

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101218\
 Data File : BN003132.D
 Acq On : 12 Oct 2018 08:20
 Operator : MJ/SJ
 Sample : J5234-17
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AK0

Quant Time: Oct 12 12:45:51 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	1840	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.39	136	6774	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.26	164	3994	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.11	188	290851	0.40	ng/ul	0.08
16) Chrysene-d12	21.31	240	750	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	2758	0.26	ng/ul	-0.03
14) Fluoranthene-d10	19.15	212	5046	0.01	ng/ul	0.09
Target Compounds				Qvalue		
3) Naphthalene	10.43	128	25940	1.492	ng/ul	99
5) 2-Methylnaphthalene	12.06	142	45710	3.820	ng/ul	97
7) Acenaphthylene	13.98	152	3758	0.206	ng/ul#	44
8) Acenaphthene	14.32	153	6735	0.486	ng/ul	93
9) Fluorene	15.31	166	29636	1.866	ng/ul#	99
12) Phenanthrene	17.05	178	55394	0.068	ng/ul	98
17) Pyrene	19.44	202	30021	10.306	ng/ul	99
18) Benzo(a)anthracene	21.26	228	8645	3.582	ng/ul#	68
19) Chrysene	21.34	228	292	0.119	ng/ul#	1

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

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