

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101218\
 Data File : BN003108.D
 Acq On : 11 Oct 2018 18:16
 Operator : MJ/SJ
 Sample : J5368-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A3Z47

Manual Integrations
 APPROVED

Sohil
 10/12/2018 6:38:17 PM

Quant Time: Oct 11 19:05:07 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 : Thu Oct 11 16:42:09 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1863	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.39	136	7831	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.26	164	4453	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	9136m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.22	240	5396m	0.40	ng/ul	-0.01
20) Perylene-d12	23.44	264	4849m	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	2672	0.22	ng/ul	-0.02
14) Fluoranthene-d10	19.05	212	4901m	0.19	ng/ul	-0.01
Target Compounds					Ovalue	
3) Naphthalene	10.44	128	2169	0.108	ng/ul	95
5) 2-Methylnaphthalene	12.08	142	1278	0.092	ng/ul	99
7) Acenaphthylene	13.98	152	3291	0.162	ng/ul#	35
8) Acenaphthene	14.32	153	4432	0.287	ng/ul	98
9) Fluorene	15.32	166	5149	0.291	ng/ul	95
12) Phenanthrene	17.06	178	127500m	5.006	ng/ul	
13) Anthracene	17.15	178	35245m	1.465	ng/ul	
15) Fluoranthene	19.08	202	223646m	7.485	ng/ul	
17) Pyrene	19.45	202	189753	9.054	ng/ul	96
18) Benzo(a)anthracene	21.21	228	109024m	6.279	ng/ul	
19) Chrysene	21.26	228	109148m	6.190	ng/ul	
21) Benzo(b)fluoranthene	22.78	252	195764m	11.210	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	56133m	3.101	ng/ul	
23) Benzo(a)pyrene	23.34	252	129671m	8.285	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.64	276	101884m	5.665	ng/ul	
25) Dibenzo(a,h)anthracene	25.64	278	28441m	1.976	ng/ul	
26) Benzo(a,h,i)perylene	26.32	276	87063m	5.741	ng/ul	

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

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