

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101119\
 Data File : BN008107.D
 Acq On : 12 Oct 2019 11:10
 Operator : JU
 Sample : K5124-18
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 JLMA7

Manual Integrations
 APPROVED

mohammad
 10/15/2019 7:58:28 AM

Quant Time: Oct 13 02:48:14 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 11 19:05:54 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	2801	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	13438	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.26	164	8712	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	19952m	0.40	ng/ul	0.00
16) Chrysene-d12	21.21	240	20219m	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	21010	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	3953	0.17	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	13083m	0.19	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.44	128	1058	0.028	ng/ul#	74
5) 2-Methylnaphthalene	12.07	142	1045	0.040	ng/ul	98
7) Acenaphthylene	13.98	152	2589	0.062	ng/ul#	84
12) Phenanthrene	17.05	178	2817m	0.048	ng/ul	
15) Fluoranthene	19.07	202	4049m	0.052	ng/ul	
17) Pyrene	19.44	202	4074	0.051	ng/ul#	80
18) Benzo(a)anthracene	21.20	228	1779	0.024	ng/ul#	1
19) Chrysene	21.25	228	2834	0.031	ng/ul#	35
21) Benzo(b)fluoranthene	22.76	252	3863m	0.056	ng/ul	
23) Benzo(a)pyrene	23.31	252	2029	0.028	ng/ul#	1

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

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