

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110118\
 Data File : BN003394.D
 Acq On : 02 Nov 2018 00:16
 Operator : MJ/SJ
 Sample : J5704-22
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A40J0

Manual Integrations
 APPROVED

Sohil
 11/2/2018 4:14:22 PM

Quant Time: Nov 02 06:07:17 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 04:49:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.86	152	10581	0.40	ng/ul	0.00
2) Naphthalene-d8	10.65	136	45699	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	28274	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	65175m	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	58000m	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	46221m	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.24	152	21559	0.29	ng/ul	0.00
14) Fluoranthene-d10	19.25	212	56917m	0.31	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.21	152	6099	0.050	ng/ul	95
12) Phenanthrene	17.28	178	48394m	0.226	ng/ul	
13) Anthracene	17.36	178	11407m	0.065	ng/ul	
15) Fluoranthene	19.29	202	124816m	0.520	ng/ul	
17) Pyrene	19.65	202	115248m	0.546	ng/ul	
18) Benzo(a)anthracene	21.39	228	58257m	0.320	ng/ul	
19) Chrysene	21.45	228	55949m	0.271	ng/ul	
21) Benzo(b)fluoranthene	23.02	252	69443m	0.349	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	26169m	0.136	ng/ul	
23) Benzo(a)pyrene	23.62	252	48964m	0.294	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.07	276	31091m	0.175	ng/ul	
25) Dibenzo(a,h)anthracene	26.09	278	7634m	0.053	ng/ul	
26) Benzo(g,h,i)perylene	26.80	276	23291m	0.153	ng/ul	

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

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