

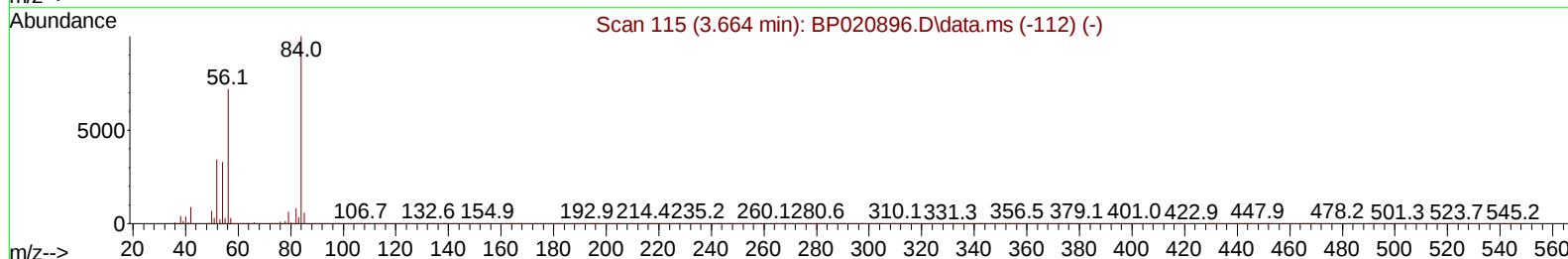
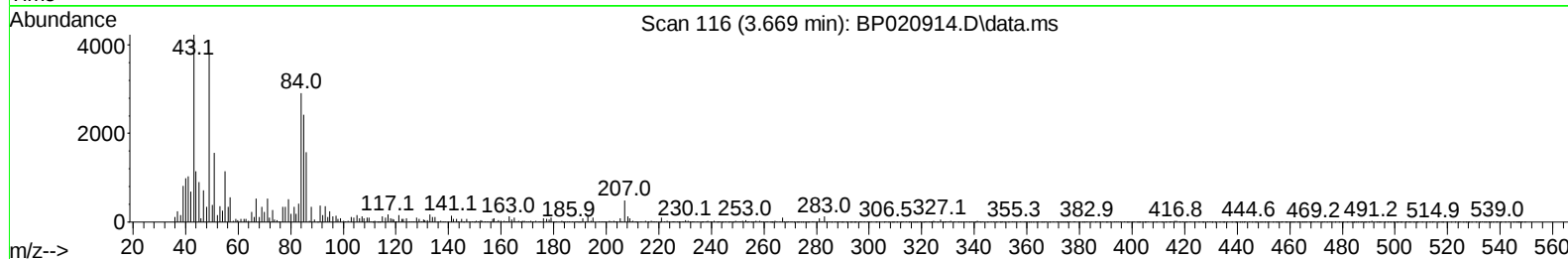
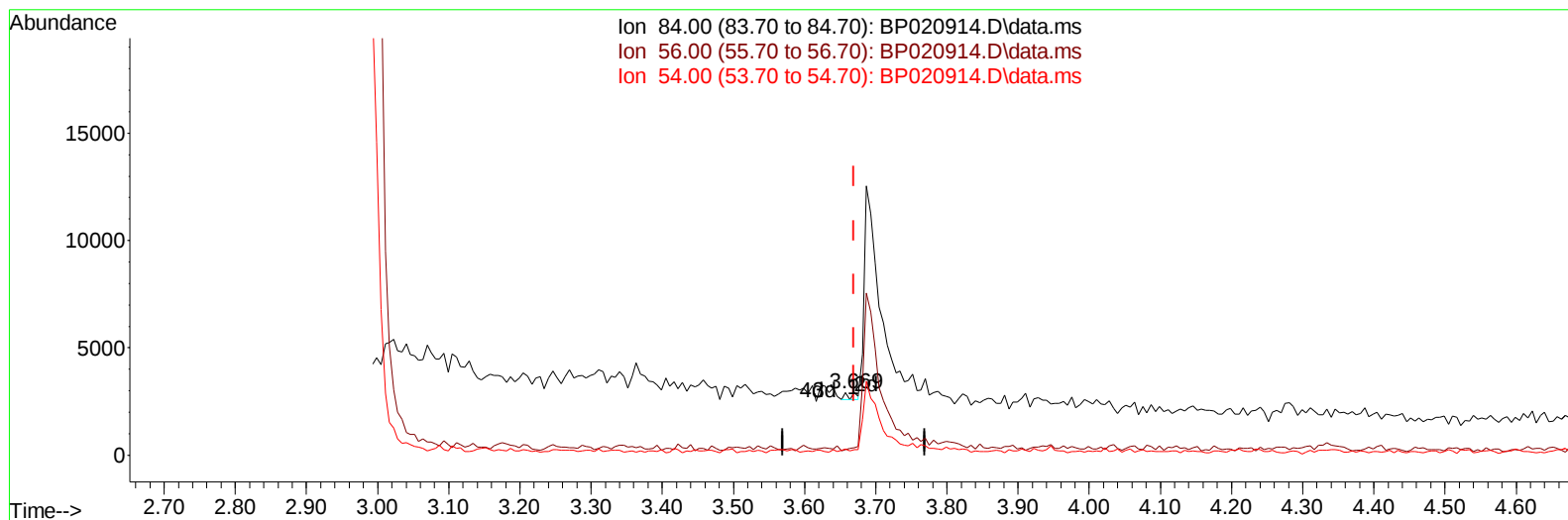
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071124\
 Data File : BP020914.D
 Acq On : 11 Jul 2024 23:14
 Operator : MA/JU
 Sample : P3087-06
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4CB1

Manual IntegrationsAPPROVED

Quant Time: Jul 11 23:57:05 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jul 10 18:38:19 2024
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 07/12/2024
 Supervised By :mohammad ahmed 07/13/2024



TIC: BP020914.D\data.ms

(4) Pyridine-d5 (S)

3.669min (-0.000) 0.01 ng/u1

response	282	
Ion	Exp%	Act%
84.00	100.00	100.00
56.00	70.20	12.02#
54.00	34.30	9.10#
0.00	0.00	0.00

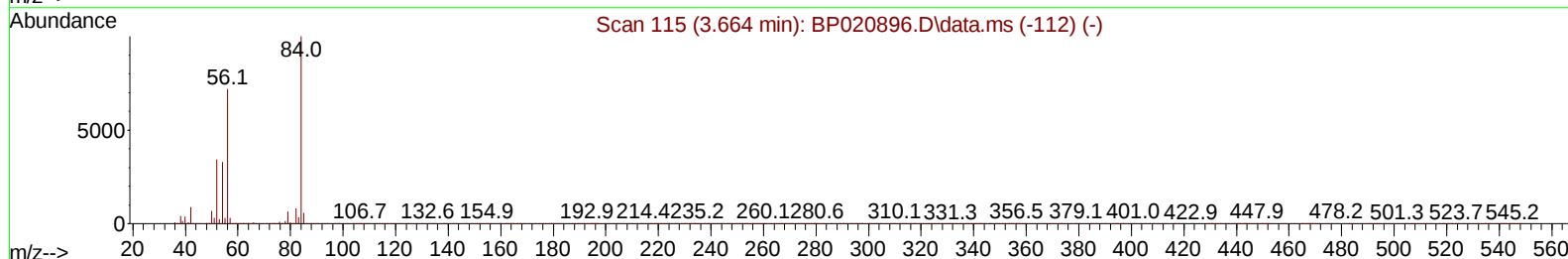
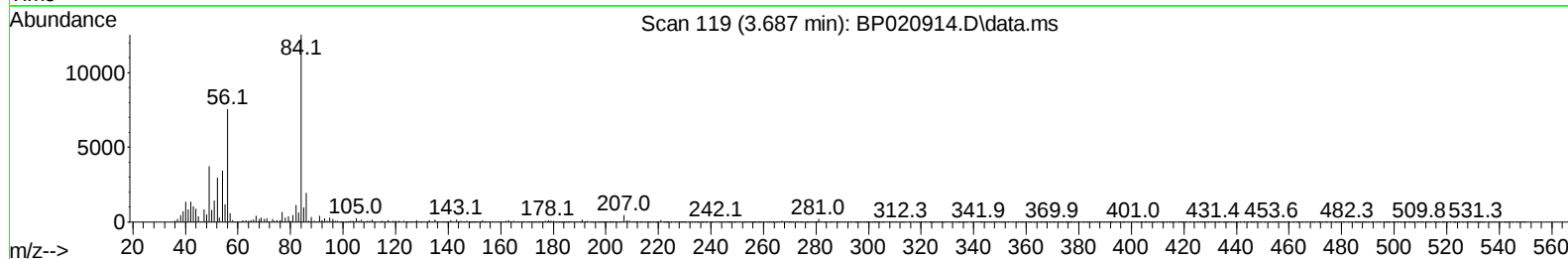
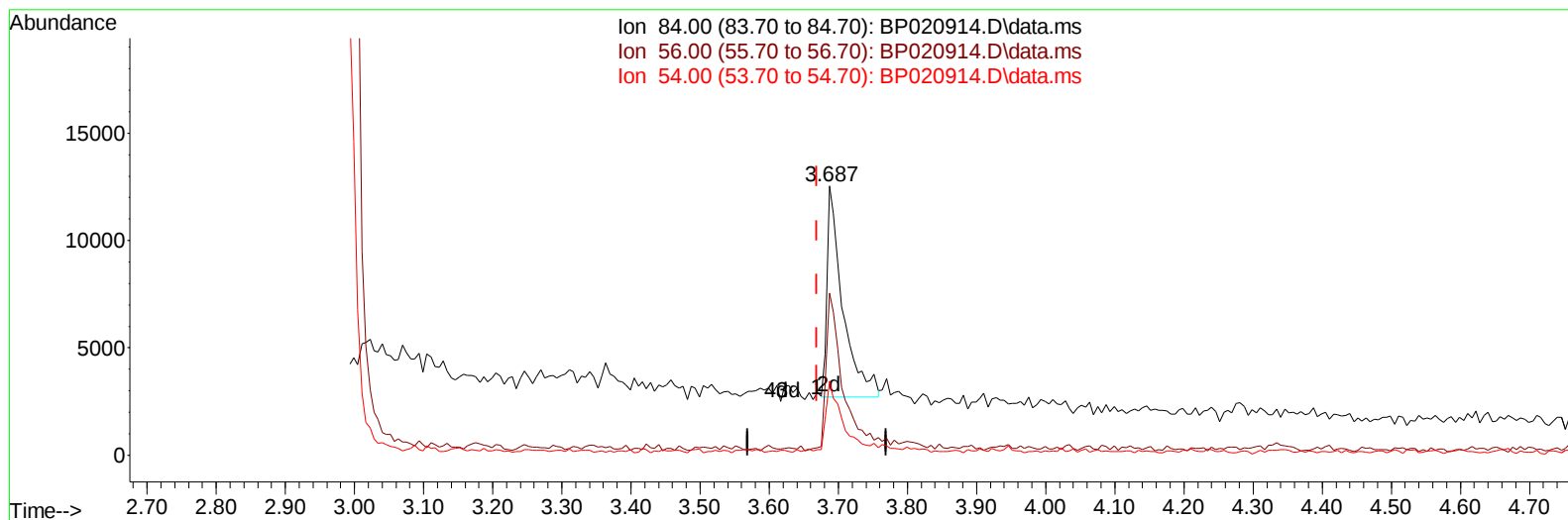
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071124\
 Data File : BP020914.D
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 ALS Vial : 21 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Manual IntegrationsAPPROVED

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TIC: BP020914.D\data.ms

(4) Pyridine-d5 (S)

3.687min (+ 0.018) 0.75 ng/ul m

response 15402

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	70.20	60.32
54.00	34.30	27.44#
0.00	0.00	0.00

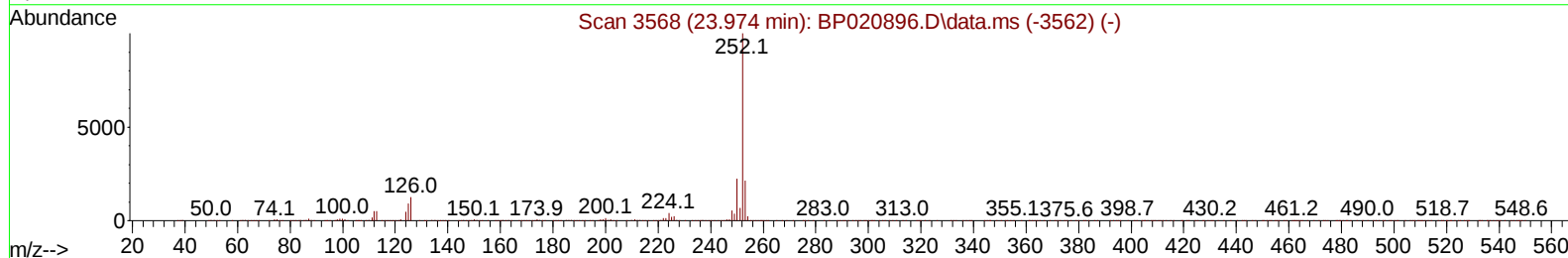
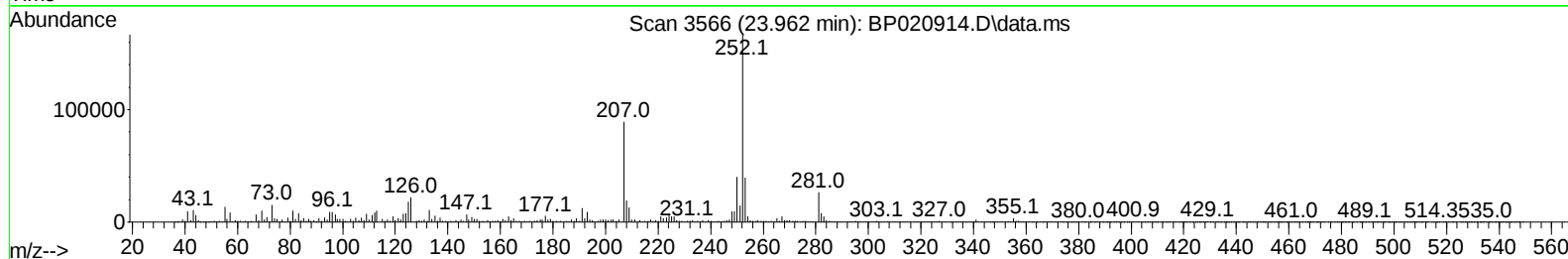
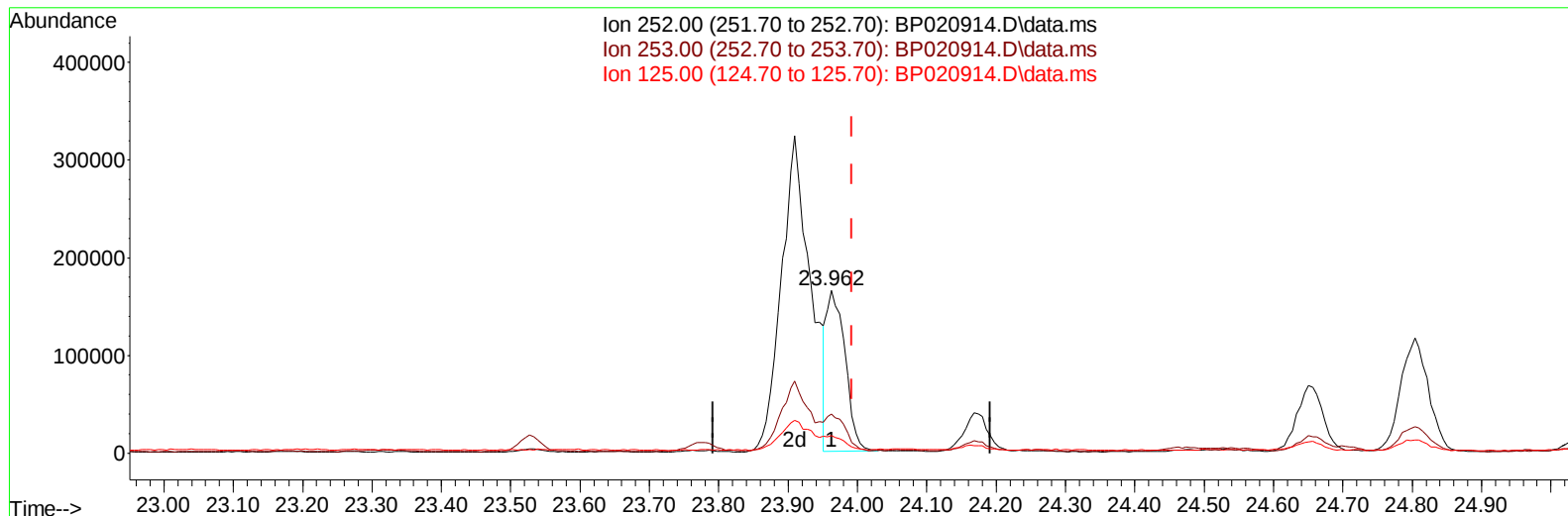
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071124\
 Data File : BP020914.D
 Acq On : 11 Jul 2024 23:14
 Operator : MA/JU
 Sample : P3087-06
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4CB1

Manual IntegrationsAPPROVED

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TIC: BP020914.D\data.ms

(91) Benzo(k)fluoranthene

23.962min (-0.029) 3.07 ng/u1

response 302391

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	21.80	23.81
125.00	9.20	10.75
0.00	0.00	0.00

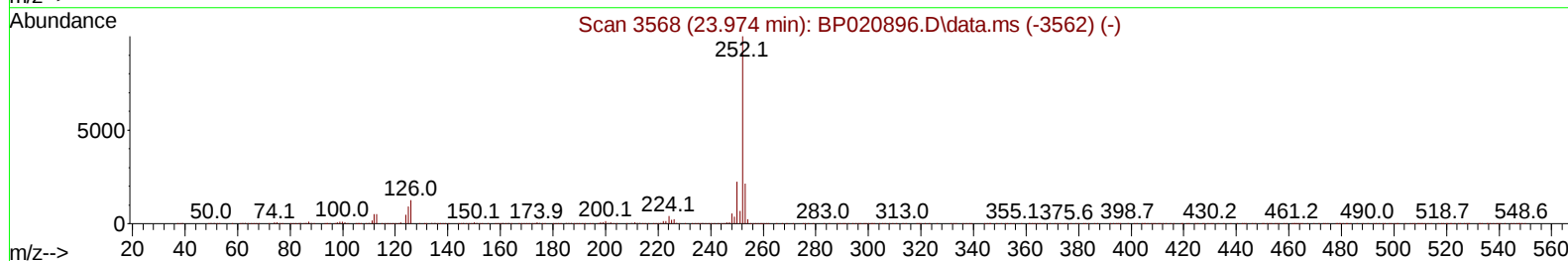
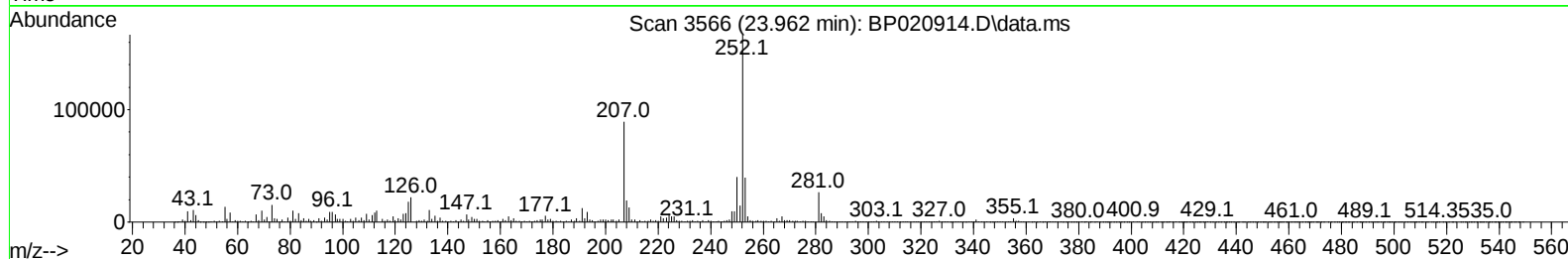
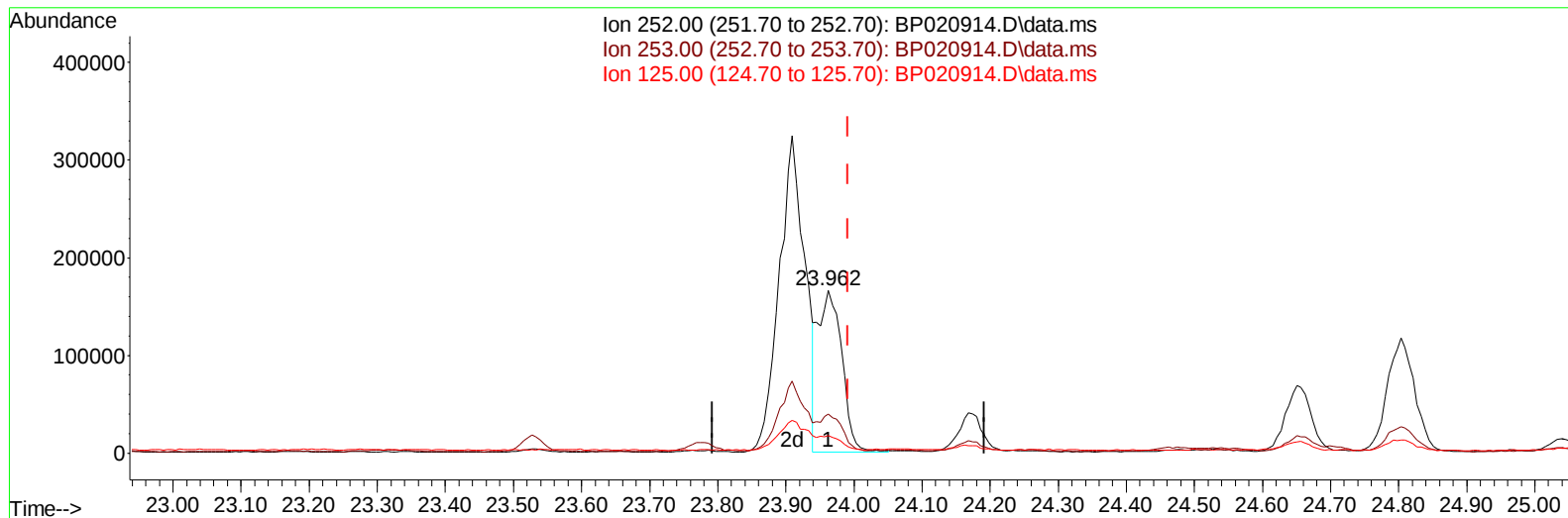
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071124\
Data File : BP020914.D
Acq On : 11 Jul 2024 23:14
Operator : MA/JU
Sample : P3087-06
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CB1

Manual IntegrationsAPPROVED

Quant Time: Jul 11 23:57:05 2024
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TIC: BP020914.D\data.ms

(91) Benzo(k)fluoranthene

23.962min (-0.029) 4.11 ng/u1 m

response 404199

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	21.80	23.81
125.00	9.20	10.75
0.00	0.00	0.00

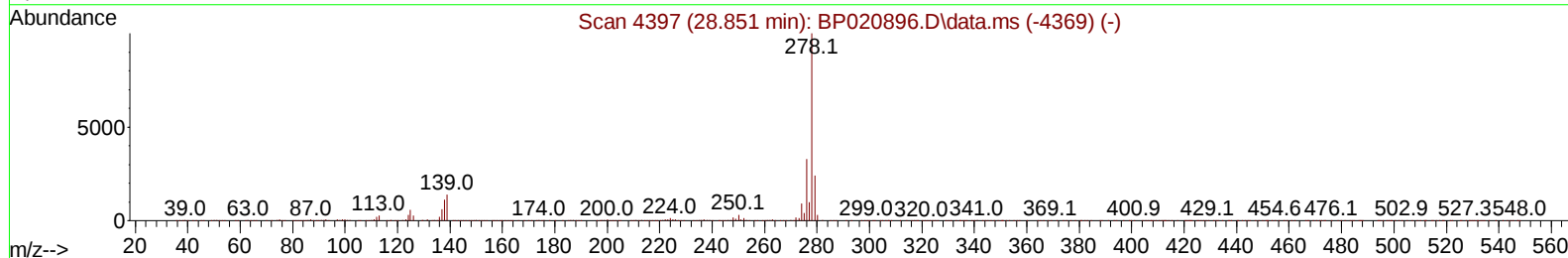
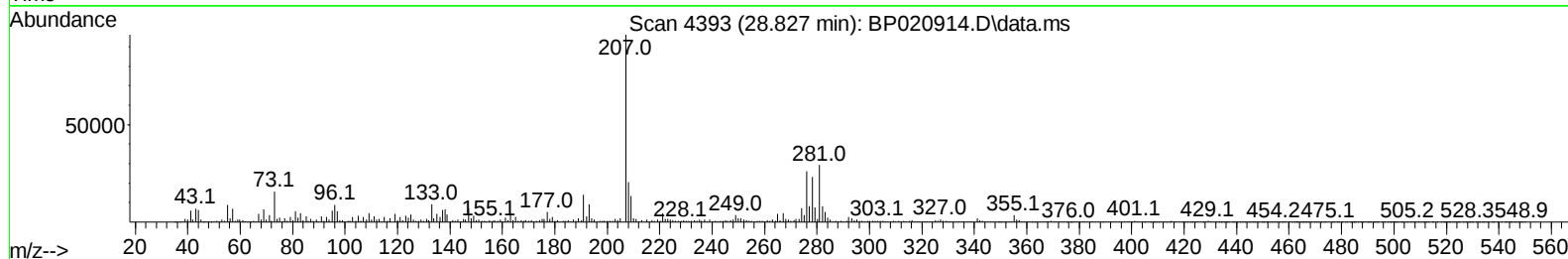
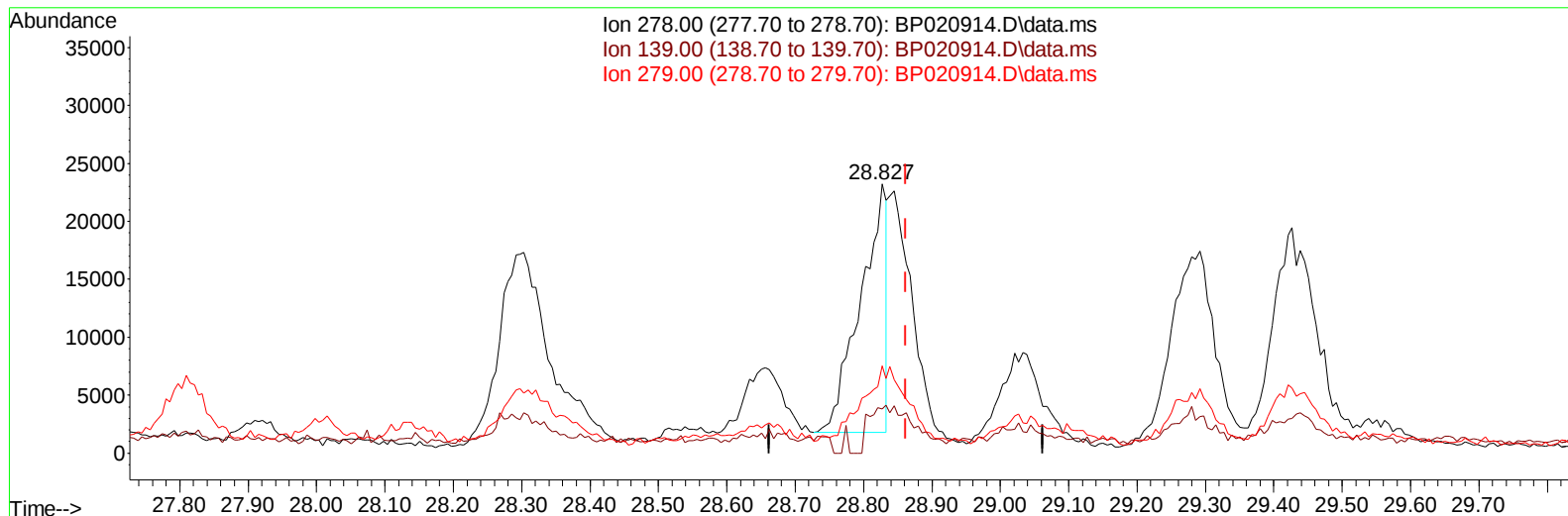
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071124\
Data File : BP020914.D
Acq On : 11 Jul 2024 23:14
Operator : MA/JU
Sample : P3087-06
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Manual IntegrationsAPPROVED

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TIC: BP020914.D\data.ms

(95) Dibenzo(a,h)anthracene

28.827min (-0.035) 0.57 ng/u1

response 56979

Ion	Exp%	Act%
278.00	100.00	100.00
139.00	13.60	16.45#
279.00	24.20	32.36#
0.00	0.00	0.00

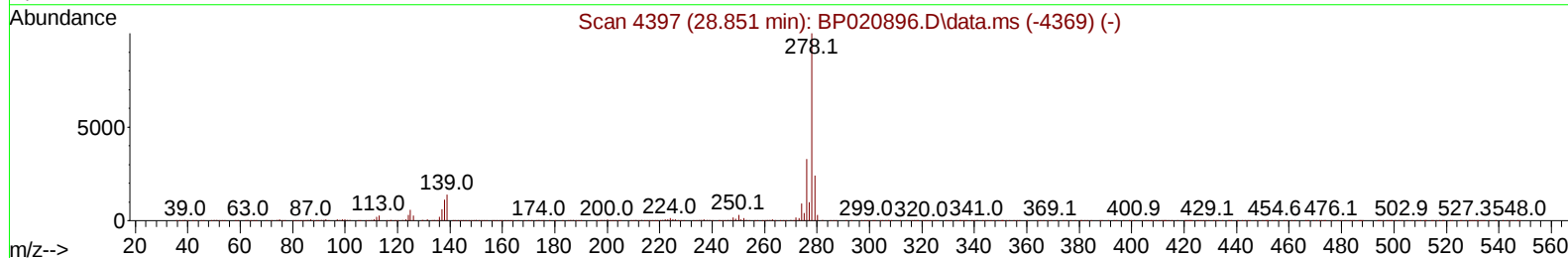
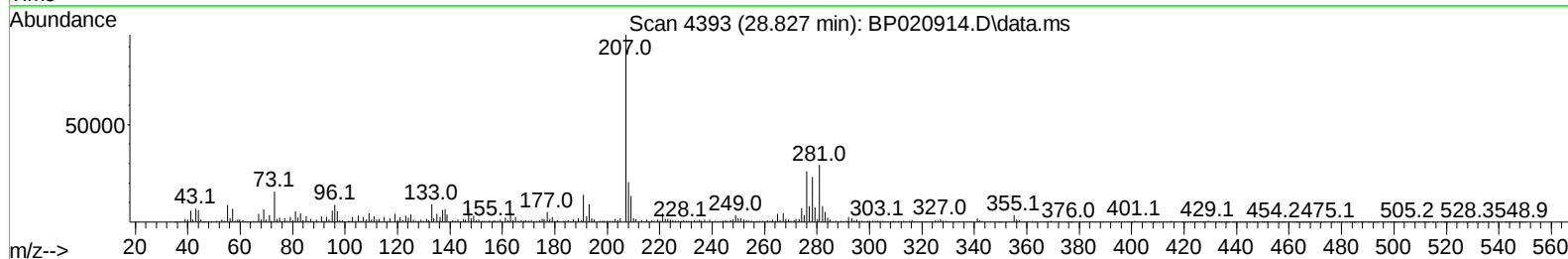
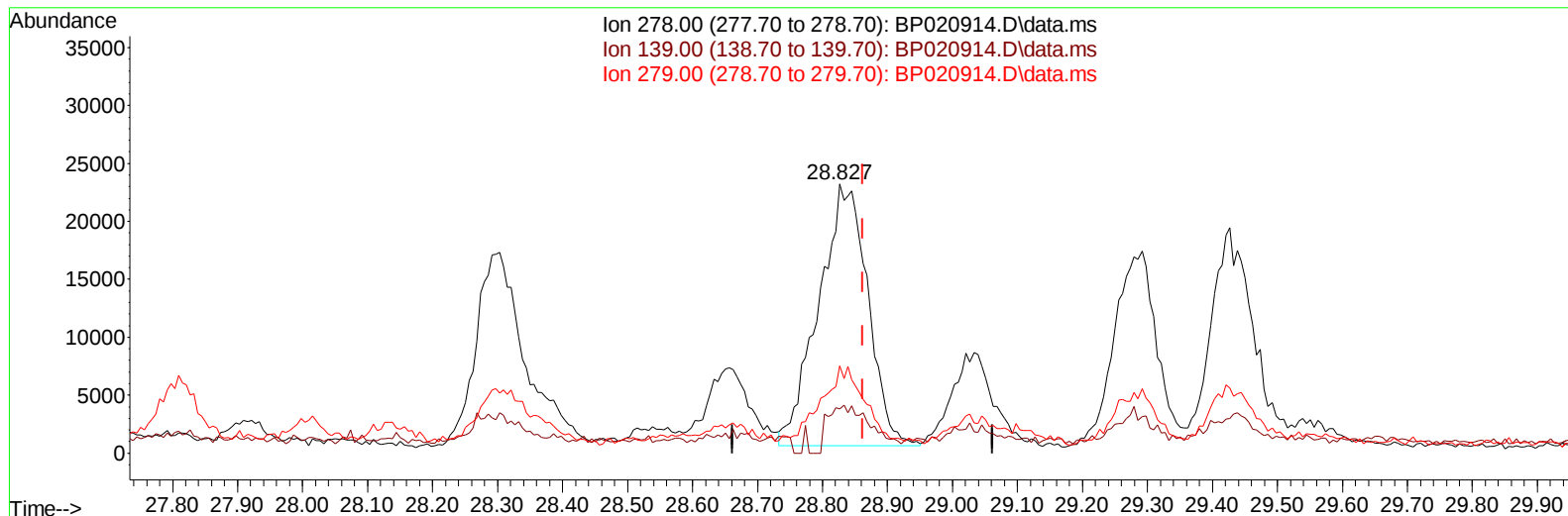
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071124\
Data File : BP020914.D
Acq On : 11 Jul 2024 23:14
Operator : MA/JU
Sample : P3087-06
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CB1

Manual IntegrationsAPPROVED

Quant Time: Jul 11 23:57:05 2024
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TIC: BP020914.D\data.ms

(95) Dibenzo(a,h)anthracene

28.827min (-0.035) 1.18 ng/ul m

response 117475

Ion	Exp%	Act%
278.00	100.00	100.00
139.00	13.60	16.45#
279.00	24.20	32.36#
0.00	0.00	0.00

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 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

Quant Time: Jul 12 00:00:00 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
 Quant Title : SVOA CALIBRATION
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.716	152	320970	20.000	ng/ul	0.00
20) Naphthalene-d8	10.487	136	1342651	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.351	164	788124	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.163	188	1436669	20.000	ng/ul	-0.02
79) Chrysene-d12	21.621	240	1403002	20.000	ng/ul	0.00
88) Perylene-d12	24.968	264	1624411	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	22560	3.198	ng/uL	0.00
4) Pyridine-d5	3.687	84	15402m	0.754	ng/ul	0.02
7) Phenol-d5	6.916	99	491499	18.783	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.069	67	317372	20.013	ng/ul	0.00
11) 2-Chlorophenol-d4	7.263	132	430968	19.989	ng/ul	0.00
15) 4-Methylphenol-d8	8.440	113	362107	16.661	ng/ul	0.00
21) Nitrobenzene-d5	8.875	128	207628	19.910	ng/ul	0.00
24) 2-Nitrophenol-d4	9.587	143	237572	19.904	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.122	165	400864	18.573	ng/ul	0.00
31) 4-Chloroaniline-d4	10.640	131	248125	8.678	ng/ul	0.00
46) Dimethylphthalate-d6	13.757	166	1163267	20.304	ng/ul	-0.01
49) Acenaphthylene-d8	14.034	160	1293861	20.112	ng/ul	-0.02
54) 4-Nitrophenol-d4	14.616	143	111889	13.726	ng/ul	0.00
60) Fluorene-d10	15.363	176	939004	19.978	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.504	200	126156	14.513	ng/ul	0.00
73) Anthracene-d10	17.269	188	1305455	20.459	ng/ul	0.00
81) Pyrene-d10	19.657	212	1331827	15.354	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.733	264	1061952	13.255	ng/ul	-0.02
Target Compounds						
					Qvalue	
72) Phenanthrene	17.204	178	532432	7.144	ng/ul	99
74) Anthracene	17.304	178	88452	1.164	ng/ul	99
78) Di-n-butylphthalate	18.175	149	1893151	22.093	ng/ul	100
80) Fluoranthene	19.304	202	1390836	13.278	ng/ul	100
82) Pyrene	19.686	202	919405	8.625	ng/ul	100
85) Benzo(a)anthracene	21.604	228	595218	6.056	ng/ul	98
86) Bis(2-ethylhexyl)phtha...	21.527	149	907374	14.197	ng/ul	99
87) Chrysene	21.668	228	684315	7.493	ng/ul	97
90) Benzo(b)fluoranthene	23.910	252	844376	8.566	ng/ul	97
91) Benzo(k)fluoranthene	23.962	252	404199m	4.109	ng/ul	
93) Benzo(a)pyrene	24.804	252	326174	3.501	ng/ul	97
94) Indeno(1,2,3-cd)pyrene	28.774	276	326785	2.722	ng/ul	98
95) Dibenzo(a,h)anthracene	28.827	278	117475m	1.182	ng/ul	

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

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