

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN082318\
 Data File : BN002499.D
 Acq On : 24 Aug 2018 12:58
 Operator : JU/SJ
 Sample : J4424-15RE
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A3Z61RE

Manual Integrations
 APPROVED

Sohil
 8/24/2018 5:38:33 PM

Quant Time: Aug 24 15:37:53 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	1948	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.44	136	10040	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.31	164	7105	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	10455m	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	9855m	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	9093m	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	2581	0.16	ng/ul	-0.02
14) Fluoranthene-d10	19.09	212	3507m	0.11	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.51	128	3556	0.138	ng/ul#	1
7) Acenaphthylene	14.04	152	22582	0.679	ng/ul#	1
8) Acenaphthene	14.37	153	4616m	0.182	ng/ul	
9) Fluorene	15.36	166	23150	0.796	ng/ul#	55
15) Fluoranthene	19.12	202	14509m	0.399	ng/ul	
17) Pyrene	19.48	202	39098m	1.163	ng/ul	
18) Benzo(a)anthracene	21.24	228	3441m	0.117	ng/ul	
19) Chrysene	21.29	228	1936m	0.058	ng/ul	
26) Benzo(g,h,i)perylene	26.40	276	1085m	0.040	ng/ul	

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

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