

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN082318\
 Data File : BN002471.D
 Acq On : 23 Aug 2018 18:18
 Operator : JU/SJ
 Sample : J4421-05
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AA4

Manual Integrations
 APPROVED

Sohil
 8/24/2018 5:11:43 PM

Quant Time: Aug 23 19:04:10 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN082318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 23 14:05:54 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	1522	0.40	ng/ul	0.00
2) Naphthalene-d8	10.45	136	6226	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	3447	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	8233m	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	8818m	0.40	ng/ul	0.00
20) Perylene-d12	23.47	264	8200m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.06	152	1629	0.17	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	4621m	0.19	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.02	152	2510	0.156	ng/ul	98
12) Phenanthrene	17.09	178	4068m	0.167	ng/ul	
13) Anthracene	17.19	178	1265m	0.060	ng/ul	
15) Fluoranthene	19.12	202	12191m	0.426	ng/ul	
17) Pyrene	19.48	202	23041	0.766	ng/ul	96
18) Benzo(a)anthracene	21.24	228	10219m	0.388	ng/ul	
19) Chrysene	21.29	228	12178m	0.408	ng/ul	
21) Benzo(b)fluoranthene	22.82	252	12818m	0.435	ng/ul	
22) Benzo(k)fluoranthene	22.85	252	4710m	0.155	ng/ul	
23) Benzo(a)pyrene	23.37	252	11982m	0.433	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.70	276	5633m	0.195	ng/ul	
25) Dibenzo(a,h)anthracene	25.70	278	1587m	0.070	ng/ul	
26) Benzo(g,h,i)perylene	26.37	276	7727m	0.313	ng/ul	

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

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