

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN101018\
 Data File : BN003035.D
 Acq On : 09 Oct 2018 20:06
 Operator : MJ/SJ
 Sample : J5157-05
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AC7

Manual Integrations
 APPROVED

Sohil
 10/10/2018 4:57:46 PM

Quant Time: Oct 10 01:56:10 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 : Wed Oct 10 01:34:31 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1438	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.39	136	6166	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.26	164	3745	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	8555m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.23	240	8490m	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	7599	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	3244	0.34	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	8849m	0.36	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.44	128	465	0.029	ng/ul#	65
7) Acenaphthylene	13.98	152	849	0.050	ng/ul#	84
12) Phenanthrene	17.06	178	1062m	0.045	ng/ul	
13) Anthracene	17.15	178	1548m	0.069	ng/ul	
15) Fluoranthene	19.09	202	2378m	0.085	ng/ul	
17) Pyrene	19.45	202	3224m	0.098	ng/ul	
18) Benzo(a)anthracene	21.22	228	2832m	0.104	ng/ul	
19) Chrysene	21.27	228	6353	0.229	ng/ul	99
21) Benzo(b)fluoranthene	22.78	252	6123m	0.224	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	2320m	0.082	ng/ul	
23) Benzo(a)pyrene	23.34	252	3625	0.148	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	25.68	276	2003	0.071	ng/ul#	89
25) Dibenzo(a,h)anthracene	25.67	278	657	0.029	ng/ul#	19
26) Benzo(g,h,i)perylene	26.36	276	1803	0.076	ng/ul#	80

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

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