

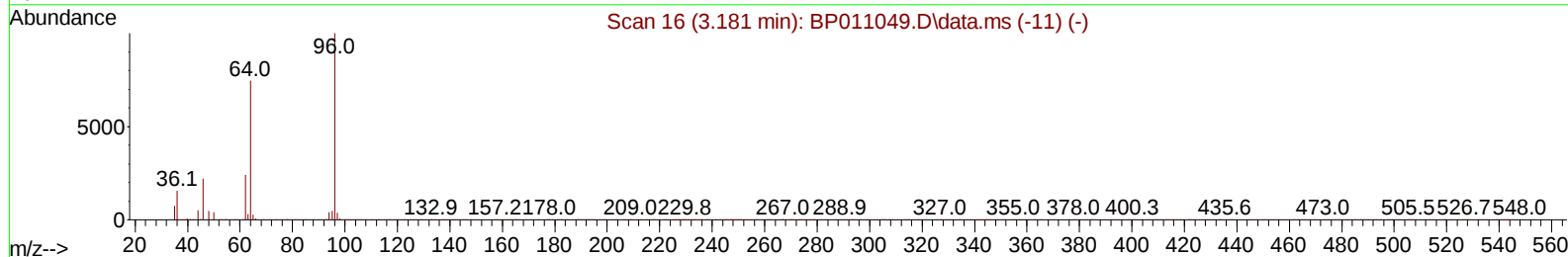
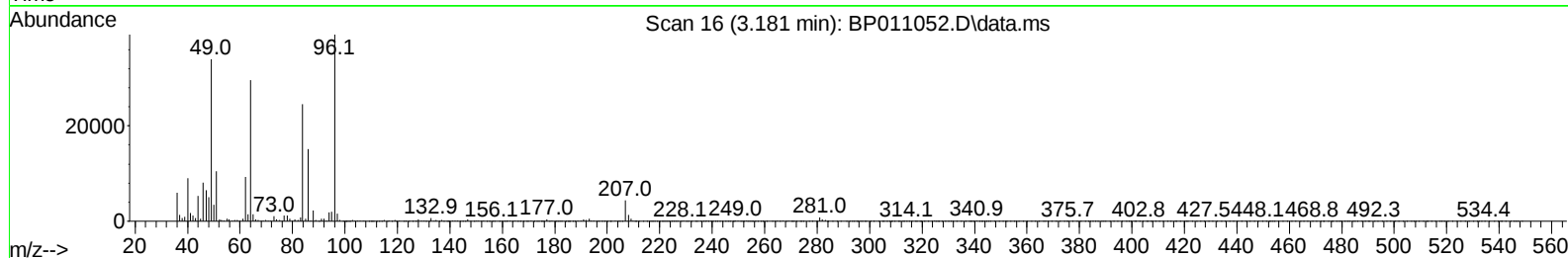
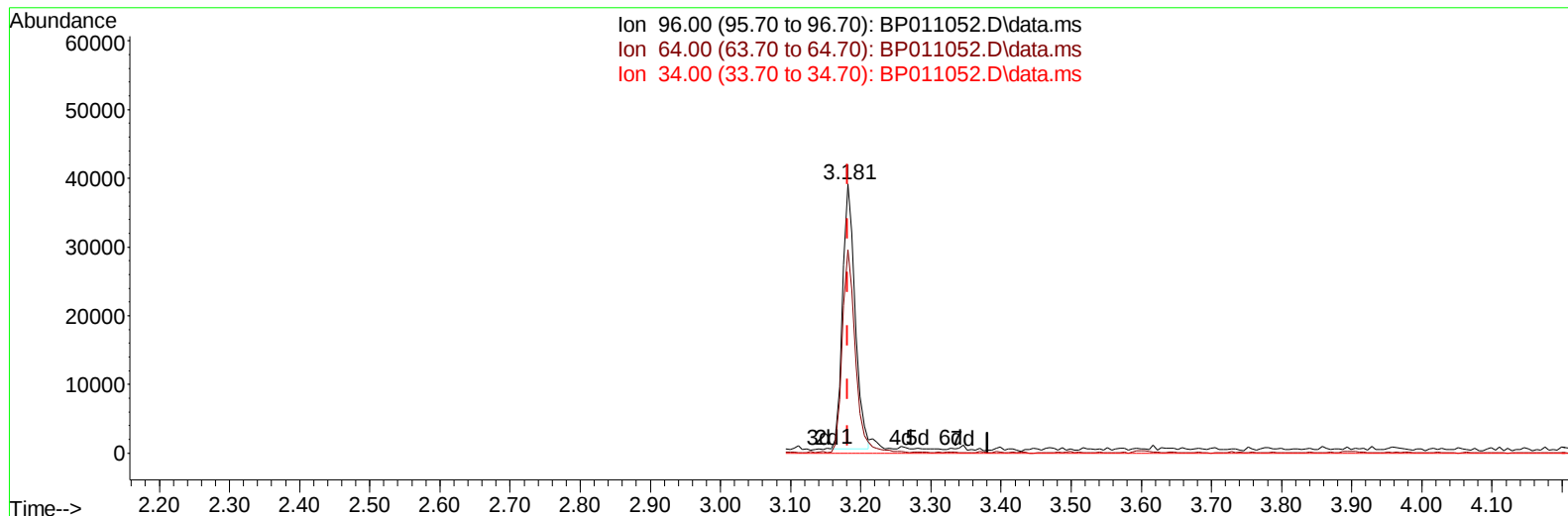
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072022\
 Data File : BP011052.D
 Acq On : 20 Jul 2022 11:10
 Operator : CG/JU
 Sample : N3753-15
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GBHQ7

Manual IntegrationsAPPROVED

Quant Time: Jul 21 00:42:41 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 19 03:25:54 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 07/21/2022
 Supervised By :mohammad ahmed 07/21/2022



TIC: BP011052.D\data.ms

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

3.181min (+ 0.000) 5.40 ng/uL

response 47930

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.90	75.54
34.00	0.00	0.00
0.00	0.00	0.00

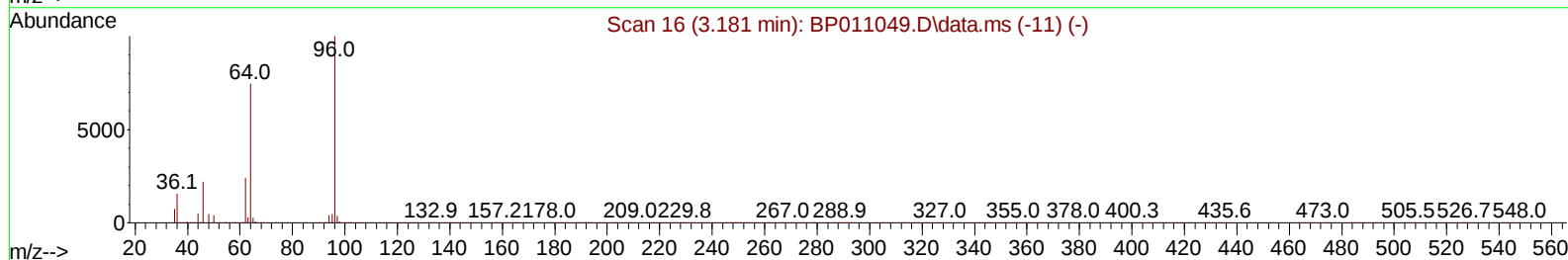
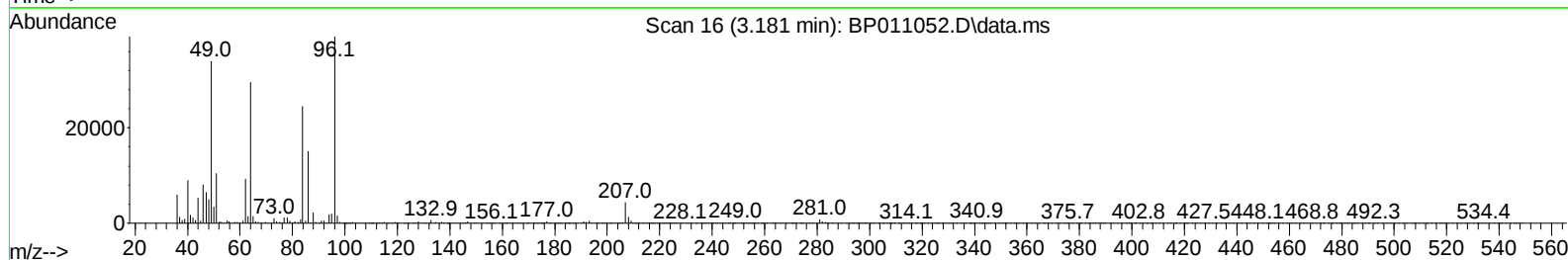
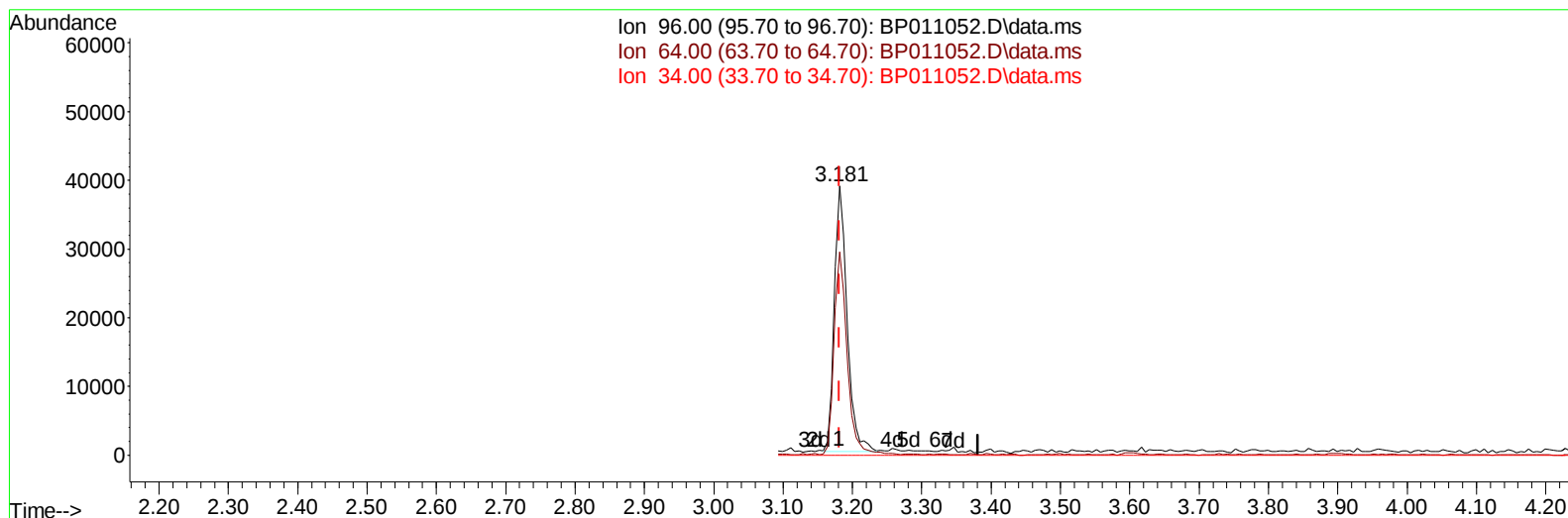
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TIC: BP011052.D\data.ms

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

3.181min (+ 0.000) 5.55 ng/uL m

response 49249

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.90	75.54
34.00	0.00	0.00
0.00	0.00	0.00

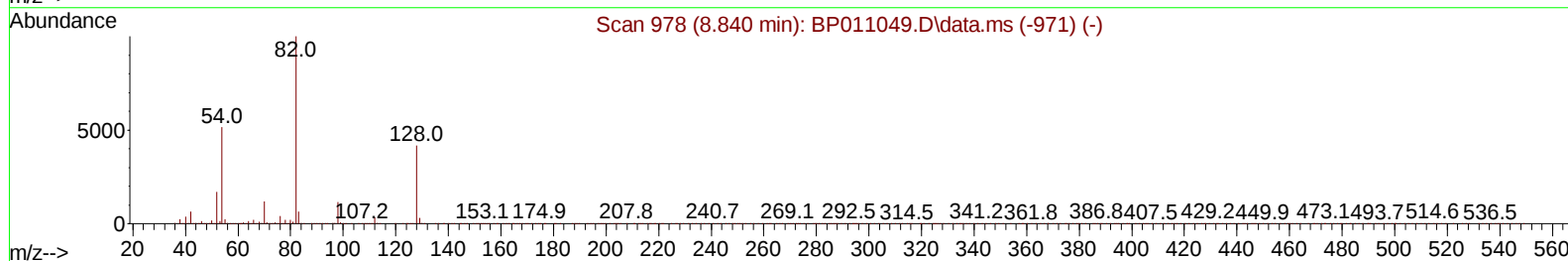
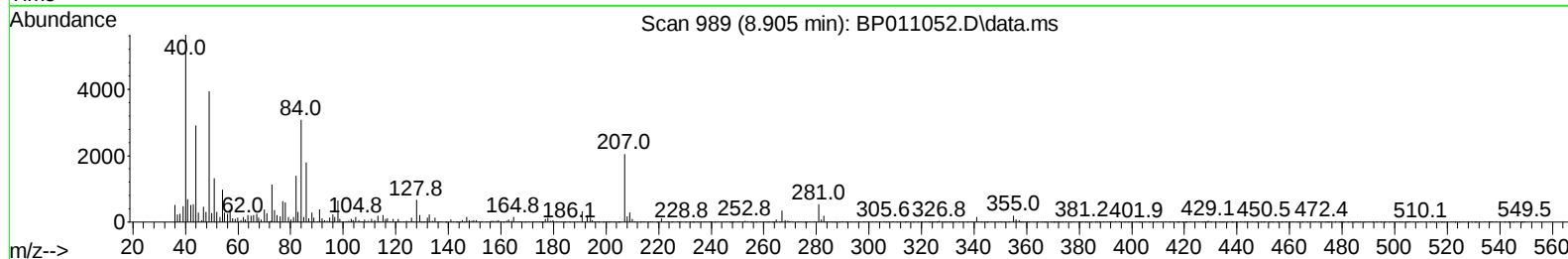
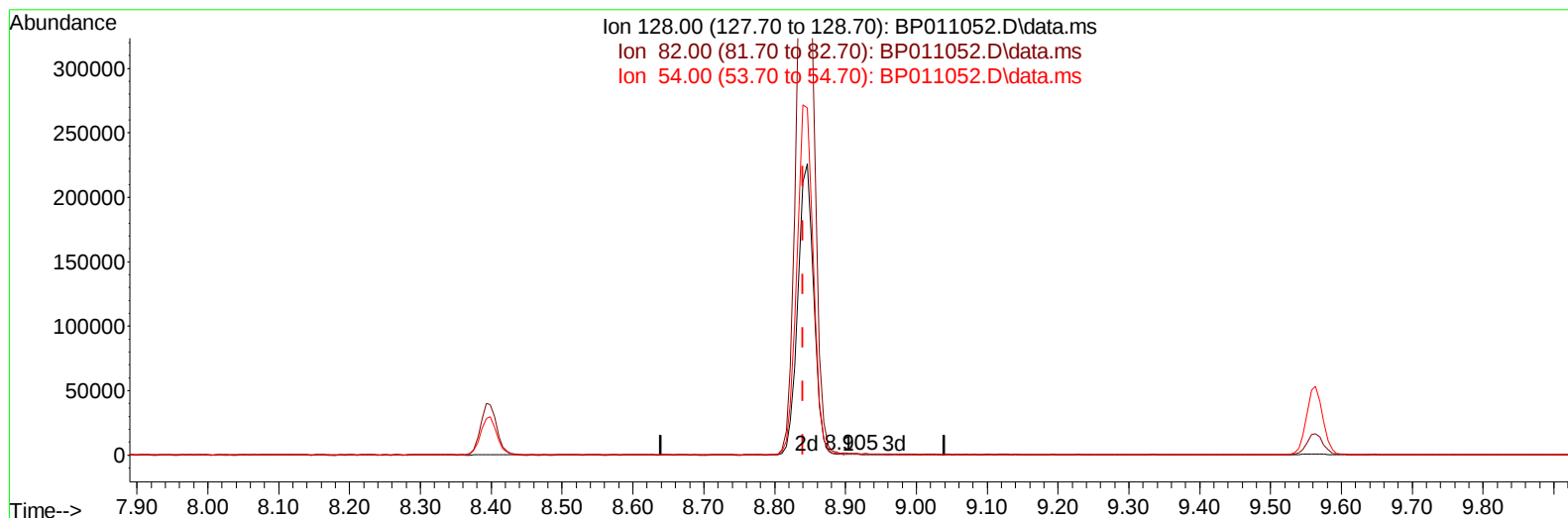
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TIC: BP011052.D\data.ms

(21) Nitrobenzene-d5 (S)

8.905min (+ 0.065) 0.06 ng/u1

response 586

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	249.70	207.92
54.00	155.90	144.69
0.00	0.00	0.00

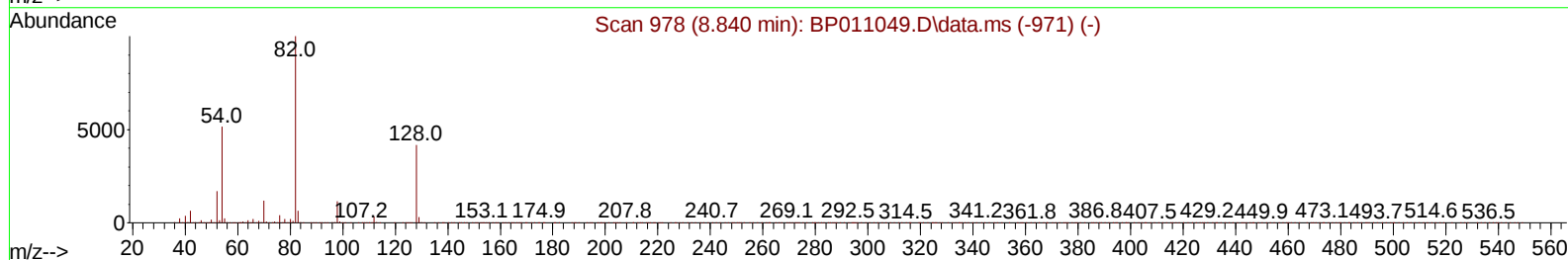
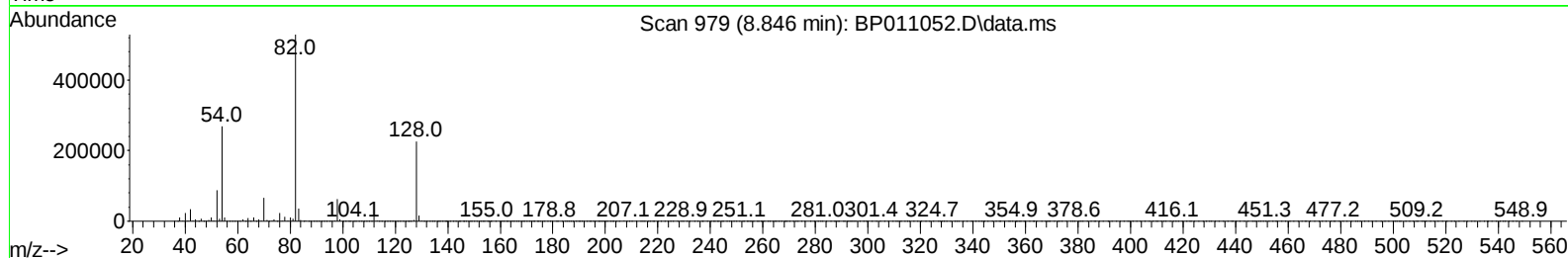
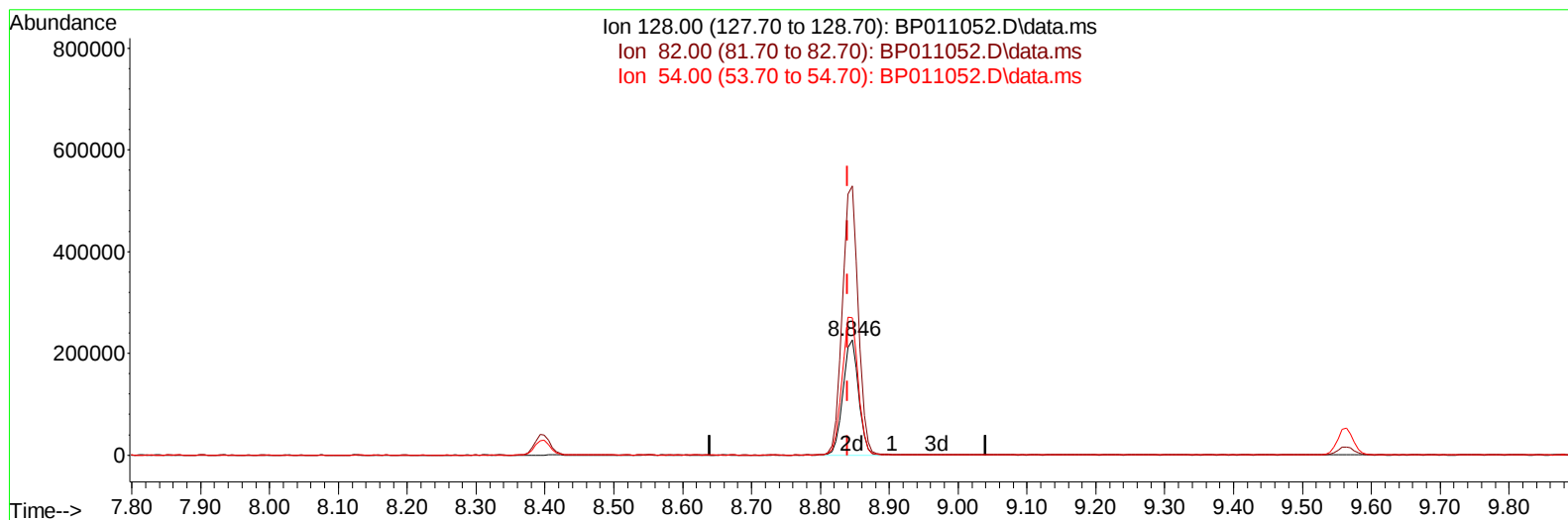
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TIC: BP011052.D\data.ms

(21) Nitrobenzene-d5 (S)

8.846min (+ 0.006) 36.84 ng/ul m

response 355059

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	249.70	234.12
54.00	155.90	119.30#
0.00	0.00	0.00

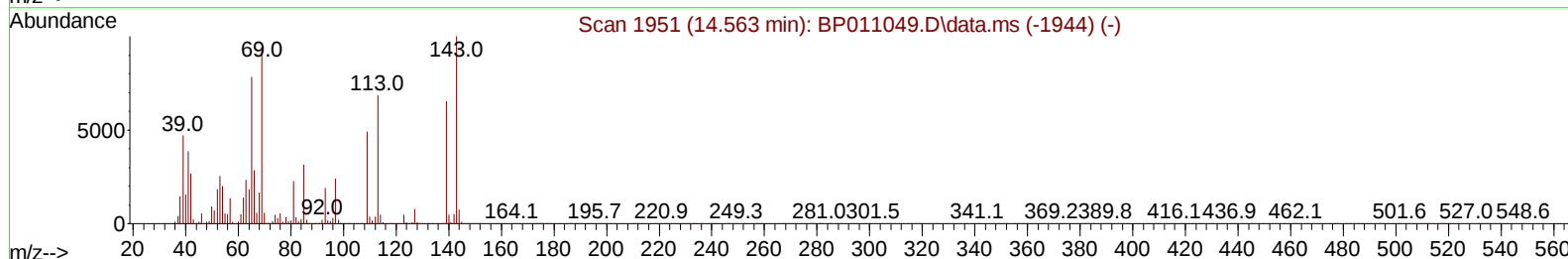
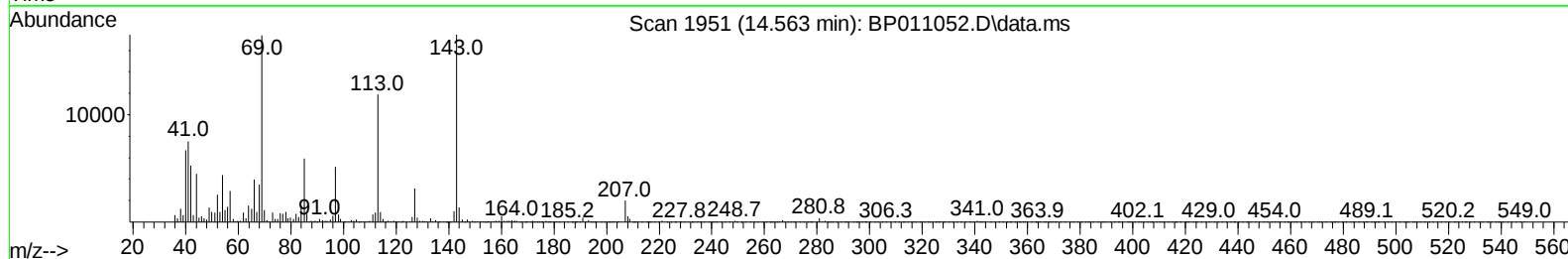
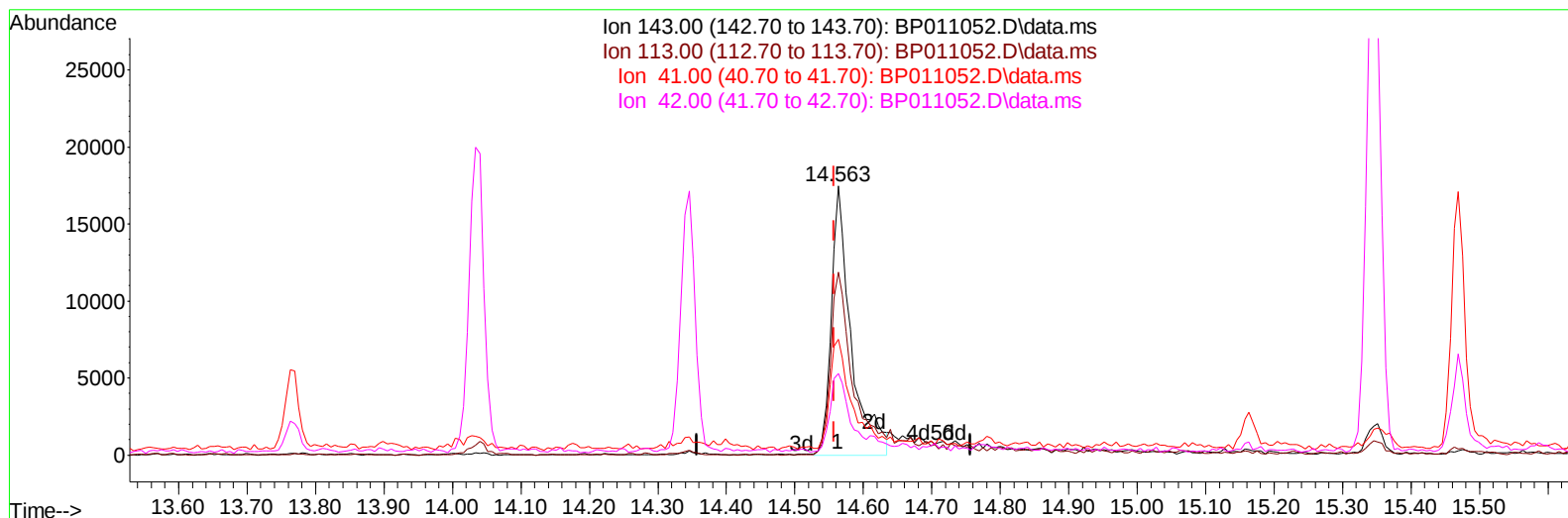
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 ALS Vial : 5 Sample Multiplier: 1

Instrument :
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TIC: BP011052.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.563min (+ 0.006) 3.40 ng/ul m

response 35057

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.50	68.06#
41.00	59.00	43.10#
42.00	41.10	30.22#

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 Sample : N3753-15
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.681	152	320095	20.000	ng/ul	0.00
20) Naphthalene-d8	10.475	136	1322523	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.346	164	744691	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.116	188	1498217	20.000	ng/ul	# 0.01
79) Chrysene-d12	21.233	240	1436023	20.000	ng/ul	# 0.00
88) Perylene-d12	23.592	264	1454371	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	49249m	5.547	ng/uL	0.00
4) Pyridine-d5	3.599	84	159704	6.959	ng/ul	0.00
7) Phenol-d5	6.858	99	178019	6.775	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.022	67	606877	35.918	ng/ul	0.00
11) 2-Chlorophenol-d4	7.217	132	558912	26.325	ng/ul	0.00
15) 4-Methylphenol-d8	8.393	113	341571	16.238	ng/ul	0.00
21) Nitrobenzene-d5	8.846	128	355059m	36.840	ng/ul	0.00
24) 2-Nitrophenol-d4	9.563	143	307598	32.254	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.099	165	524765	30.089	ng/ul	0.00
31) 4-Chloroaniline-d4	10.622	131	768660	26.734	ng/ul	0.00
46) Dimethylphthalate-d6	13.769	166	1968280	38.672	ng/ul	0.00
49) Acenaphthylene-d8	14.034	160	2348732	39.112	ng/ul	0.00
54) 4-Nitrophenol-d4	14.563	143	35057m	3.403	ng/ul	0.00
60) Fluorene-d10	15.346	176	1676736	38.757	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.469	200	200444	27.037	ng/ul	0.00
73) Anthracene-d10	17.216	188	2861808	44.931	ng/ul	0.01
81) Pyrene-d10	19.469	212	3276810	43.447	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.445	264	3142342	44.986	ng/ul	0.02

Target Compounds Qvalue

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

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