

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110218\
 Data File : BN003409.D
 Acq On : 02 Nov 2018 18:40
 Operator : MJ/SJ
 Sample : J5704-08
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A40J9

Manual Integrations
 APPROVED

Sohil
 11/5/2018 4:09:14 PM

Quant Time: Nov 02 23:27:05 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN110118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 02 23:49:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.86	152	10956	0.40	ng/ul	0.00
2) Naphthalene-d8	10.65	136	45553	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	26412	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	59156m	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	45495m	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	36022m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.24	152	20862	0.29	ng/ul	0.00
14) Fluoranthene-d10	19.26	212	48516m	0.29	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.28	178	33346	0.172	ng/ul	99
13) Anthracene	17.36	178	6888m	0.044	ng/ul	
15) Fluoranthene	19.28	202	66935m	0.307	ng/ul	
17) Pyrene	19.65	202	53337m	0.322	ng/ul	
18) Benzo(a)anthracene	21.39	228	22854m	0.160	ng/ul	
19) Chrysene	21.44	228	24479m	0.151	ng/ul	
21) Benzo(b)fluoranthene	23.02	252	28238m	0.182	ng/ul	
22) Benzo(k)fluoranthene	23.07	252	9635m	0.064	ng/ul	
23) Benzo(a)pyrene	23.62	252	17544m	0.135	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.08	276	12237m	0.088	ng/ul	
25) Dibenzo(a,h)anthracene	26.08	278	3239m	0.029	ng/ul	
26) Benzo(g,h,i)perylene	26.81	276	9140m	0.077	ng/ul	

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

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