

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN082318\
 Data File : BN002482.D
 Acq On : 24 Aug 2018 00:50
 Operator : JU/SJ
 Sample : J4421-09
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AA7

Manual Integrations
 APPROVED

Sohil
 8/24/2018 5:11:53 PM

Quant Time: Aug 24 04:24:42 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 : Fri Aug 24 02:20:09 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	1982	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	8278	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.31	164	4716	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	11209m	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	11246m	0.40	ng/ul	0.00
20) Perylene-d12	23.47	264	9877m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.05	152	2254	0.17	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	6882m	0.21	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.49	128	552	0.026	ng/ul#	70
7) Acenaphthylene	14.02	152	8948	0.405	ng/ul	97
9) Fluorene	15.36	166	1001	0.052	ng/ul#	86
12) Phenanthrene	17.09	178	32397m	0.977	ng/ul	
13) Anthracene	17.19	178	5271m	0.185	ng/ul	
15) Fluoranthene	19.12	202	89948m	2.307	ng/ul	
17) Pyrene	19.48	202	162073	4.225	ng/ul	96
18) Benzo(a)anthracene	21.23	228	51785m	1.540	ng/ul	
19) Chrysene	21.29	228	55314	1.455	ng/ul	97
21) Benzo(b)fluoranthene	22.81	252	45526m	1.283	ng/ul	
22) Benzo(k)fluoranthene	22.85	252	13638m	0.372	ng/ul	
23) Benzo(a)pyrene	23.37	252	44095	1.322	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.69	276	19699	0.567	ng/ul#	79
25) Dibenzo(a,h)anthracene	25.69	278	5244	0.191	ng/ul	96
26) Benzo(g,h,i)perylene	26.36	276	28657	0.962	ng/ul	95

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

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