

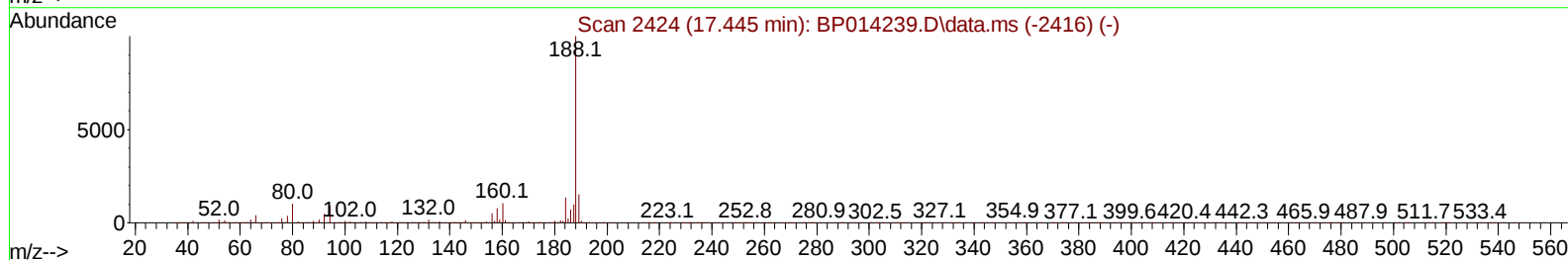
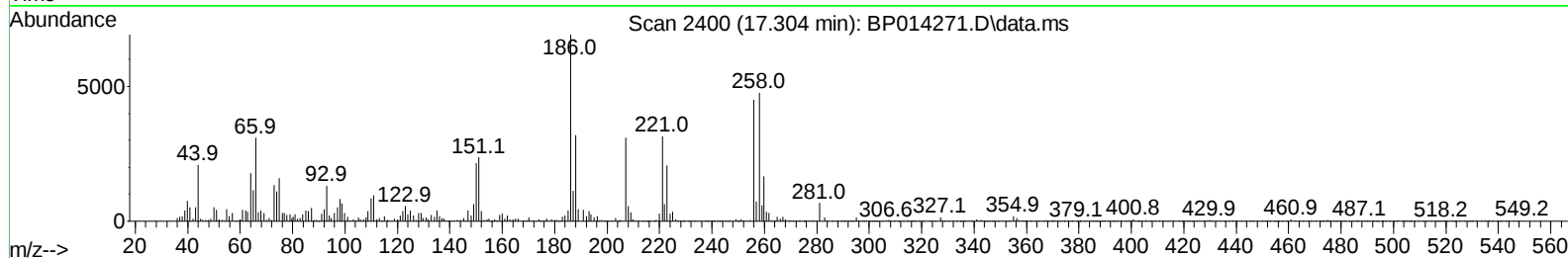
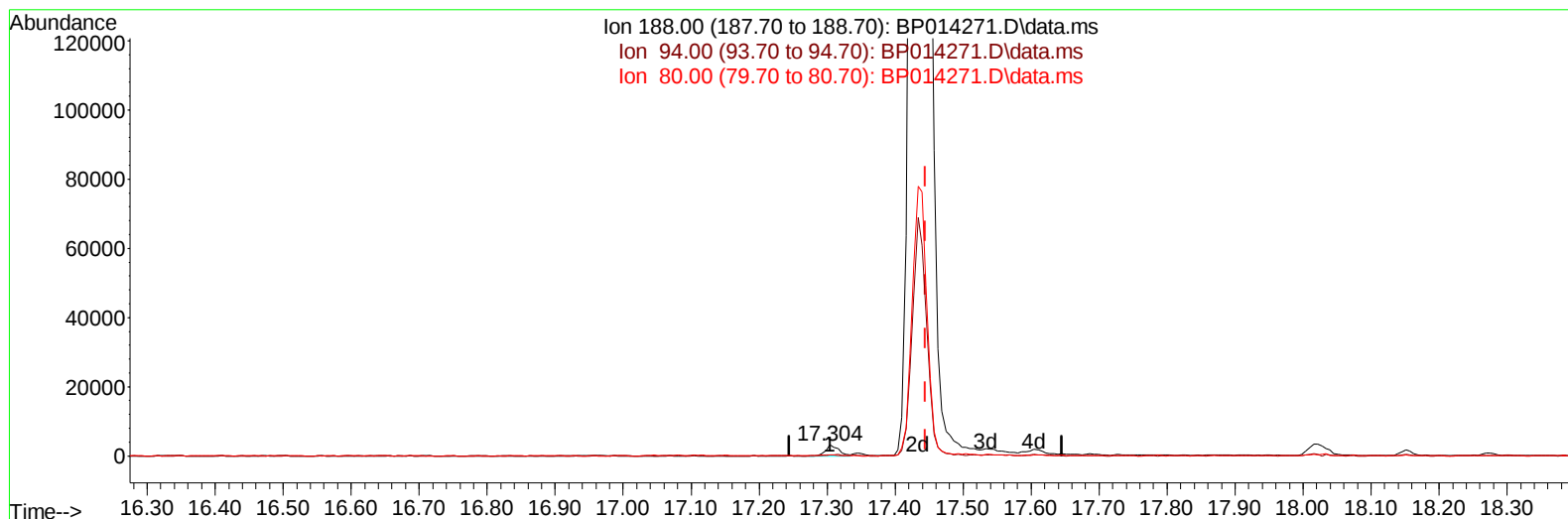
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032223\
Data File : BP014271.D
Acq On : 23 Mar 2023 23:16
Operator : CG/JU
Sample : 01990-13
Misc :
ALS Vial : 51 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E0269

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 03/24/2023
Supervised By :Jagrut Upadhyay 03/24/2023

Quant Time: Mar 24 00:54:42 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Mar 23 04:56:06 2023
Response via : Initial Calibration



TIC: BP014271.D\data.ms

(64) Phenanthrene-d10 (I)

17.304min (-0.141) 20.00 ng/u1

response 4147

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	7.10	6.98
80.00	7.80	9.33
0.00	0.00	0.00

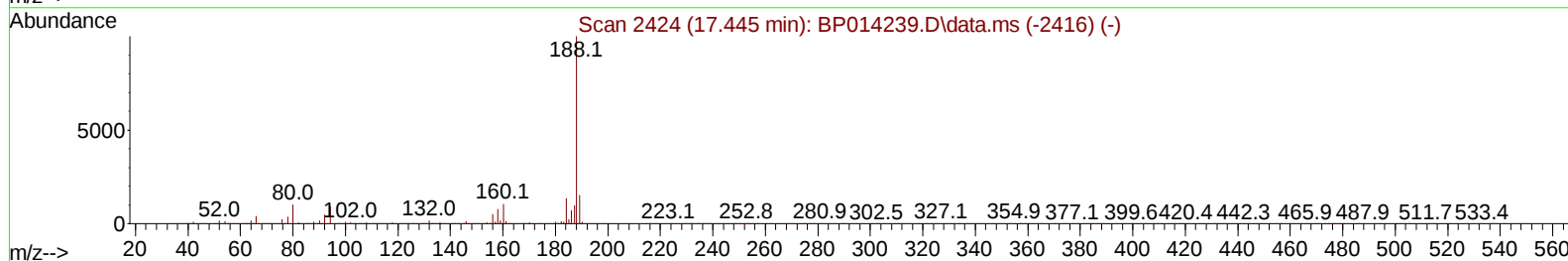
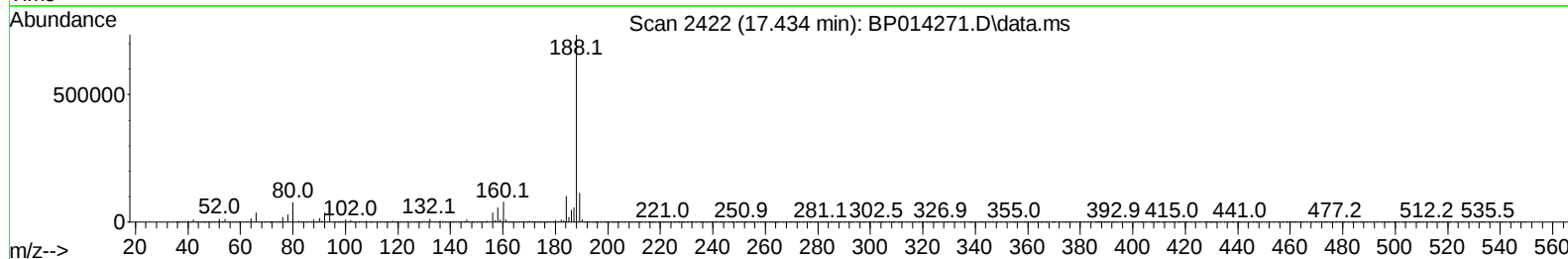
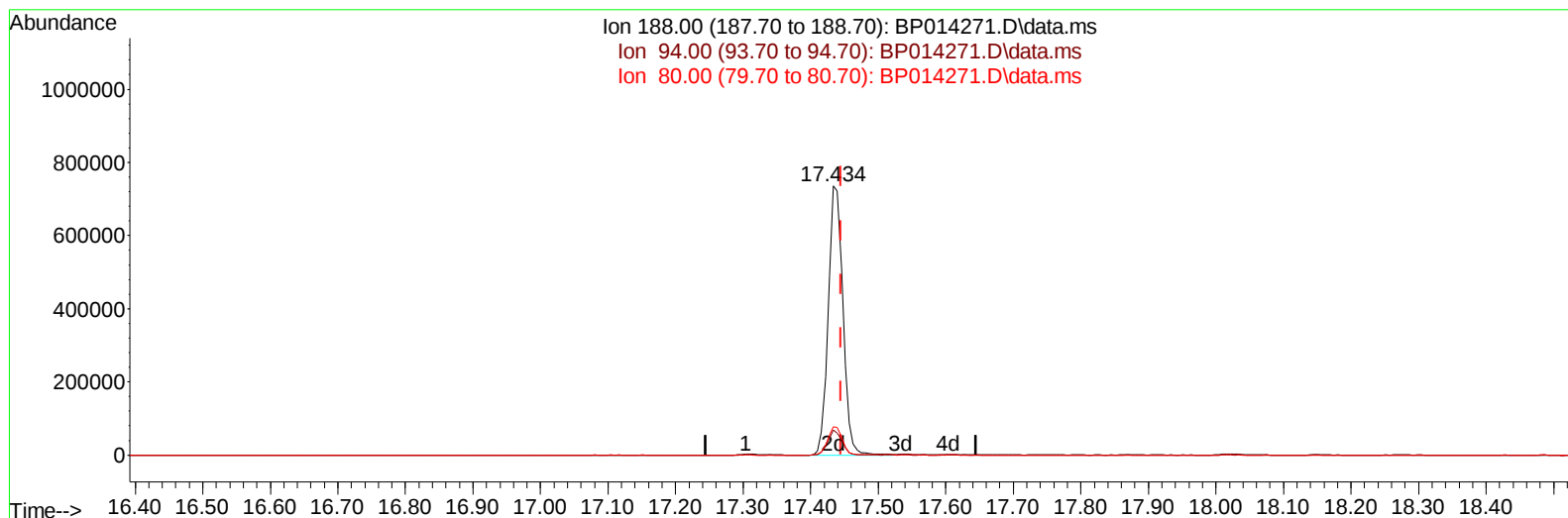
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TIC: BP014271.D\data.ms

(64) Phenanthrene-d10 (I)

17.434min (-0.012) 20.00 ng/ul m

response 1114069

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	7.10	9.37#
80.00	7.80	10.59#
0.00	0.00	0.00

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.034	152	156892	20.000	ng/ul	0.00
20) Naphthalene-d8	10.857	136	614585	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.681	164	453775	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.434	188	1114069m	20.000	ng/ul	-0.01
79) Chrysene-d12	21.516	240	792321	20.000	ng/ul	# 0.00
88) Perylene-d12	24.033	264	951400	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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