

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN101018\
 Data File : BN003031.D
 Acq On : 09 Oct 2018 17:46
 Operator : MJ/SJ
 Sample : J5157-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AA0

Manual Integrations
 APPROVED

Sohil
 10/10/2018 4:57:38 PM

Quant Time: Oct 09 18:47:10 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 08:56:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1462	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.39	136	6261	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.26	164	3847	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	9017m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.23	240	8660m	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	7825	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	2884	0.30	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	12631m	0.49	ng/ul	0.00
Target Compounds				Qvalue		
7) Acenaphthylene	13.98	152	675	0.038	ng/ul#	79
12) Phenanthrene	17.06	178	564m	0.022	ng/ul	
13) Anthracene	17.15	178	848m	0.036	ng/ul	
15) Fluoranthene	19.09	202	2367m	0.080	ng/ul	
17) Pyrene	19.45	202	3320m	0.099	ng/ul	
18) Benzo(a)anthracene	21.22	228	2449m	0.088	ng/ul	
19) Chrysene	21.27	228	3949	0.140	ng/ul	95
21) Benzo(b)fluoranthene	22.78	252	5583m	0.198	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	2132m	0.073	ng/ul	
23) Benzo(a)pyrene	23.34	252	3124	0.124	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.68	276	1664	0.057	ng/ul#	97
26) Benzo(g,h,i)perylene	26.36	276	1539	0.063	ng/ul#	70

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

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