

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110118\
 Data File : BN003387.D
 Acq On : 01 Nov 2018 20:10
 Operator : MJ/SJ
 Sample : J5635-27
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A40H7

Manual Integrations
 APPROVED

Sohil
 11/2/2018 4:14:09 PM

Quant Time: Nov 02 05:15:58 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	10978	0.40	ng/ul	0.00
2) Naphthalene-d8	10.65	136	45774	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	27459	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	64215m	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	61530m	0.40	ng/ul	0.00
20) Perylene-d12	23.73	264	49310m	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.24	152	22627	0.31	ng/ul	0.00
14) Fluoranthene-d10	19.26	212	59902m	0.33	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.21	152	3121	0.026	ng/ul#	91
9) Fluorene	15.54	166	3791	0.030	ng/ul#	95
11) Pentachlorophenol	16.88	266	1442m	0.112	ng/ul	
12) Phenanthrene	17.28	178	33119m	0.157	ng/ul	
13) Anthracene	17.36	178	10384m	0.060	ng/ul	
15) Fluoranthene	19.29	202	98968m	0.418	ng/ul	
17) Pyrene	19.65	202	81384m	0.364	ng/ul	
18) Benzo(a)anthracene	21.39	228	45787m	0.237	ng/ul	
19) Chrysene	21.45	228	55322m	0.252	ng/ul	
21) Benzo(b)fluoranthene	23.02	252	69273m	0.326	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	33378m	0.162	ng/ul	
23) Benzo(a)pyrene	23.62	252	41505m	0.234	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.08	276	35659m	0.188	ng/ul	
25) Dibenzo(a,h)anthracene	26.09	278	14881m	0.096	ng/ul	
26) Benzo(g,h,i)perylene	26.80	276	28115m	0.173	ng/ul	

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

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