

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN101018\
 Data File : BN003032.D
 Acq On : 09 Oct 2018 18:21
 Operator : MJ/SJ
 Sample : J5157-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AC4

Manual Integrations
 APPROVED

Sohil
 10/10/2018 4:57:39 PM

Quant Time: Oct 09 19:01:41 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	1514	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.39	136	6418	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.26	164	3867	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	8981m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.23	240	8487m	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	7203	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	3691	0.37	ng/ul	-0.01
14) Fluoranthene-d10	19.06	212	9798m	0.38	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.44	128	736	0.045	ng/ul#	80
7) Acenaphthylene	13.98	152	1219	0.069	ng/ul#	87
12) Phenanthrene	17.06	178	933m	0.037	ng/ul	
13) Anthracene	17.15	178	1698m	0.072	ng/ul	
15) Fluoranthene	19.09	202	3906m	0.133	ng/ul	
17) Pyrene	19.45	202	6044m	0.183	ng/ul	
18) Benzo(a)anthracene	21.22	228	3944m	0.144	ng/ul	
19) Chrysene	21.27	228	7336	0.265	ng/ul	97
21) Benzo(b)fluoranthene	22.78	252	9847m	0.380	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	3863m	0.144	ng/ul	
23) Benzo(a)pyrene	23.34	252	5836	0.251	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.67	276	3216	0.120	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.67	278	1029	0.048	ng/ul#	36
26) Benzo(g,h,i)perylene	26.35	276	2898	0.129	ng/ul#	88

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

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