

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN110518\
 Data File : BN003424.D
 Acq On : 05 Nov 2018 11:56
 Operator : MJ/SJ
 Sample : J5704-04
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A40G1

Manual Integrations
 APPROVED

Sohil
 11/6/2018 1:23:45 PM

Quant Time: Nov 05 14:05:54 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 : Mon Nov 05 13:38:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.86	152	8793	0.40	ng/ul	0.00
2) Naphthalene-d8	10.65	136	36926	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	22168	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	47781m	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	35330m	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	28850m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.24	152	20531	0.35	ng/ul	0.00
14) Fluoranthene-d10	19.25	212	51334m	0.38	ng/ul	0.00
Target Compounds				Qvalue		
12) Phenanthrene	17.27	178	3475	0.022	ng/ul	93
15) Fluoranthene	19.29	202	10032m	0.057	ng/ul	
17) Pyrene	19.65	202	9472	0.074	ng/ul#	95
18) Benzo(a)anthracene	21.39	228	4345m	0.039	ng/ul	
19) Chrysene	21.44	228	4583m	0.036	ng/ul	
21) Benzo(b)fluoranthene	23.02	252	5395m	0.043	ng/ul	
23) Benzo(a)pyrene	23.62	252	3443m	0.033	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.07	276	2547m	0.023	ng/ul	
26) Benzo(g,h,i)perylene	26.80	276	2320m	0.024	ng/ul	

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

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