

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110118\
 Data File : BN003392.D
 Acq On : 01 Nov 2018 23:06
 Operator : MJ/SJ
 Sample : J5704-18
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A40H6

Manual Integrations
 APPROVED

Sohil
 11/2/2018 4:14:17 PM

Quant Time: Nov 02 05:44:52 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	9447	0.40	ng/ul	0.00
2) Naphthalene-d8	10.65	136	41351	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	27307	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	74362m	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	55488m	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	45463m	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.24	152	22403	0.34	ng/ul	0.00
14) Fluoranthene-d10	19.26	212	61169m	0.29	ng/ul	0.00
Target Compounds				Ovalue		
9) Fluorene	15.54	166	41517	0.333	ng/ul#	67
12) Phenanthrene	17.27	178	23594	0.097	ng/ul	97
13) Anthracene	17.36	178	5266m	0.026	ng/ul	
15) Fluoranthene	19.29	202	46403m	0.169	ng/ul	
17) Pyrene	19.65	202	45819m	0.227	ng/ul	
18) Benzo(a)anthracene	21.39	228	21122m	0.121	ng/ul	
19) Chrysene	21.45	228	19802m	0.100	ng/ul	
21) Benzo(b)fluoranthene	23.02	252	23380m	0.119	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	8989m	0.047	ng/ul	
23) Benzo(a)pyrene	23.63	252	16597m	0.101	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.08	276	10322m	0.059	ng/ul	
26) Benzo(g,h,i)perylene	26.80	276	7840m	0.052	ng/ul	

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

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