

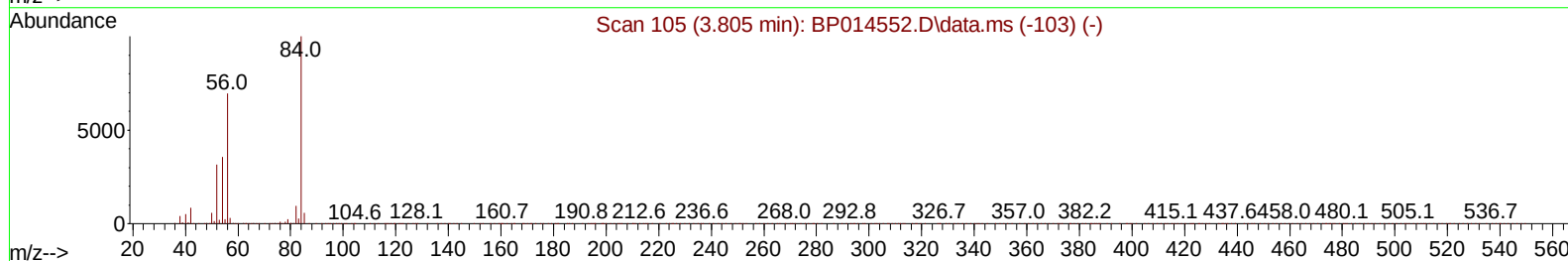
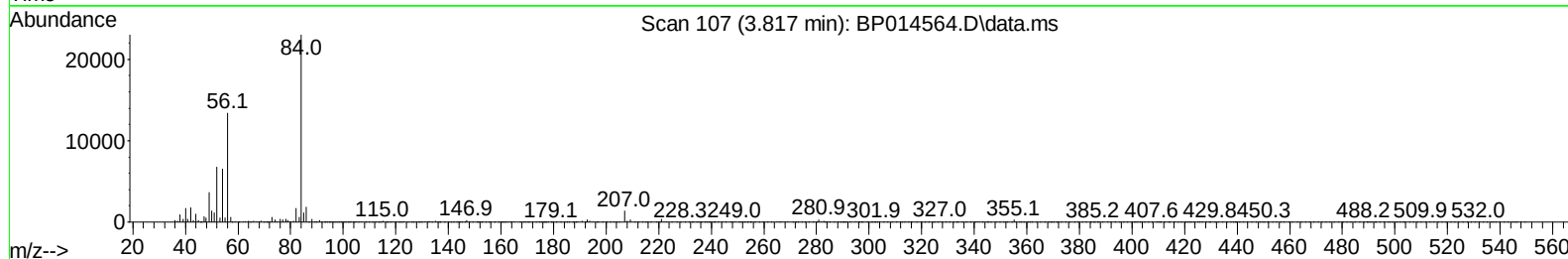
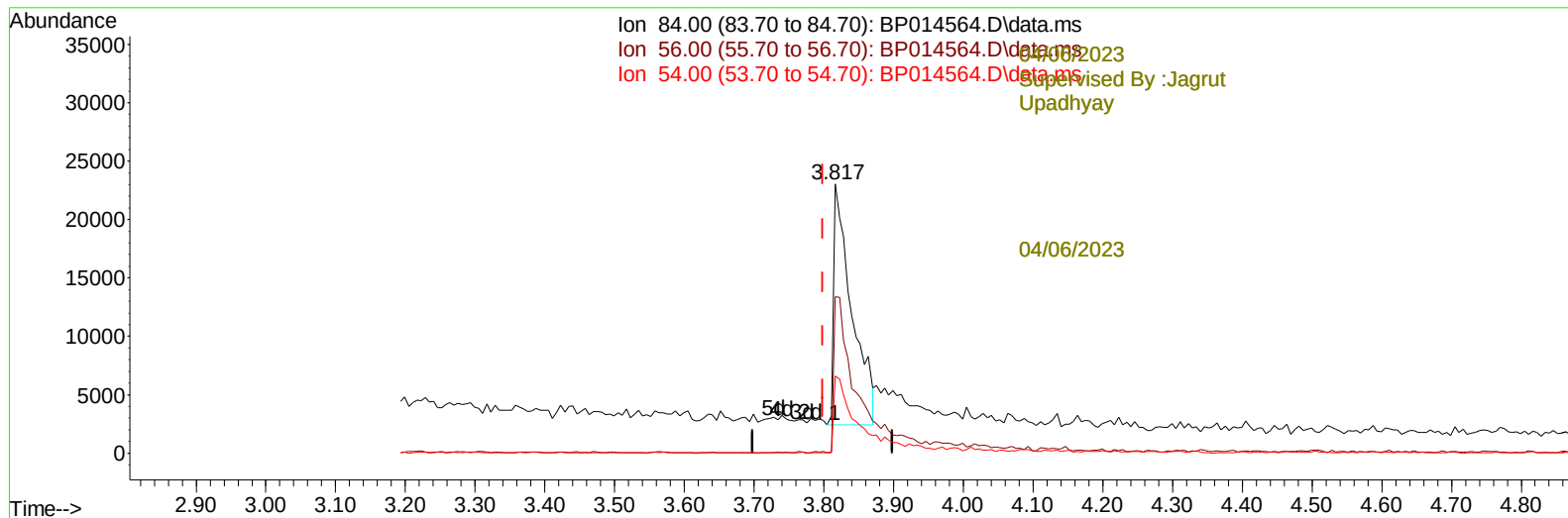
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040523\
 Data File : BP014564.D
 Acq On : 05 Apr 2023 18:51
 Operator : CG/JU
 Sample : 02122-14
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB9C8

Manual IntegrationsAPPROVED

Quant Time: Apr 06 00:58:48 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 06 00:55:13 2023
 Response via : Initial Calibration

Reviewed By :Christian
 Giraldo



TIC: BP014564.D\data.ms

(4) Pyridine-d5 (S)

3.817min (+ 0.018) 4.87 ng/uI

response 37027

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	58.25
54.00	32.60	28.64
0.00	0.00	0.00

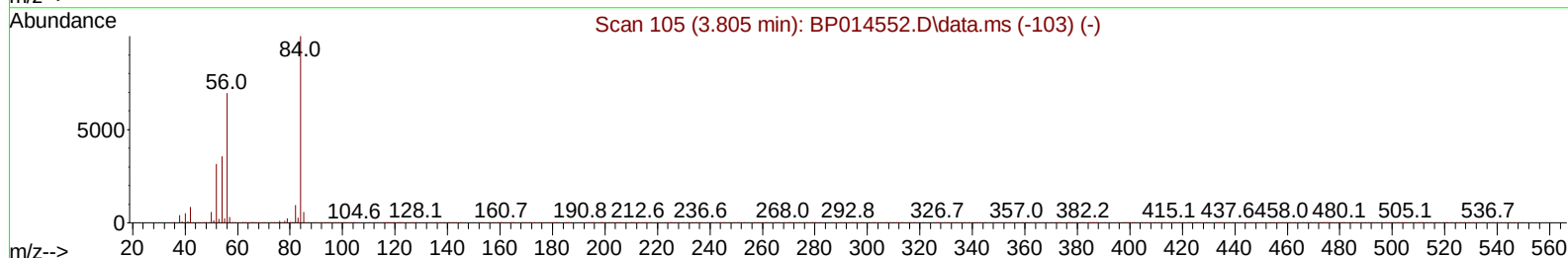
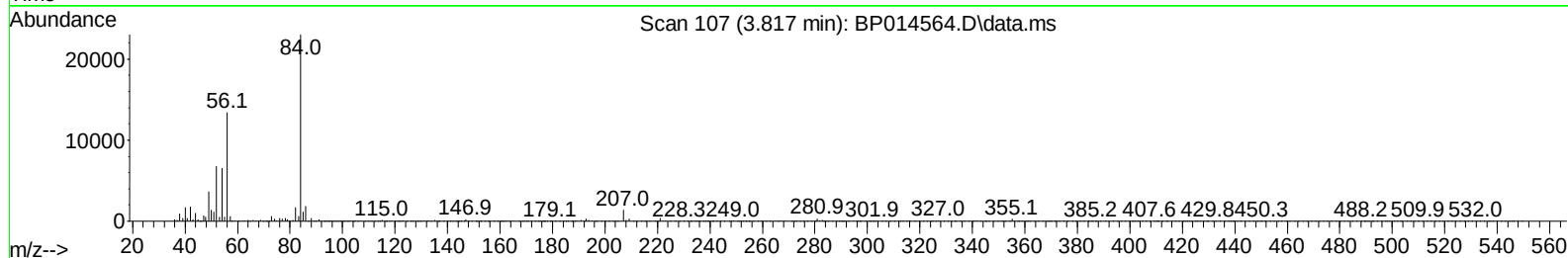
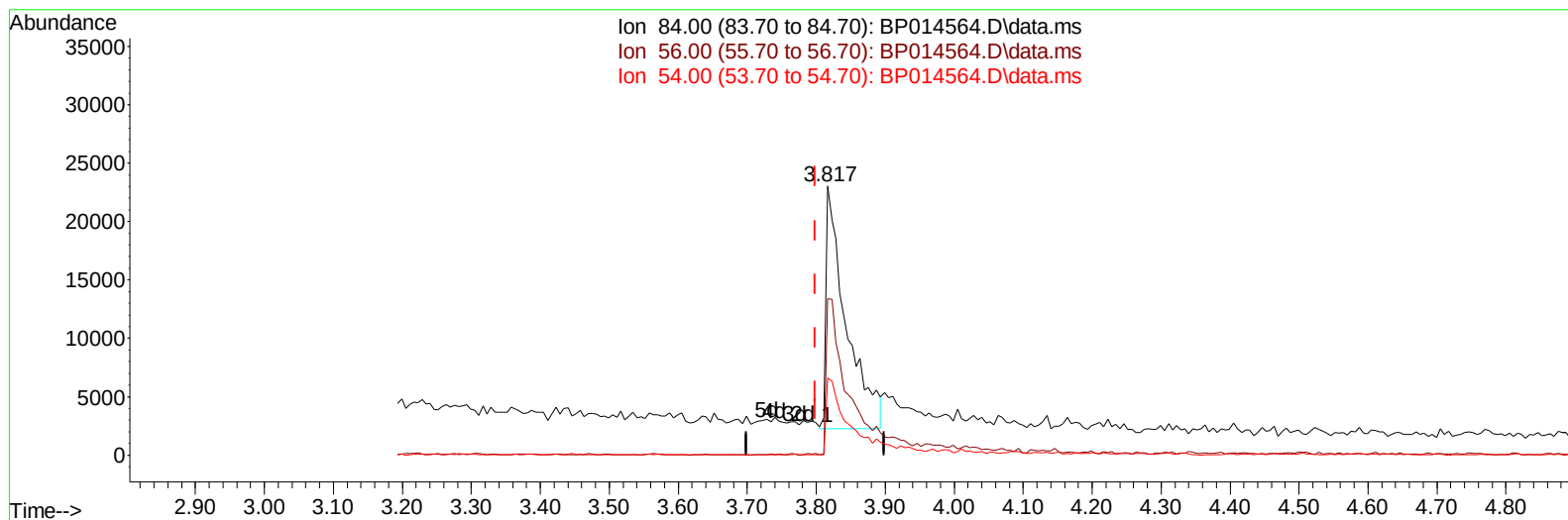
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Supervised By :Jagrut Upadhyay 04/06/2023



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(4) Pyridine-d5 (S)

3.817min (+ 0.018) 5.52 ng/ul m

response 41947

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	58.25
54.00	32.60	28.64
0.00	0.00	0.00

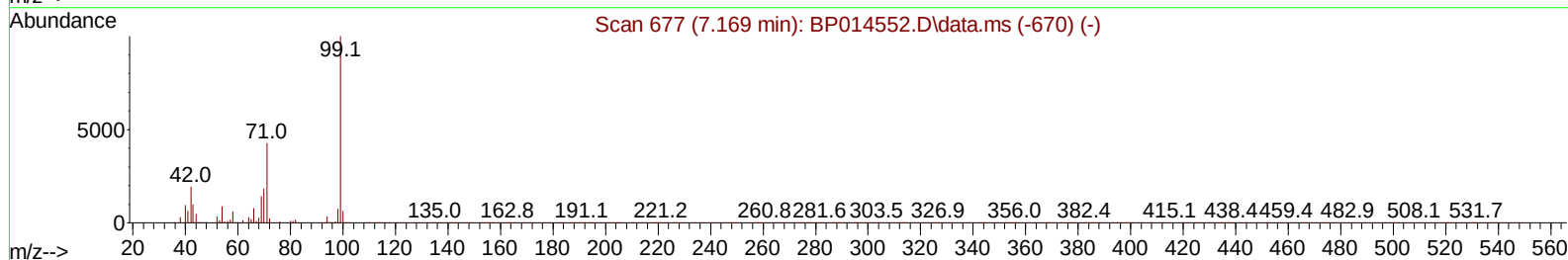
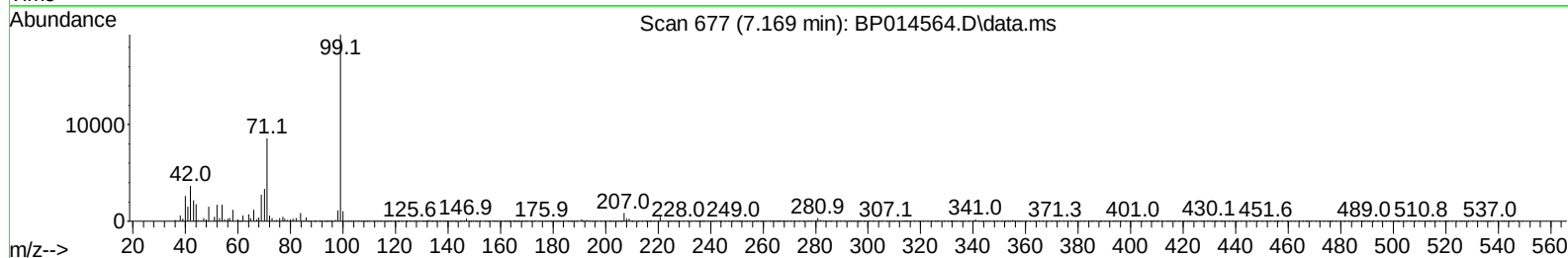
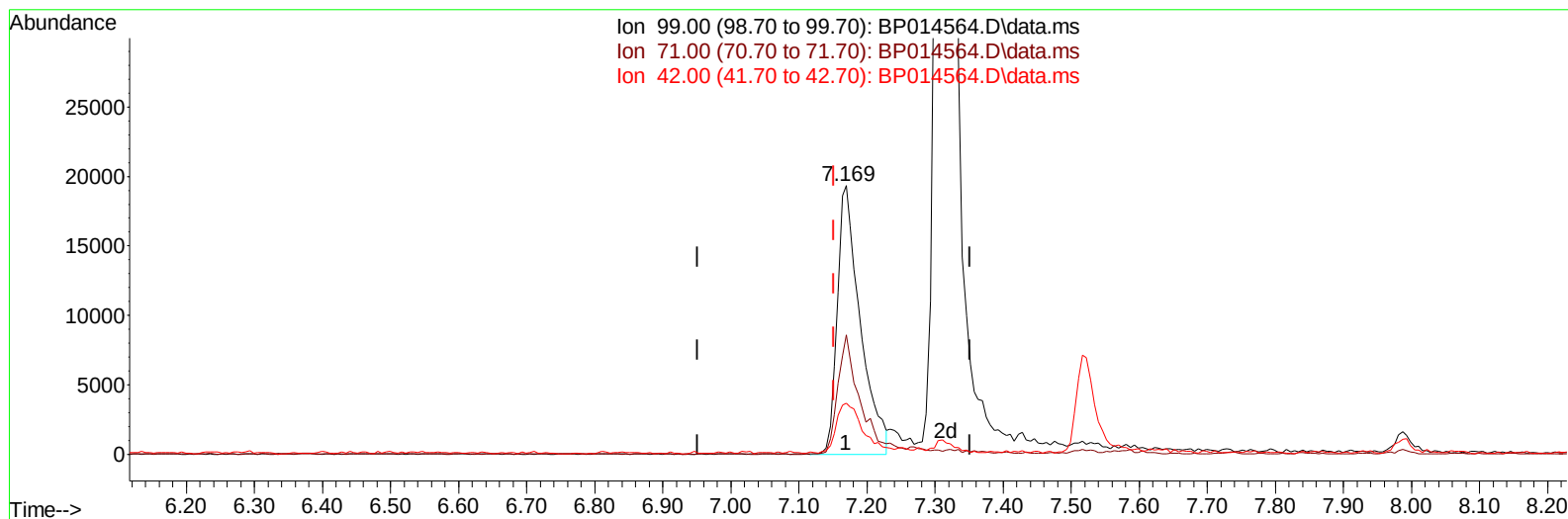
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Instrument :
BNA_P
ClientSampleId :
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TIC: BP014564.D\data.ms

(7) Phenol-d5 (S)

7.169min (+ 0.018) 4.29 ng/u1

response 45490

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	40.30	44.41
42.00	18.90	19.06
0.00	0.00	0.00

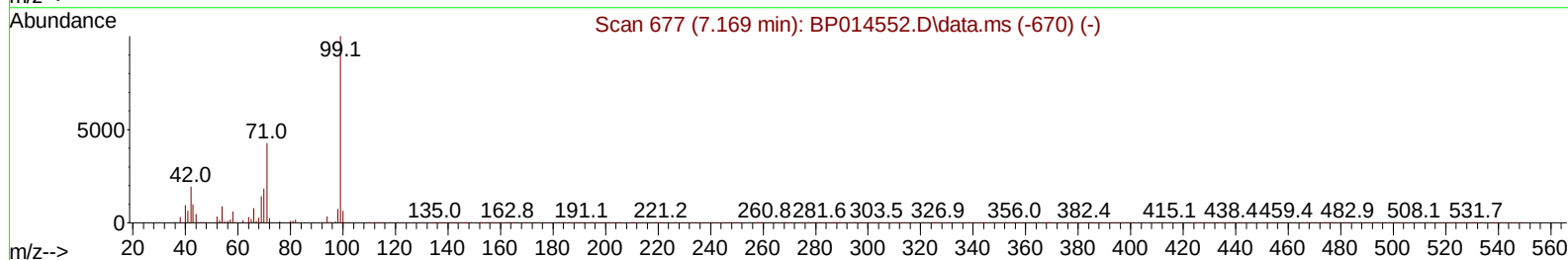
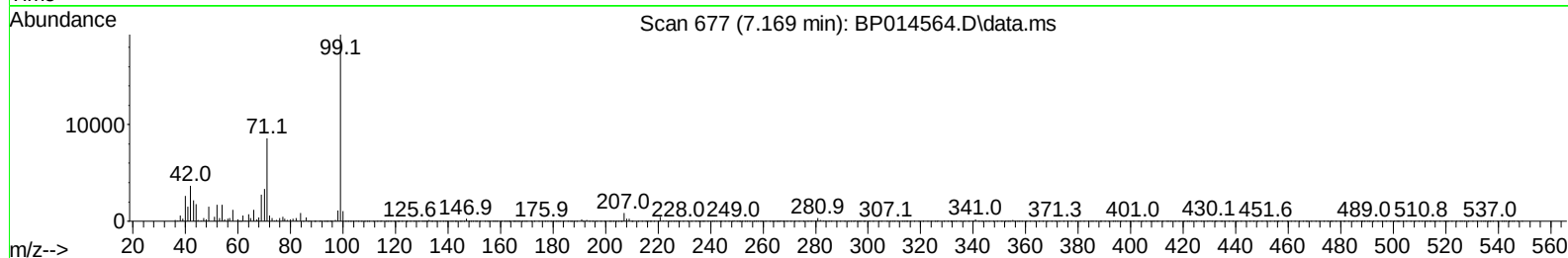
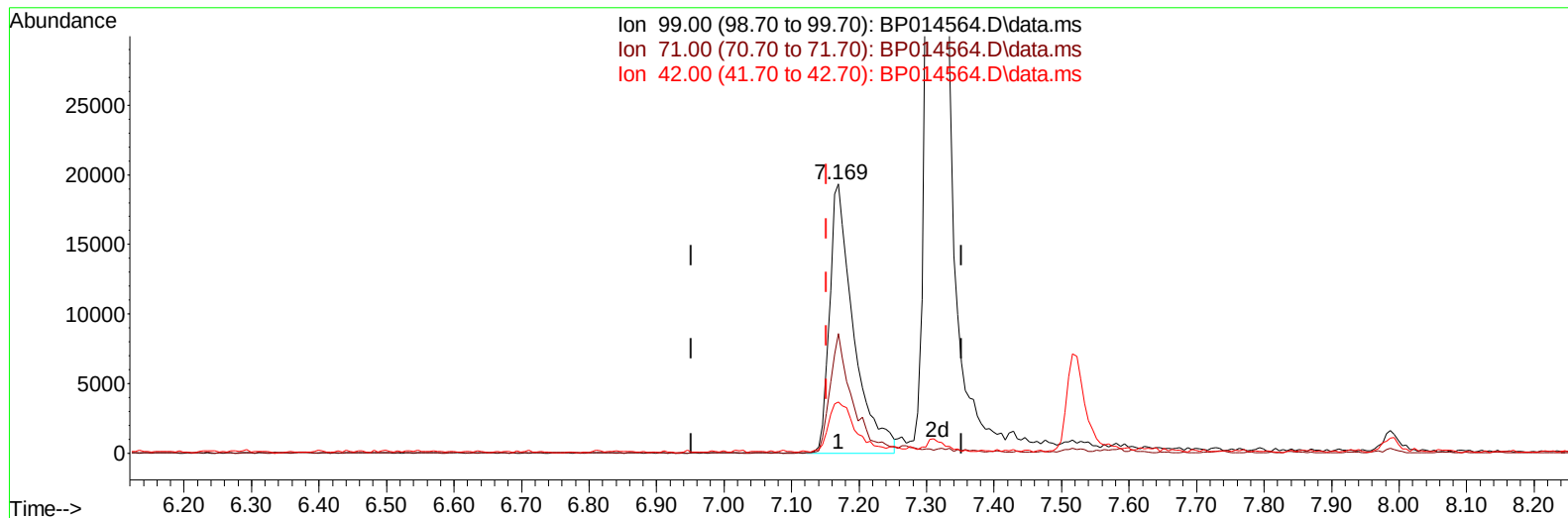
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Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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TIC: BP014564.D\data.ms

(7) Phenol-d5 (S)

7.169min (+ 0.018) 4.48 ng/ul m

response 47604

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	40.30	44.41
42.00	18.90	19.06
0.00	0.00	0.00

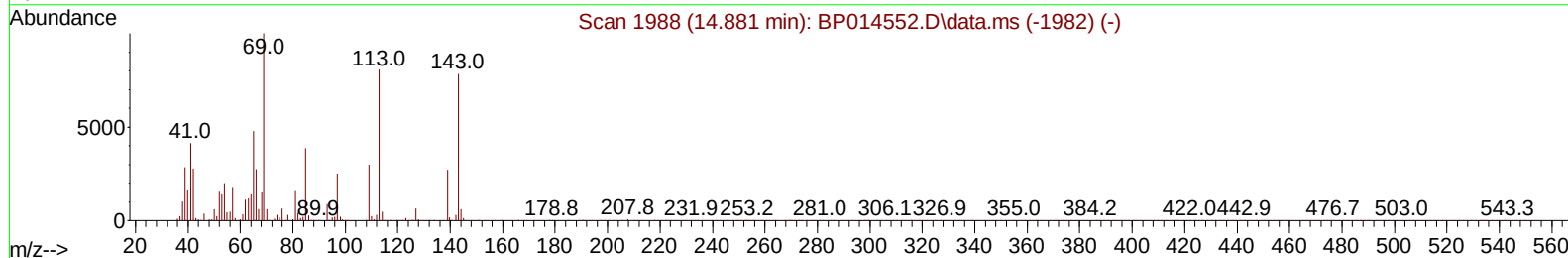
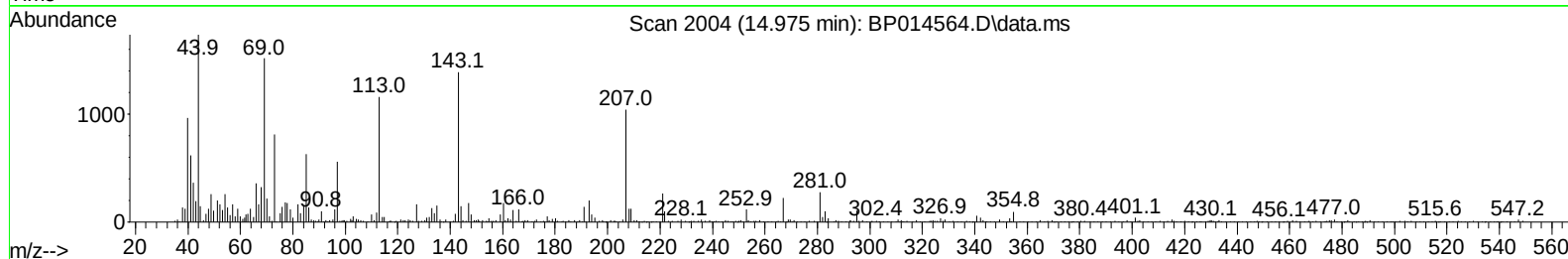
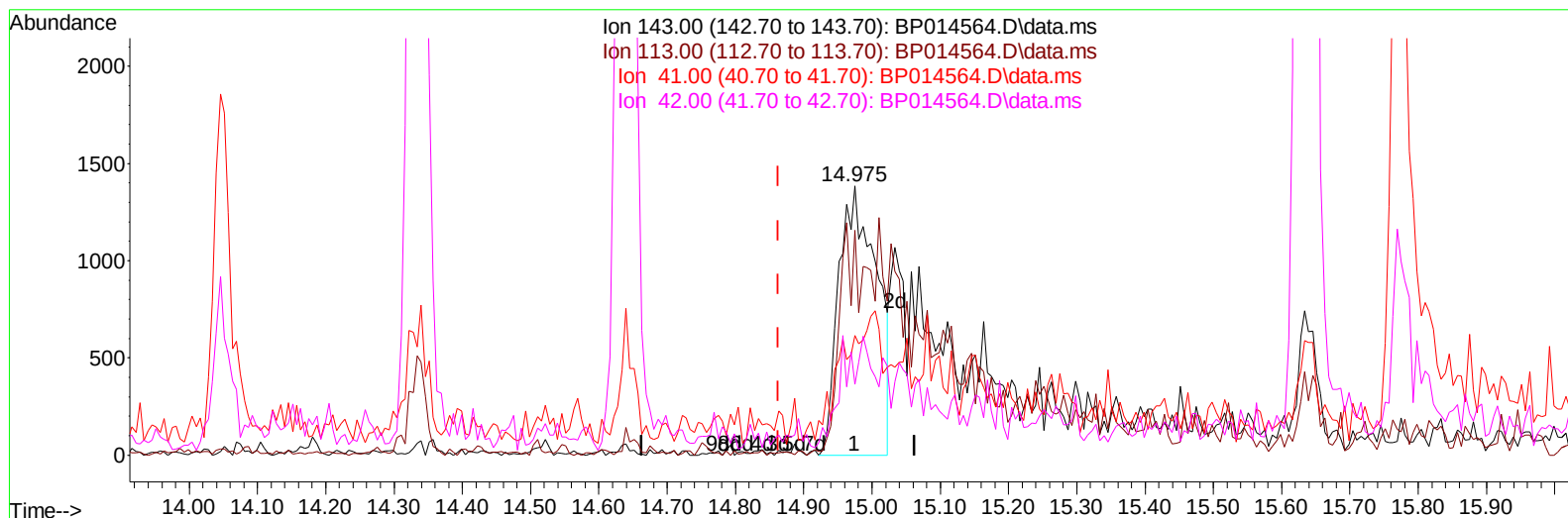
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TIC: BP014564.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.975min (+ 0.112) 1.35 ng/ul

response 5273

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	83.60
41.00	44.20	44.51
42.00	31.50	26.23

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 Sample : 02122-14
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.987	152		115093	20.000	ng/ul	0.00
20) Naphthalene-d8	10.810	136		498402	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.640	164		318668	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.398	188		676706	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240		587328	20.000	ng/ul	0.00
88) Perylene-d12	23.992	264		686075	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.364	96		15144	5.103	ng/uL	0.00
4) Pyridine-d5	3.817	84		41947m	5.515	ng/ul	0.02
7) Phenol-d5	7.169	99		47604m	4.485	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	7.316	67		201496	29.587	ng/ul	0.00
11) 2-Chlorophenol-d4	7.522	132		158714	20.688	ng/ul	0.00
15) 4-Methylphenol-d8	8.716	113		93098	10.873	ng/ul	0.00
21) Nitrobenzene-d5	9.169	128		109286	27.162	ng/ul	0.00
24) 2-Nitrophenol-d4	9.893	143		97871	21.243	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.440	165		171557	20.446	ng/ul	0.00
31) 4-Chloroaniline-d4	10.963	131		226648	20.507	ng/ul	0.00
46) Dimethylphthalate-d6	14.051	166		793991	30.729	ng/ul	0.00
49) Acenaphthylene-d8	14.334	160		866913	30.237	ng/ul	0.00
54) 4-Nitrophenol-d4	14.975	143		9621m	2.465	ng/ul	0.11
60) Fluorene-d10	15.634	176		663997	30.816	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.775	200		65979	13.251	ng/ul	0.01
73) Anthracene-d10	17.498	188		1098799	33.552	ng/ul	0.00
81) Pyrene-d10	19.733	212		1250188	31.690	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.827	264		1220654	33.488	ng/ul	0.00

Target Compounds Qvalue

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

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