

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101218\
 Data File : BN003136.D
 Acq On : 12 Oct 2018 10:41
 Operator : MJ/SJ
 Sample : J5234-12
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AL0

Quant Time: Oct 12 13:28:29 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN100918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 12 07:27:48 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1910	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.39	136	7776	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.26	164	4628	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.11	188	378890	0.40	ng/ul	0.08
16) Chrysene-d12	21.31	240	174	0.40	ng/ul	0.07
20) Perylene-d12	0.00	264	0	0.00	ng/ul	-23.44
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	2964	0.25	ng/ul	-0.03
14) Fluoranthene-d10	19.16	212	2920	0.00	ng/ul	0.09
Target Compounds				Qvalue		
3) Naphthalene	10.44	128	4517	0.226	ng/ul#	96
5) 2-Methylnaphthalene	12.06	142	13066	0.951	ng/ul	98
7) Acenaphthylene	13.98	152	1875	0.089	ng/ul#	44
8) Acenaphthene	14.32	153	3207	0.200	ng/ul	97
9) Fluorene	15.31	166	14949	0.812	ng/ul#	98
12) Phenanthrene	17.05	178	28530	0.027	ng/ul	96
17) Pyrene	19.44	202	21839	32.315	ng/ul	98
18) Benzo(a)anthracene	21.26	228	7037	12.568	ng/ul#	77
19) Chrysene	21.38	228	699	1.229	ng/ul#	1

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

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