

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN100918\
 Data File : BN003009.D
 Acq On : 09 Oct 2018 04:07
 Operator : MJ/SJ
 Sample : J5115-10
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 JLC41

Manual Integrations
 APPROVED

Sohil
 10/9/2018 5:38:03 PM

Quant Time: Oct 09 08:54:40 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 08:37:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1226	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.40	136	5110	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.26	164	2968	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	7004m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.23	240	7279m	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	7299	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	1570	0.20	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	4427m	0.22	ng/ul	0.00
Target Compounds				Qvalue		
18) Benzo(a)anthracene	21.22	228	766	0.033	ng/ul#	73
19) Chrysene	21.27	228	1675	0.070	ng/ul#	89
21) Benzo(b)fluoranthene	22.78	252	3330m	0.127	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	1137m	0.042	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.68	276	1286	0.048	ng/ul#	97
26) Benzo(g,h,i)perylene	26.35	276	1326	0.058	ng/ul#	76

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

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