

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN100918\
 Data File : BN003001.D
 Acq On : 08 Oct 2018 23:27
 Operator : MJ/SJ
 Sample : J5116-02
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 JLC53

Manual Integrations
 APPROVED

Sohil
 10/9/2018 5:37:55 PM

Quant Time: Oct 09 01:56:24 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 01:14:49 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1416	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.39	136	5893	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.26	164	3430	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	7739m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.23	240	7753m	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	7187	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	2669	0.29	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	6712m	0.30	ng/ul	0.00
Target Compounds				Qvalue		
12) Phenanthrene	17.06	178	1364m	0.063	ng/ul	
13) Anthracene	17.15	178	426m	0.021	ng/ul	
15) Fluoranthene	19.09	202	2796m	0.110	ng/ul	
17) Pyrene	19.45	202	2623m	0.087	ng/ul	
18) Benzo(a)anthracene	21.22	228	1469	0.059	ng/ul#	89
19) Chrysene	21.27	228	2242	0.088	ng/ul	94
21) Benzo(b)fluoranthene	22.78	252	3071m	0.119	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	1022m	0.038	ng/ul	
23) Benzo(a)pyrene	23.34	252	1287	0.055	ng/ul#	53
24) Indeno(1,2,3-cd)pyrene	25.69	276	1083	0.041	ng/ul#	100
26) Benzo(g,h,i)perylene	26.36	276	1093	0.049	ng/ul#	69

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

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