

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN100918\
 Data File : BN003004.D
 Acq On : 09 Oct 2018 01:12
 Operator : MJ/SJ
 Sample : J5115-06
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 JLC15

Manual Integrations
 APPROVED

Sohil
 10/9/2018 5:37:58 PM

Quant Time: Oct 09 02:05:49 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	1367	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.40	136	5795	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.26	164	3393	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	7965m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.23	240	7813m	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	6999	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	1761	0.20	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	4767m	0.21	ng/ul	0.00
Target Compounds						
5) 2-Methylnaphthalene	12.11	142	210	0.021	ng/ul	Qvalue 93
12) Phenanthrene	17.06	178	1346m	0.061	ng/ul	
15) Fluoranthene	19.09	202	1944m	0.075	ng/ul	
17) Pyrene	19.45	202	2211m	0.073	ng/ul	
18) Benzo(a)anthracene	21.22	228	1181m	0.047	ng/ul	
19) Chrysene	21.27	228	3453	0.135	ng/ul	96
21) Benzo(b)fluoranthene	22.78	252	3176m	0.126	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	840m	0.032	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.68	276	1004	0.039	ng/ul	99
26) Benzo(g,h,i)perylene	26.37	276	1044	0.048	ng/ul#	70

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

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