

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101218\
 Data File : BN003134.D
 Acq On : 12 Oct 2018 09:31
 Operator : MJ/SJ
 Sample : J5234-10
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AK8

Manual Integrations
 APPROVED

Sohil
 10/12/2018 6:38:46 PM

Quant Time: Oct 12 13:17:39 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 : Fri Oct 12 07:27:48 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	1841	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.38	136	7911	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.26	164	5066	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.01	188	8900m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.22	240	7645m	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	5666m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	3085	0.25	ng/ul	-0.03
14) Fluoranthene-d10	19.05	212	8290m	0.33	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.43	128	14171	0.698	ng/ul	99
5) 2-Methylnaphthalene	12.06	142	33554	2.401	ng/ul	97
7) Acenaphthylene	13.98	152	3205	0.138	ng/ul#	48
8) Acenaphthene	14.32	153	5363	0.305	ng/ul	96
9) Fluorene	15.31	166	29392	1.459	ng/ul	99
12) Phenanthrene	17.05	178	58375	2.353	ng/ul	97
13) Anthracene	17.15	178	16018m	0.684	ng/ul	
15) Fluoranthene	19.08	202	32116m	1.103	ng/ul	
17) Pyrene	19.45	202	37256	1.255	ng/ul	99
18) Benzo(a)anthracene	21.21	228	7089m	0.288	ng/ul	
19) Chrysene	21.26	228	11249m	0.450	ng/ul	
21) Benzo(b)fluoranthene	22.78	252	7994m	0.392	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	2411m	0.114	ng/ul	
23) Benzo(a)pyrene	23.34	252	5122m	0.280	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.65	276	3288m	0.156	ng/ul	
26) Benzo(a,h,i)perylene	26.33	276	3118m	0.176	ng/ul	

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

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