

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN100918\
 Data File : BN003002.D
 Acq On : 09 Oct 2018 00:02
 Operator : MJ/SJ
 Sample : J5115-02
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 JLC11

Manual Integrations
 APPROVED

Sohil
 10/9/2018 5:37:56 PM

Quant Time: Oct 09 01:58:36 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	7.64	152	1417	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.40	136	5956	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.26	164	3529	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	8083	0.40	ng/ul	-0.02
16) Chrysene-d12	21.23	240	8098	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	7119	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	1317	0.14	ng/ul	0.01
14) Fluoranthene-d10	19.06	212	3525	0.15	ng/ul	0.00
Target Compounds				Qvalue		
12) Phenanthrene	17.06	178	2363	0.105	ng/ul	97
13) Anthracene	17.15	178	430	0.020	ng/ul#	59
15) Fluoranthene	19.09	202	4445	0.168	ng/ul	96
17) Pyrene	19.45	202	4776	0.152	ng/ul	95
18) Benzo(a)anthracene	21.22	228	1735	0.067	ng/ul#	91
19) Chrysene	21.27	228	2891	0.109	ng/ul	94
21) Benzo(b)fluoranthene	22.79	252	3639m	0.142	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	1210m	0.046	ng/ul	
23) Benzo(a)pyrene	23.34	252	1805	0.079	ng/ul#	69
24) Indeno(1,2,3-cd)pyrene	25.68	276	1375	0.052	ng/ul	96
26) Benzo(g,h,i)perylene	26.36	276	1328	0.060	ng/ul#	73

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

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