Quantitation Report (Qedit)

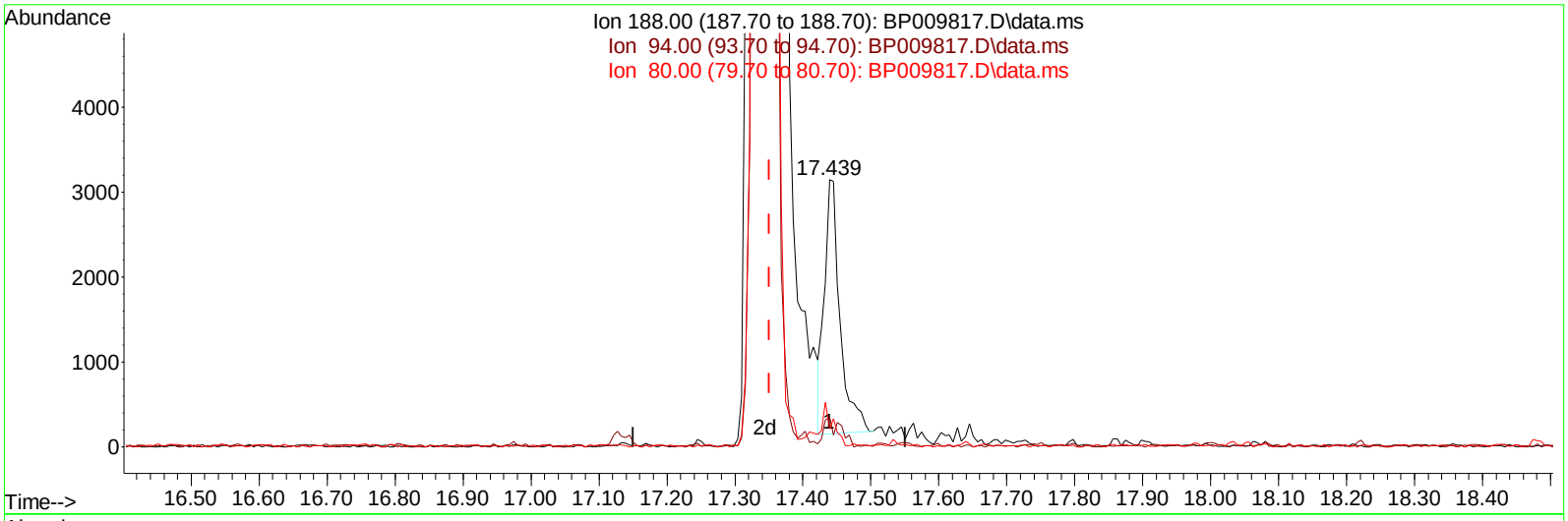
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041322\
Data File : BP009817.D
Acq On : 13 Apr 2022 14:40
Operator : CG/JU
Sample : N2353-05
Misc :
ALS Vial : 10 Sample Multiplier: 1

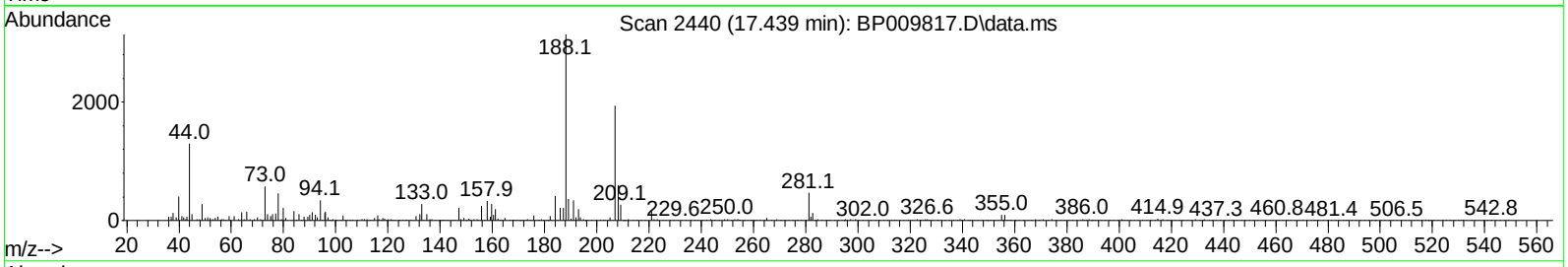
Quant Time: Apr 14 01:06:36 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Apr 13 02:15:22 2022
Response via : Initial Calibration

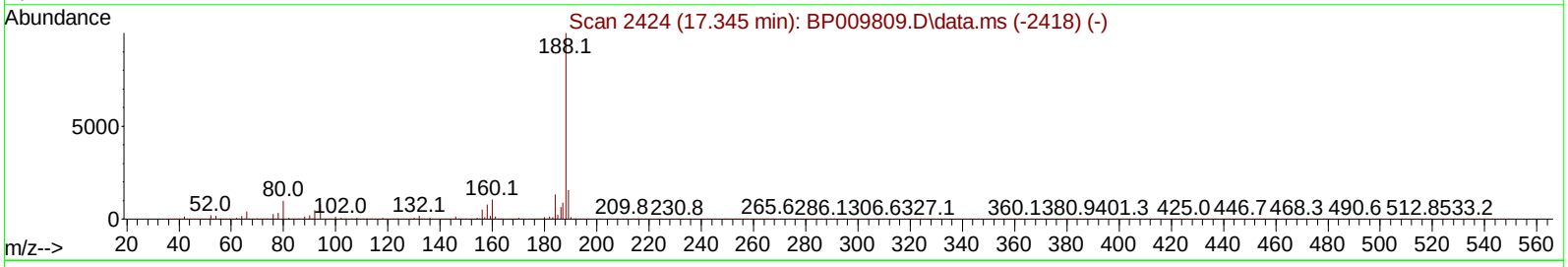
Instrument :
BNA_P
ClientSampleId :
SVOC-GPC-BLANK

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/14/2022
Supervised By :Yogesh Patel 04/14/2022







TIC: BP009817.D\data.ms

(64) Phenanthrene-d10 (I)

17.439min (+ 0.088) 20.00 ng/ul

response 4828

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	9.80	10.91
80.00	6.50	6.84
0.00	0.00	0.00

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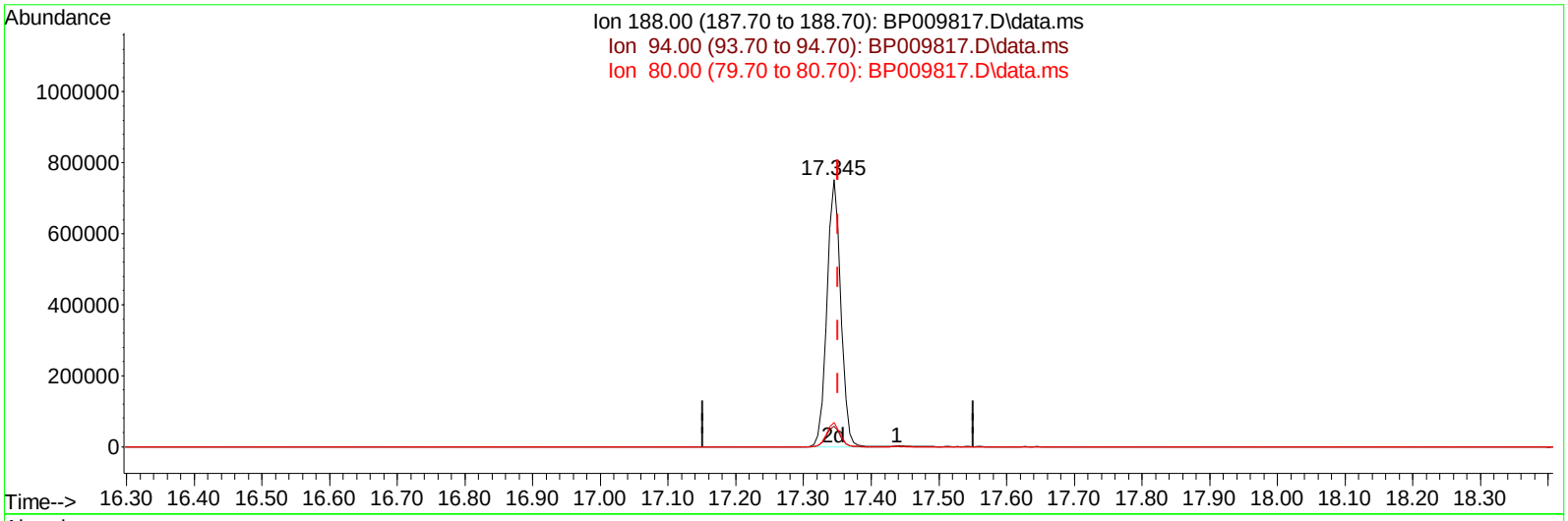

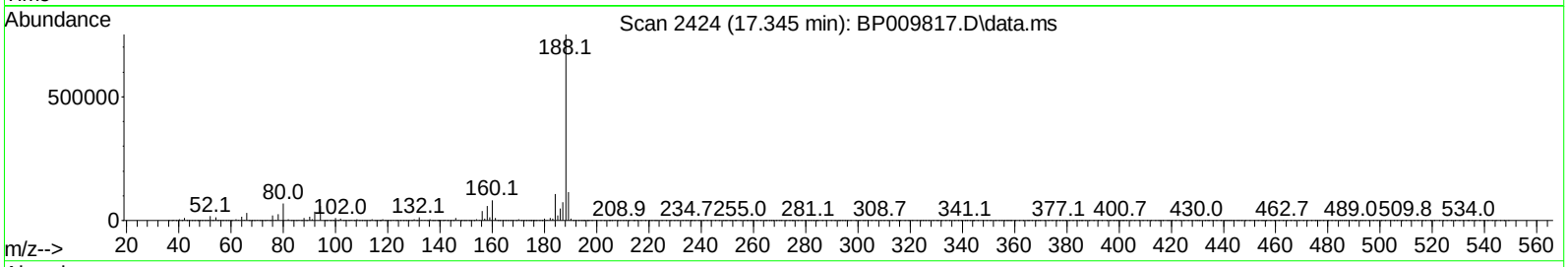

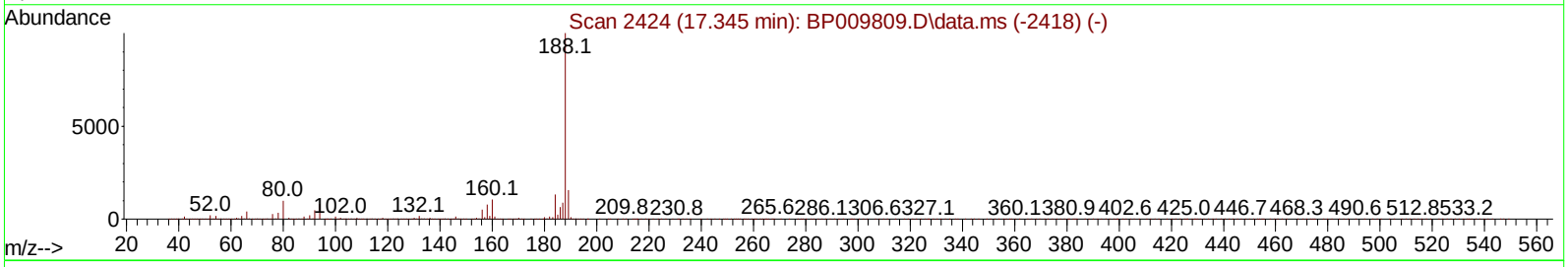

TIC: BP009817.D\data.ms

(64) Phenanthrene-d10 (I)

17.345min (-0.006) 20.00 ng/ul m

response 1063489

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	9.80	7.65#
80.00	6.50	9.10#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.963	152	140269	20.000	ng/ul	0.00
20) Naphthalene-d8	10.769	136	623678	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.592	164	473907	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.345	188	1063489m	20.000	ng/ul	0.00
79) Chrysene-d12	21.421	240	997551	20.000	ng/ul	#-0.01
88) Perylene-d12	23.886	264	845133	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.000	96	0d	0.000	ng/uL	
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	0.000	99	0d	0.000	ng/ul	
9) Bis-(2-Chloroethyl)eth...	0.000	67	0d	0.000	ng/ul	
11) 2-Chlorophenol-d4	0.000	132	0d	0.000	ng/ul	
15) 4-Methylphenol-d8	0.000	113	0d	0.000	ng/ul	
21) Nitrobenzene-d5	0.000	128	0d	0.000	ng/ul	
24) 2-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
28) 2,4-Dichlorophenol-d3	0.000	165	0d	0.000	ng/ul	
31) 4-Chloroaniline-d4	0.000	131	0d	0.000	ng/ul	
46) Dimethylphthalate-d6	0.000	166	0d	0.000	ng/ul	
49) Acenaphthylene-d8	0.000	160	0d	0.000	ng/ul	
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	0.000	176	0d	0.000	ng/ul	
65) 4,6-Dinitro-2-methylph...	0.000	200	0d	0.000	ng/ul	
73) Anthracene-d10	0.000	188	0d	0.000	ng/ul	
81) Pyrene-d10	0.000	212	0d	0.000	ng/ul	
92) Benzo(a)pyrene-d12	0.000	264	0d	0.000	ng/ul	
Target Compounds						Qvalue

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

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