2446	0.40	ng/ul	0.00
2) Naphthalene-d8	10.36	136	11529	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.23	164	6966	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.97	188	14540	0.40	ng/ul	0.00
16) Chrysene-d12	21.18	240	13478	0.40	ng/ul	0.00
20) Perylene-d12	23.36	264	12243	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	3633	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.01	212	9565	0.19	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.02	178	1372	0.032	ng/ul#	88
15) Fluoranthene	19.04	202	2951	0.052	ng/ul	89
17) Pyrene	19.40	202	3254m	0.061	ng/ul	
18) Benzo(a)anthracene	21.17	228	1521	0.031	ng/ul#	84
19) Chrysene	21.22	228	2097	0.035	ng/ul#	88
21) Benzo(b)fluoranthene	22.72	252	2340m	0.059	ng/ul	
23) Benzo(a)pyrene	23.26	252	1409	0.034	ng/ul#	1

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

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