

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100918\
 Data File : BN003026.D
 Acq On : 09 Oct 2018 14:50
 Operator : MJ/SJ
 Sample : J5116-08DL 5X
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 JLC57DL

Quant Time: Oct 09 15:41:51 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 : Tue Oct 09 08:56:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1169	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.41	136	5014	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.26	164	2913	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	89315	0.40	ng/ul	0.08
16) Chrysene-d12	0.00	240	0	0.00	ng/ul	-21.24
20) Perylene-d12	23.47	264	5479	0.40	ng/ul	0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	529	0.07	ng/ul	0.00
14) Fluoranthene-d10	0.00	212	0	0.00	ng/ul	
Target Compounds						
					Qvalue	
3) Naphthalene	10.46	128	1330	0.103	ng/ul#	87
5) 2-Methylnaphthalene	12.09	142	424	0.048	ng/ul	94
7) Acenaphthylene	13.99	152	504	0.038	ng/ul#	64
8) Acenaphthene	14.33	153	5598	0.554	ng/ul	99
9) Fluorene	15.32	166	5479	0.473	ng/ul	100
12) Phenanthrene	17.15	178	27825	0.112	ng/ul	98
21) Benzo(b)fluoranthene	22.81	252	137863	6.987	ng/ul#	1
22) Benzo(k)fluoranthene	22.81	252	137863	6.740	ng/ul#	1
23) Benzo(a)pyrene	23.37	252	41510	2.347	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	25.70	276	44138	2.172	ng/ul	97
25) Dibenzo(a,h)anthracene	25.70	278	9309	0.572	ng/ul#	1
26) Benzo(g,h,i)perylene	26.38	276	37935	2.214	ng/ul#	62

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

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