

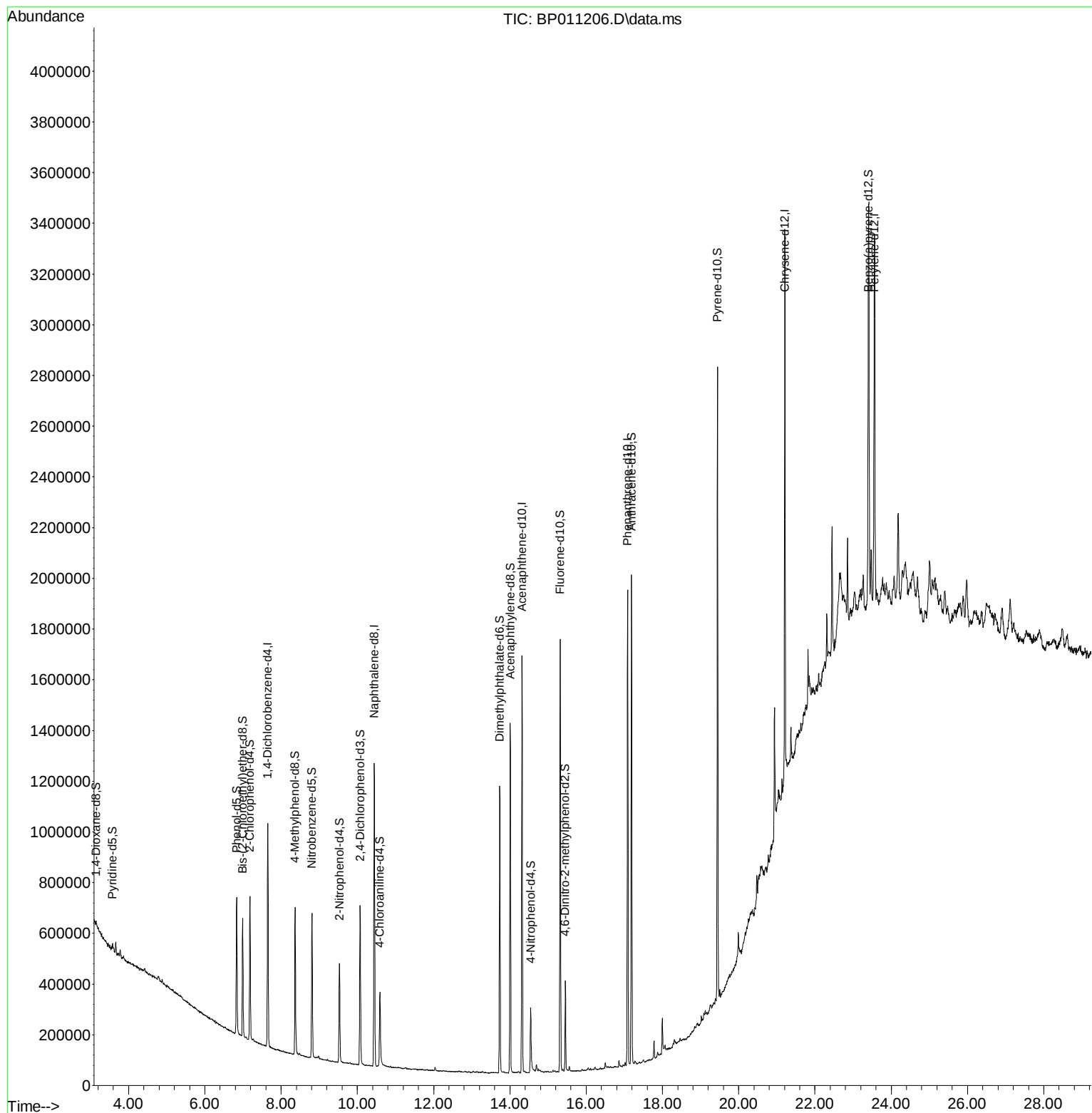
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080122\
Data File : BP011206.D
Acq On : 01 Aug 2022 18:46
Operator : CG/JU
Sample : N3790-13
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBUK7

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 08/02/2022
Supervised By :mohammad ahmed 08/03/2022

Quant Time: Aug 01 23:01:22 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP072822.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jul 28 16:05:22 2022
Response via : Initial Calibration



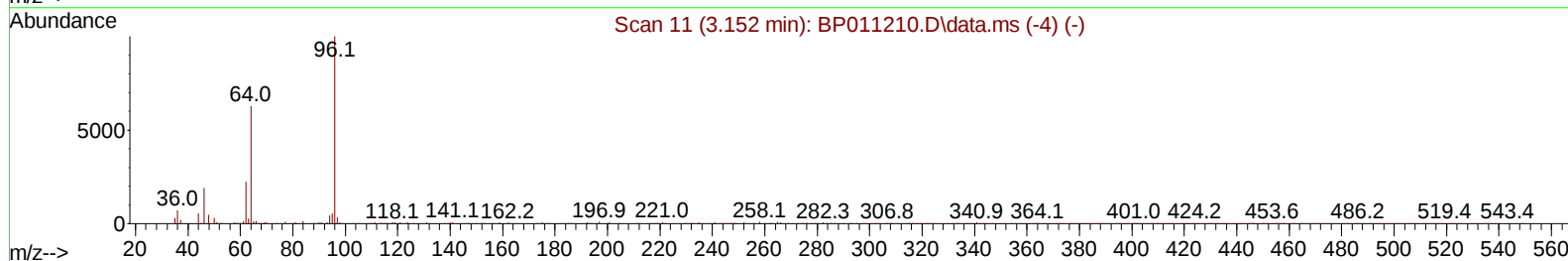
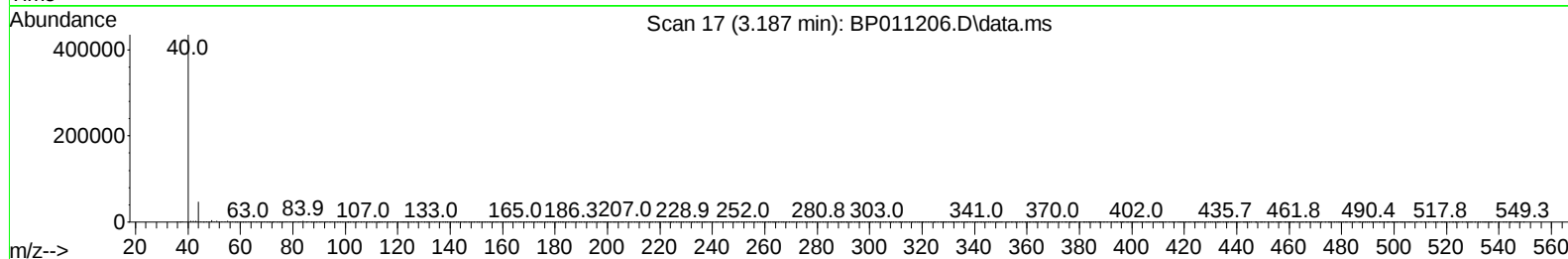
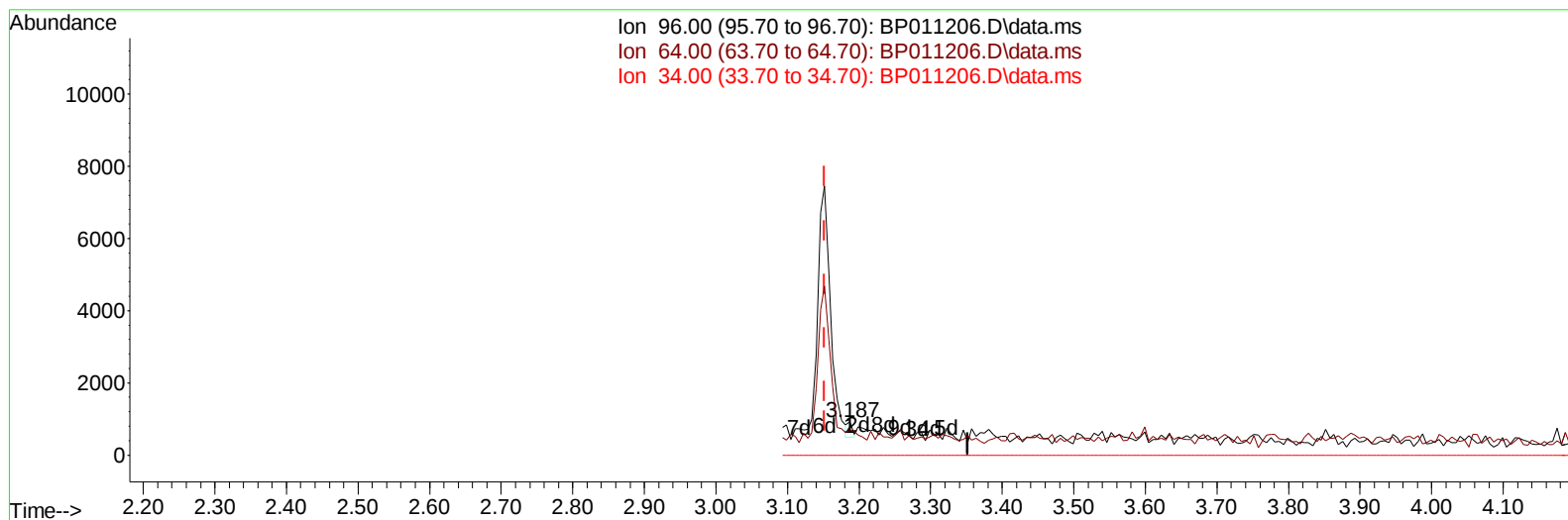
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TIC: BP011206.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.187min (+ 0.035) 0.04 ng/uL

response	173	
Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.90	74.19
34.00	0.00	0.00
0.00	0.00	0.00

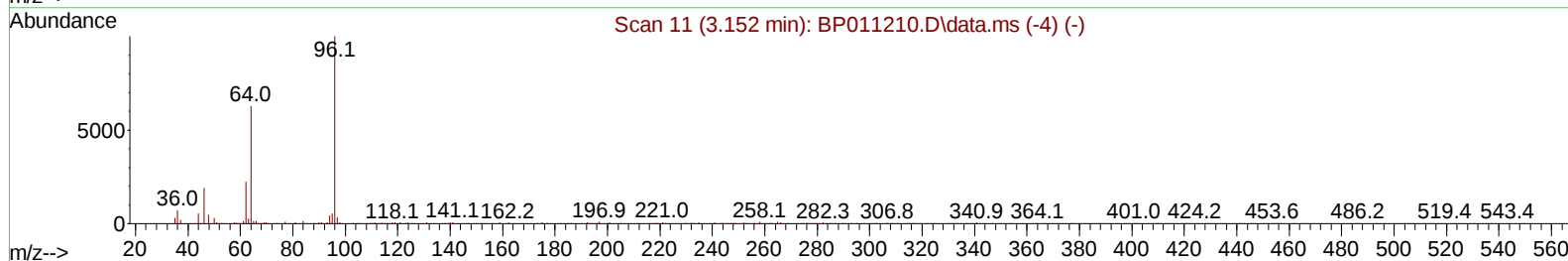
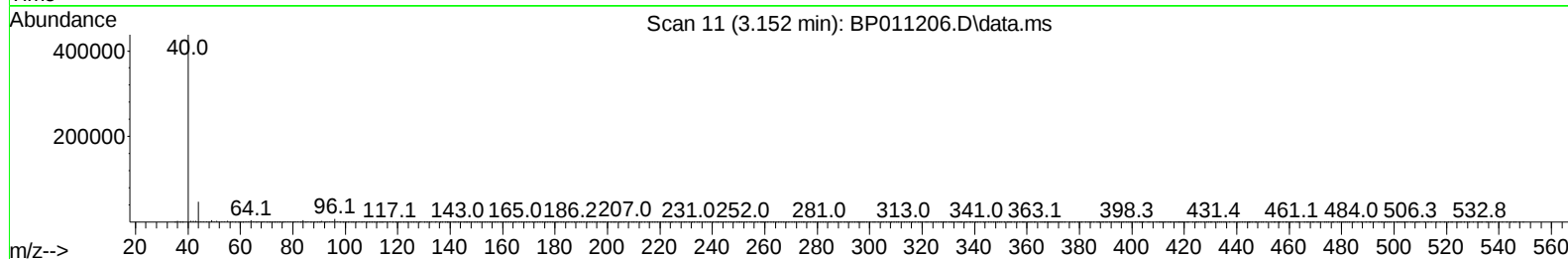
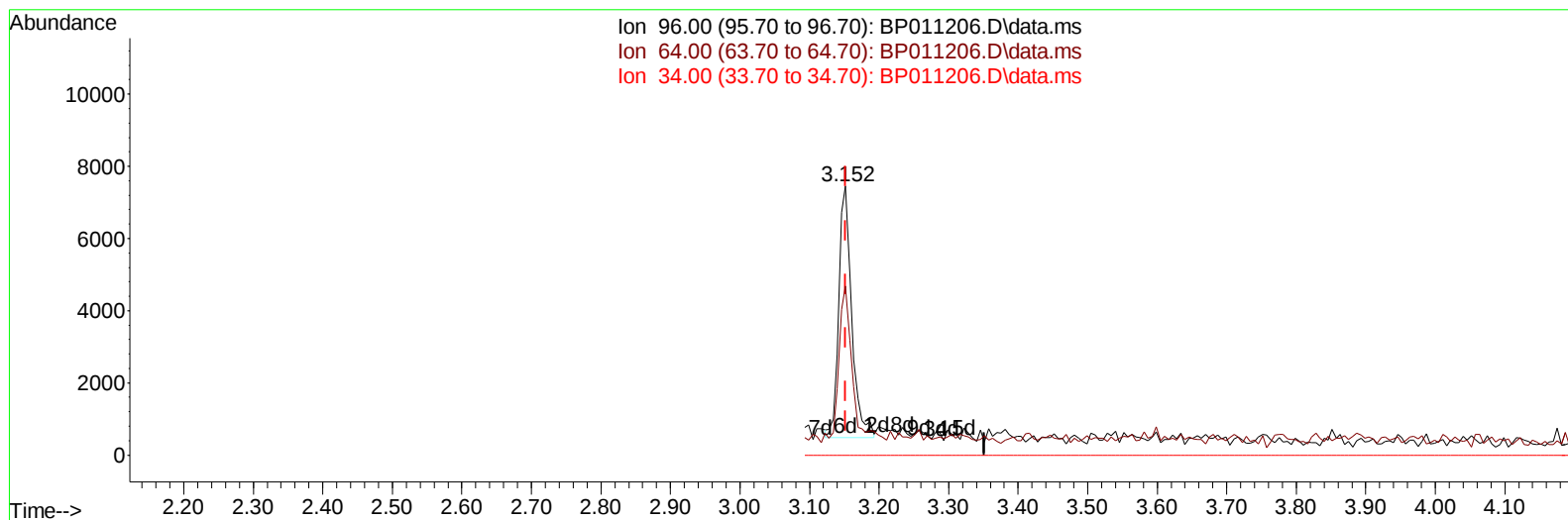
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TIC: BP011206.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.152min (-0.000) 1.96 ng/uL m

response 8967

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.90	62.90#
34.00	0.00	0.00
0.00	0.00	0.00

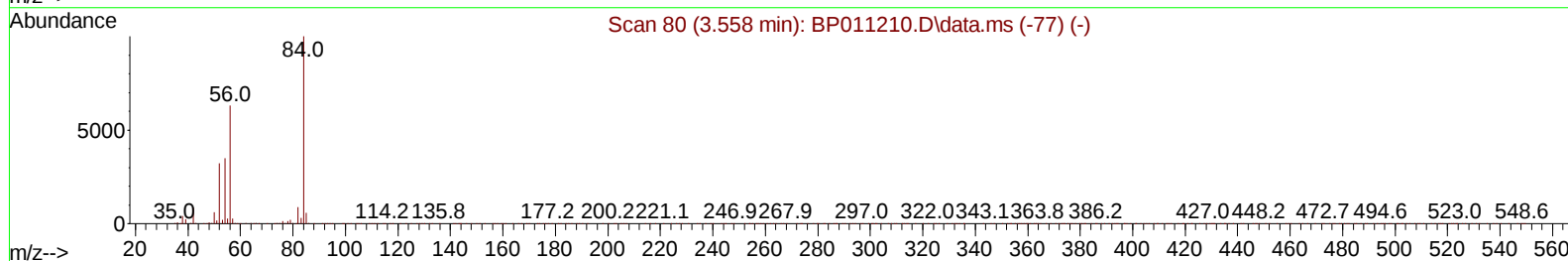
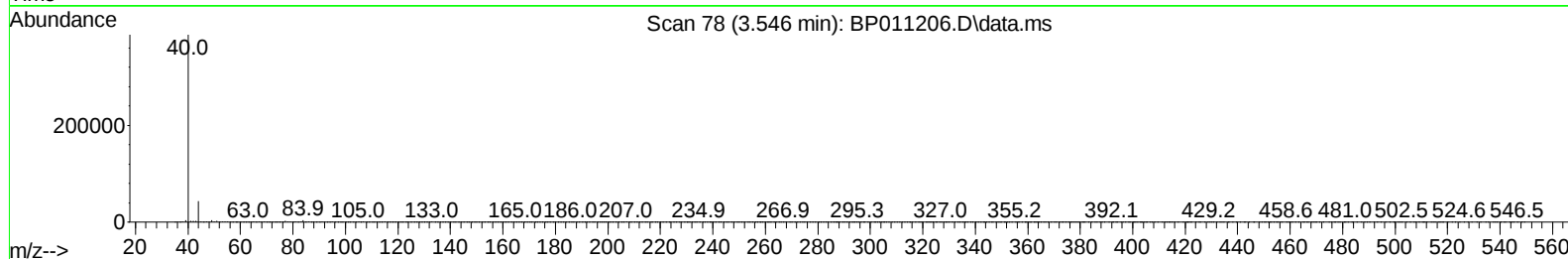
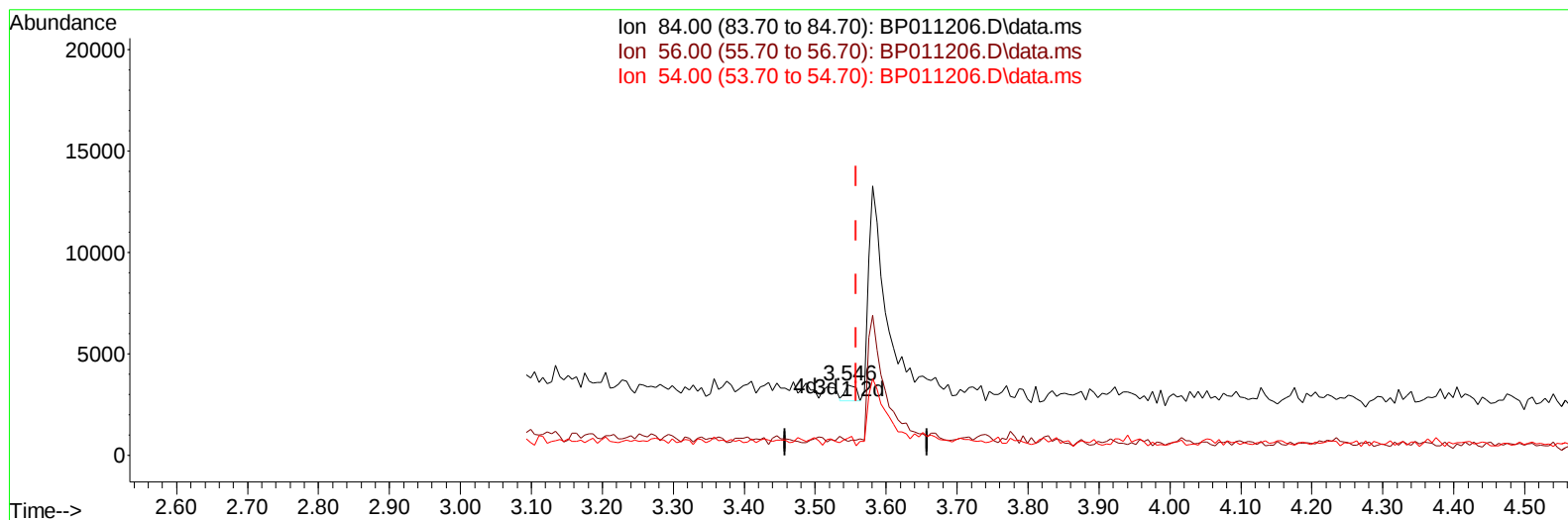
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(4) Pyridine-d5 (S)

3.546min (-0.012) 0.07 ng/u1

response 875

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	84.50	19.69#
54.00	44.80	24.50#
0.00	0.00	0.00

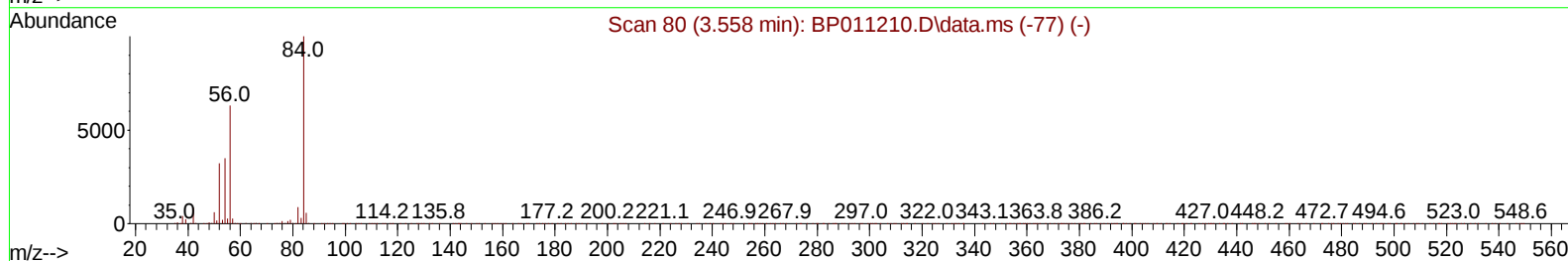
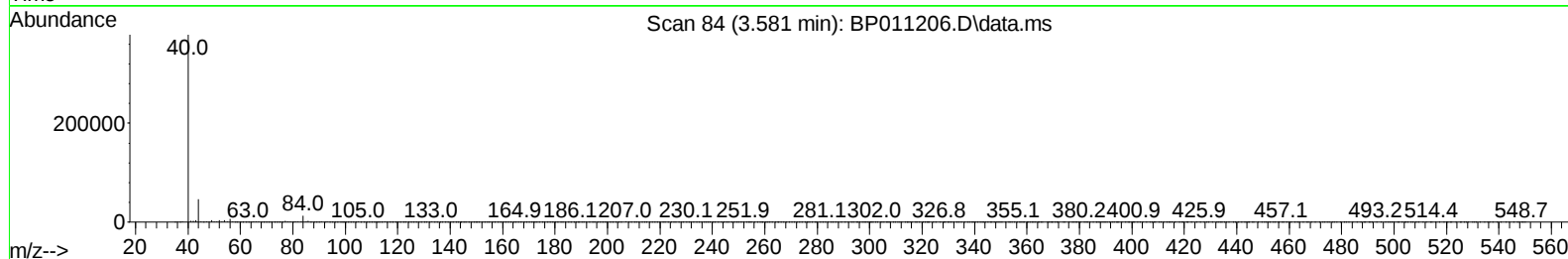
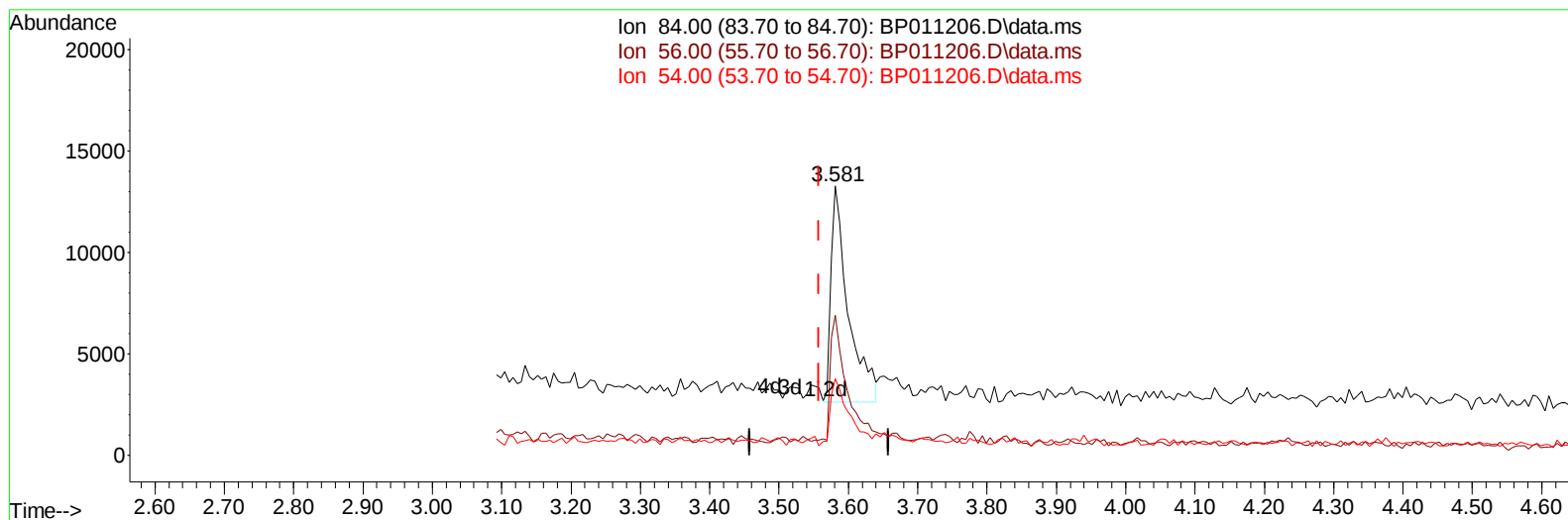
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TIC: BP011206.D\data.ms

(4) Pyridine-d5 (S)

3.581min (+ 0.023) 1.42 ng/ul m

response 18393

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	84.50	52.11#
54.00	44.80	28.46#
0.00	0.00	0.00

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 Misc :
 ALS Vial : 15 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.646	152	209901	20.000	ng/ul	0.00
20) Naphthalene-d8	10.440	136	895649	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.316	164	485598	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.086	188	990311	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.210	240	896939	20.000	ng/ul	# 0.00
88) Perylene-d12	23.557	264	940894	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.152	96	8967m	1.956	ng/uL	0.00
4) Pyridine-d5	3.581	84	18393m	1.418	ng/ul	0.02
7) Phenol-d5	6.834	99	284251	15.515	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.993	67	203561	15.713	ng/ul	0.00
11) 2-Chlorophenol-d4	7.181	132	232339	16.092	ng/ul	0.00
15) 4-Methylphenol-d8	8.369	113	207741	13.861	ng/ul	0.00
21) Nitrobenzene-d5	8.810	128	116053	16.673	ng/ul	0.00
24) 2-Nitrophenol-d4	9.534	143	107055	16.056	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.069	165	201130	17.117	ng/ul	0.00
31) 4-Chloroaniline-d4	10.593	131	167874	8.813	ng/ul	0.00
46) Dimethylphthalate-d6	13.734	166	663430	20.053	ng/ul	0.00
49) Acenaphthylene-d8	14.004	160	861346	20.718	ng/ul	0.00
54) 4-Nitrophenol-d4	14.545	143	51257	8.613	ng/ul	0.01
60) Fluorene-d10	15.316	176	610902	21.974	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.451	200	48187	9.653	ng/ul	0.00
73) Anthracene-d10	17.186	188	939377	21.528	ng/ul	0.00
81) Pyrene-d10	19.445	212	1116226	23.160	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.404	264	1031595	22.074	ng/ul	0.00

Target Compounds Qvalue

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