

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110118\
 Data File : BN003384.D
 Acq On : 01 Nov 2018 18:24
 Operator : MJ/SJ
 Sample : J5657-16
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A40F1

Manual Integrations
 APPROVED

Sohil
 11/2/2018 4:14:03 PM

Quant Time: Nov 02 05:05:58 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	11273	0.40	ng/ul	0.00
2) Naphthalene-d8	10.65	136	46719	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	28163	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	65016m	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	60671m	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	46723m	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.24	152	24909	0.33	ng/ul	0.00
14) Fluoranthene-d10	19.26	212	64804m	0.35	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.28	178	6063m	0.028	ng/ul	
15) Fluoranthene	19.29	202	19088m	0.080	ng/ul	
17) Pyrene	19.65	202	17254	0.078	ng/ul#	93
18) Benzo(a)anthracene	21.39	228	7439m	0.039	ng/ul	
19) Chrysene	21.45	228	13067m	0.060	ng/ul	
21) Benzo(b)fluoranthene	23.02	252	15092m	0.075	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	4152m	0.021	ng/ul	
23) Benzo(a)pyrene	23.62	252	6859m	0.041	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.08	276	5405m	0.030	ng/ul	
26) Benzo(g,h,i)perylene	26.80	276	5020m	0.033	ng/ul	

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

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