

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN101018\
 Data File : BN003049.D
 Acq On : 10 Oct 2018 04:20
 Operator : MJ/SJ
 Sample : J5115-17
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 JLC60

Manual Integrations
 APPROVED

Sohil
 10/10/2018 4:58:19 PM

Quant Time: Oct 10 09:02:30 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 : Wed Oct 10 08:36:11 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1832	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.39	136	7622	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.26	164	4404	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	9564m	0.40	ng/ul	-0.01
16) Chrysene-d12	21.23	240	6812m	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	5671	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	3000	0.25	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	7188m	0.26	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.44	128	1169	0.060	ng/ul#	87
5) 2-Methylnaphthalene	12.08	142	531	0.039	ng/ul	97
7) Acenaphthylene	13.98	152	583	0.029	ng/ul#	65
12) Phenanthrene	17.06	178	4817m	0.181	ng/ul	
13) Anthracene	17.15	178	743m	0.030	ng/ul	
15) Fluoranthene	19.09	202	8367m	0.268	ng/ul	
17) Pyrene	19.45	202	8793m	0.332	ng/ul	
18) Benzo(a)anthracene	21.21	228	3698	0.169	ng/ul#	80
19) Chrysene	21.26	228	4475	0.201	ng/ul#	84
21) Benzo(b)fluoranthene	22.78	252	5455m	0.267	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	1528m	0.072	ng/ul	
23) Benzo(a)pyrene	23.33	252	2864	0.156	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	25.65	276	2697	0.128	ng/ul#	94
26) Benzo(g,h,i)perylene	26.33	276	3299	0.186	ng/ul#	70

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

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