

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN101018\
 Data File : BN003034.D
 Acq On : 09 Oct 2018 19:31
 Operator : MJ/SJ
 Sample : J5157-04
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AC6

Manual Integrations
 APPROVED

Sohil
 10/10/2018 4:57:44 PM

Quant Time: Oct 10 01:50:36 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 : 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	1648	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.39	136	6888	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.26	164	4164	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	9341m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.23	240	8408m	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	7316	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	4403	0.41	ng/ul	-0.01
14) Fluoranthene-d10	19.06	212	10790m	0.40	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.44	128	632	0.036	ng/ul#	76
7) Acenaphthylene	13.98	152	1179	0.062	ng/ul#	88
12) Phenanthrene	17.06	178	643m	0.025	ng/ul	
13) Anthracene	17.15	178	1287m	0.052	ng/ul	
15) Fluoranthene	19.09	202	5167m	0.169	ng/ul	
17) Pyrene	19.45	202	9104m	0.279	ng/ul	
18) Benzo(a)anthracene	21.22	228	2680m	0.099	ng/ul	
19) Chrysene	21.27	228	9349	0.340	ng/ul	98
21) Benzo(b)fluoranthene	22.78	252	12952m	0.492	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	5490m	0.201	ng/ul	
23) Benzo(a)pyrene	23.34	252	7802	0.330	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.67	276	4439	0.164	ng/ul#	93
25) Dibenzo(a,h)anthracene	25.66	278	1356	0.062	ng/ul#	56
26) Benzo(g,h,i)perylene	26.34	276	3990	0.174	ng/ul	95

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

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