

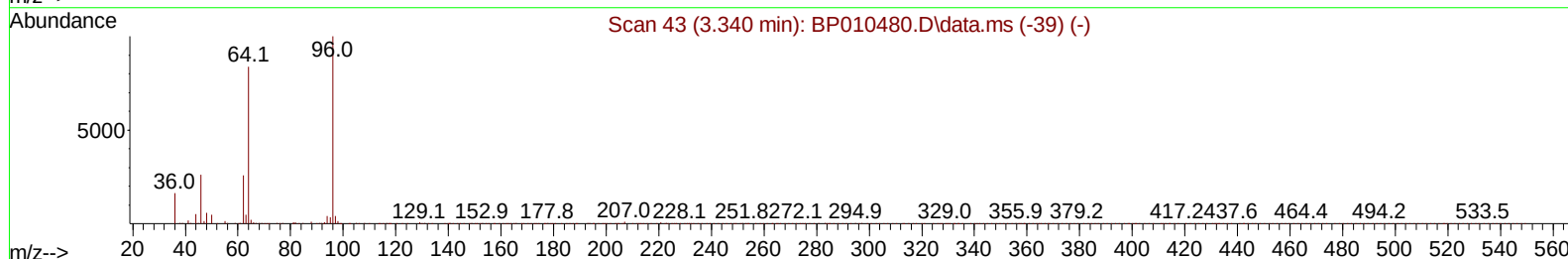
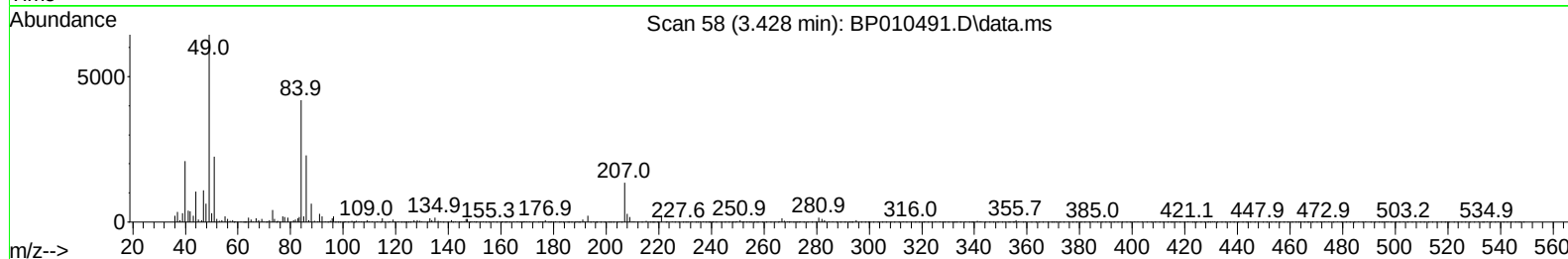
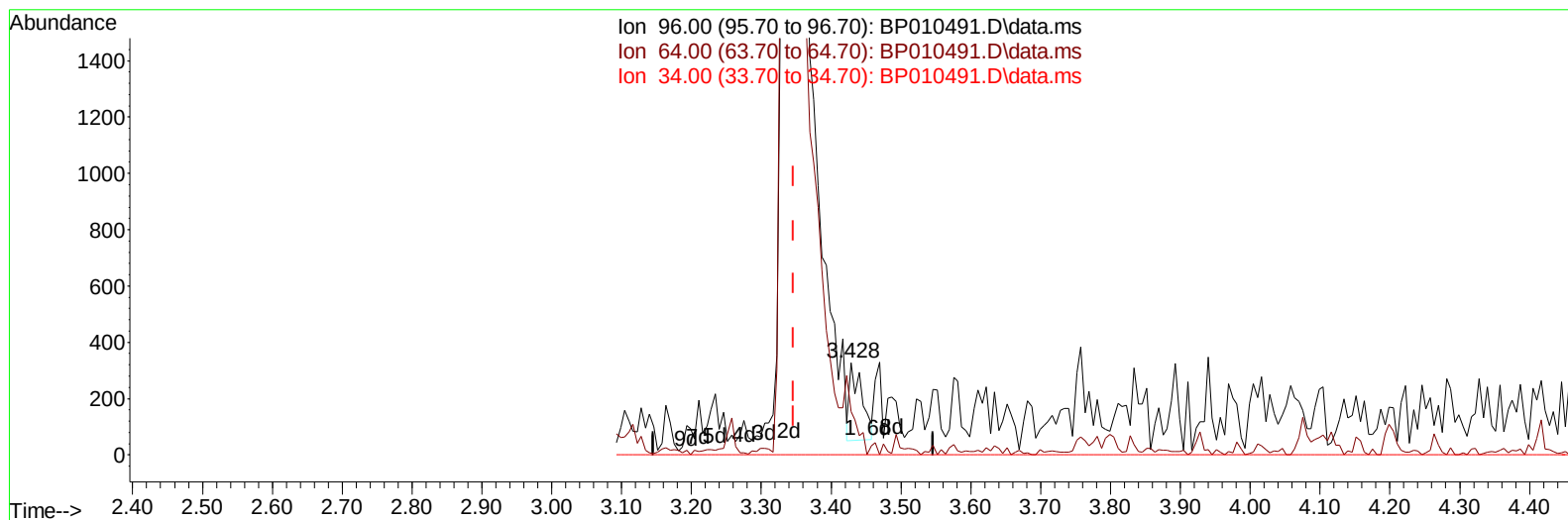
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052322\
 Data File : BP010491.D
 Acq On : 23 May 2022 23:24
 Operator : CG/JU
 Sample : N2955-03
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AA2

Manual IntegrationsAPPROVED

Quant Time: May 24 00:52:43 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri May 13 15:04:31 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/24/2022
 Supervised By :Yogesh Patel 05/26/2022



TIC: BP010491.D\data.ms

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

3.428min (+ 0.082) 0.10 ng/uL

response 336

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	46.95
34.00	0.00	0.00
0.00	0.00	0.00

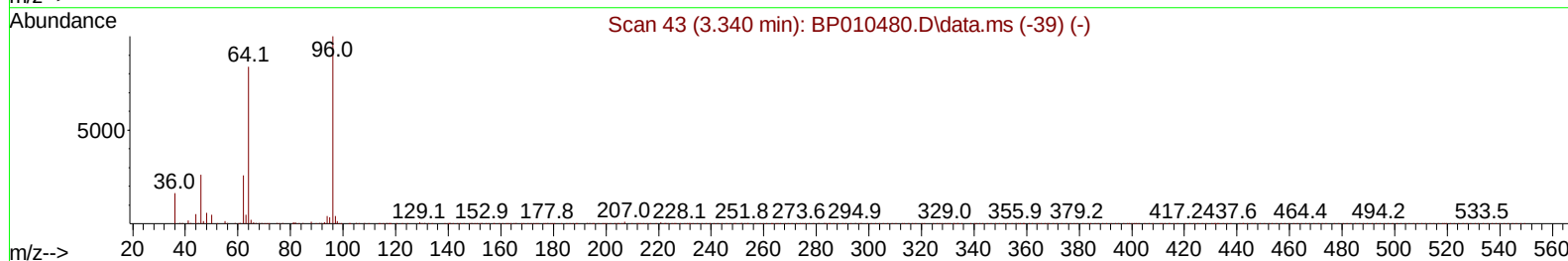
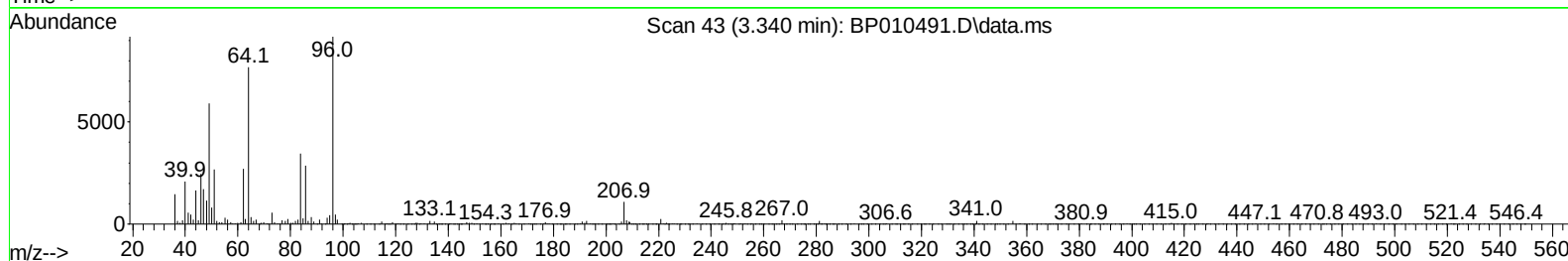
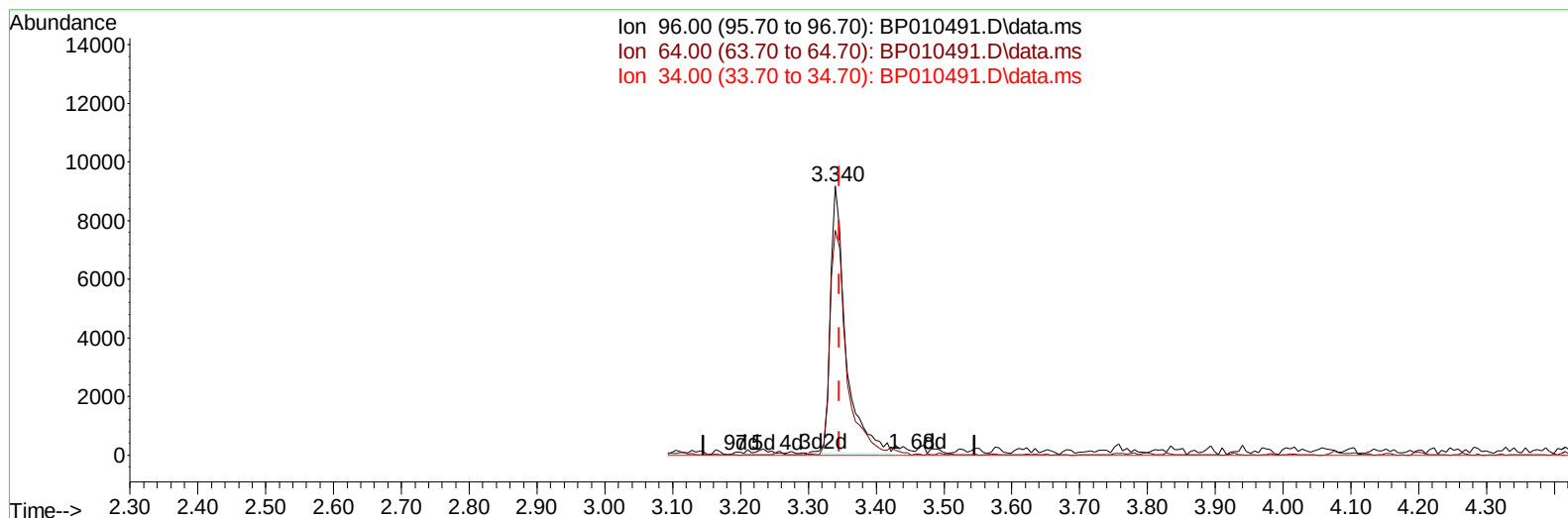
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(3) 1,4-Dioxane-d8 (S)

3.340min (-0.006) 4.31 ng/uL m

response 14376

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	83.58#
34.00	0.00	0.00
0.00	0.00	0.00

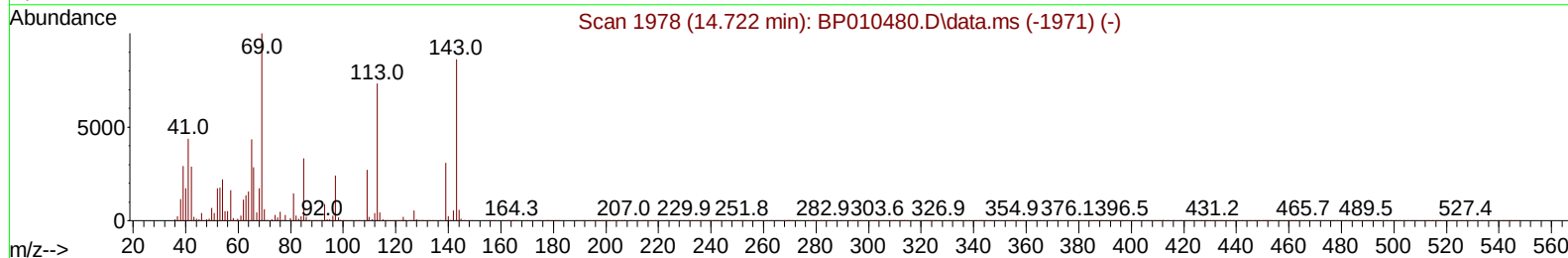
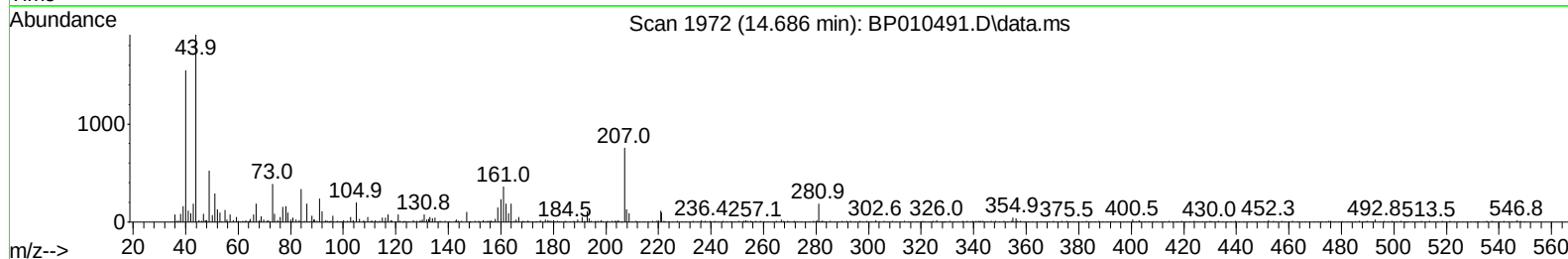
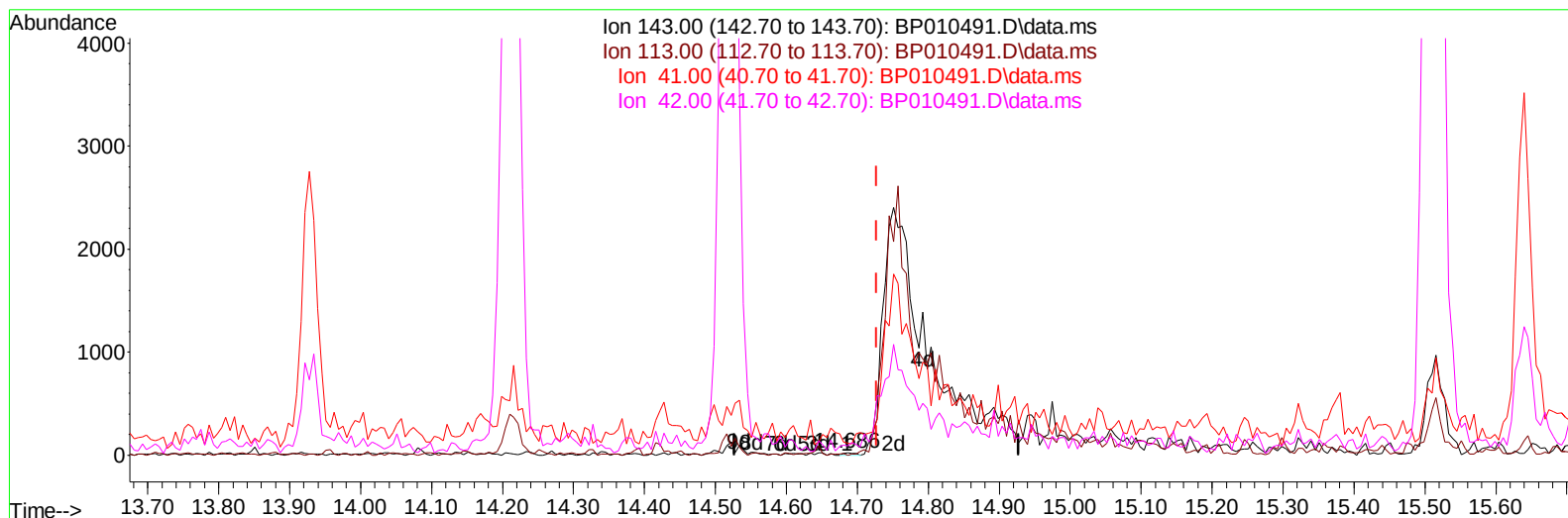
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TIC: BP010491.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.686min (-0.041) 0.00 ng/ul

response 24

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	0.00#
41.00	23.20	487.50#
42.00	18.00	366.67#

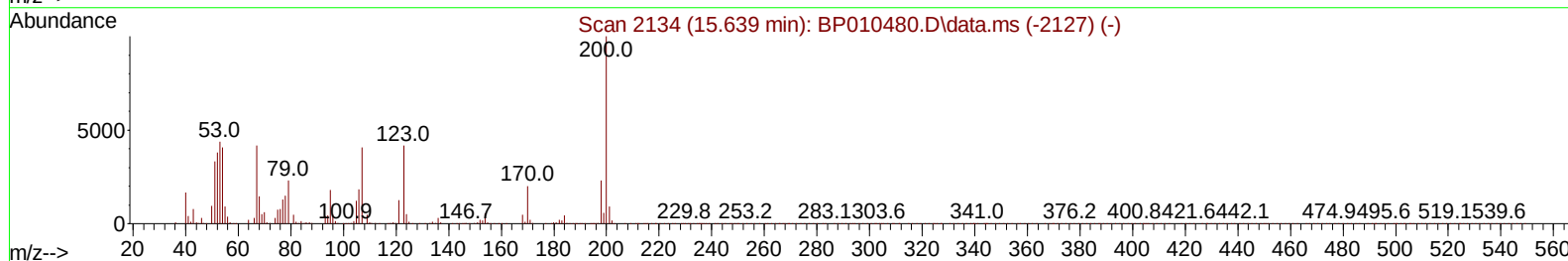
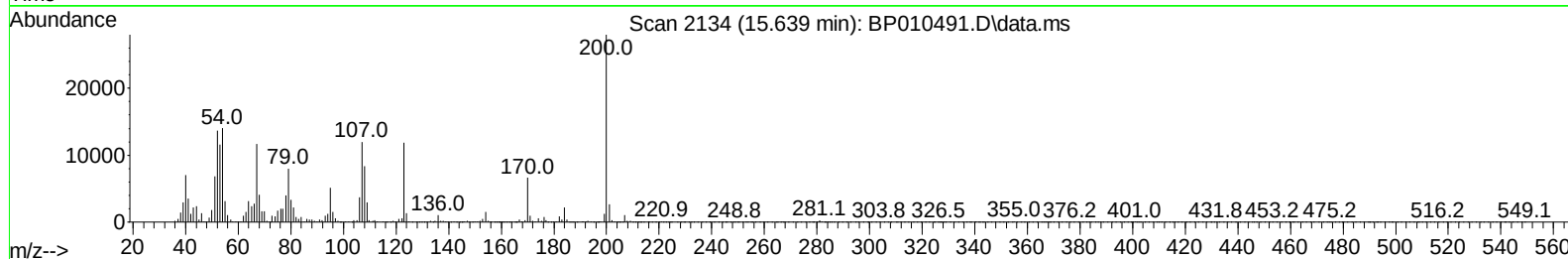
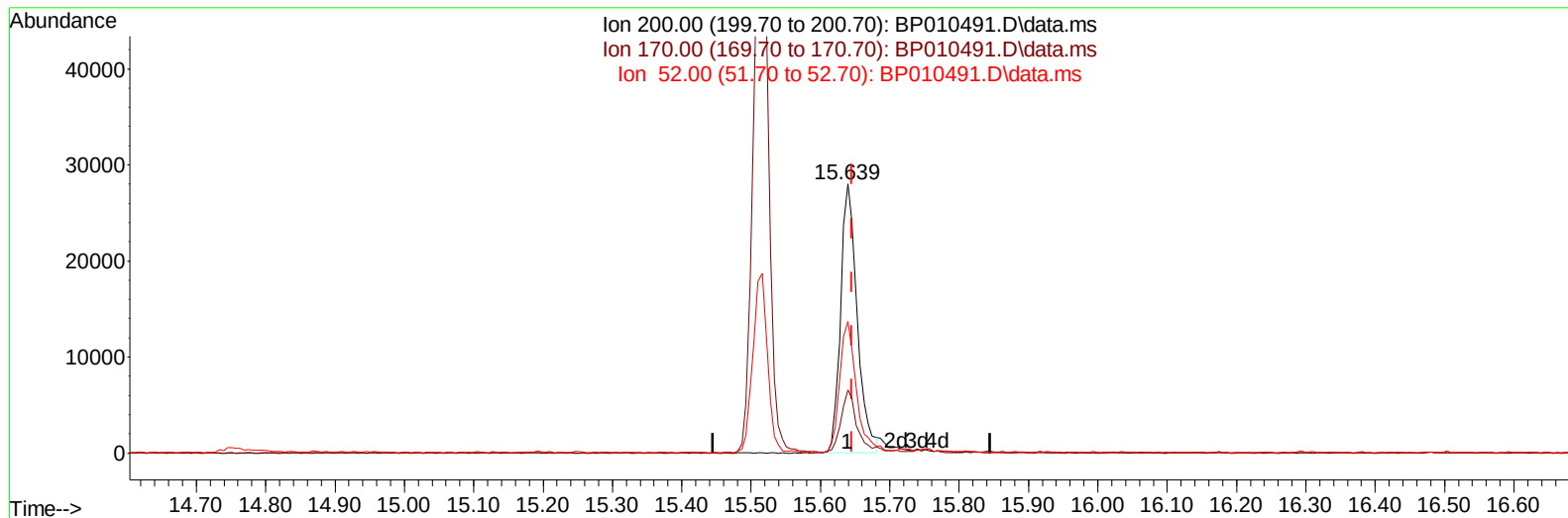
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TIC: BP010491.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.639min (-0.006) 8.46 ng/ul

response 46048

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.00	23.57
52.00	21.20	48.91#
0.00	0.00	0.00

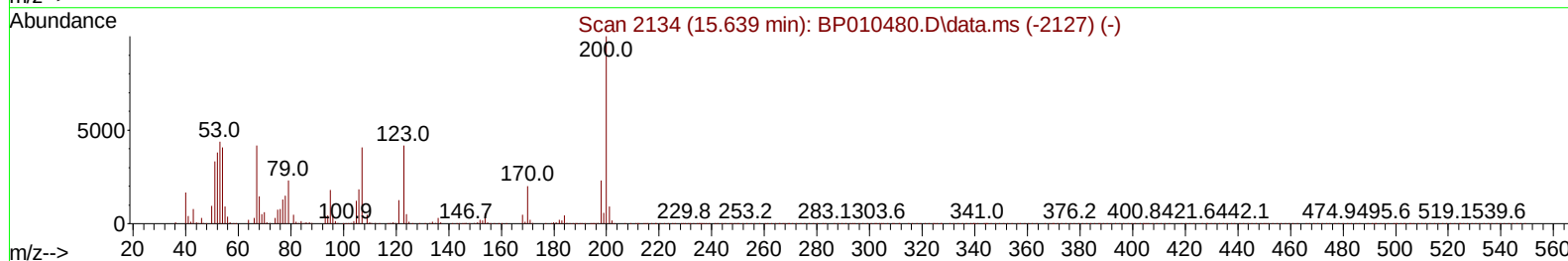
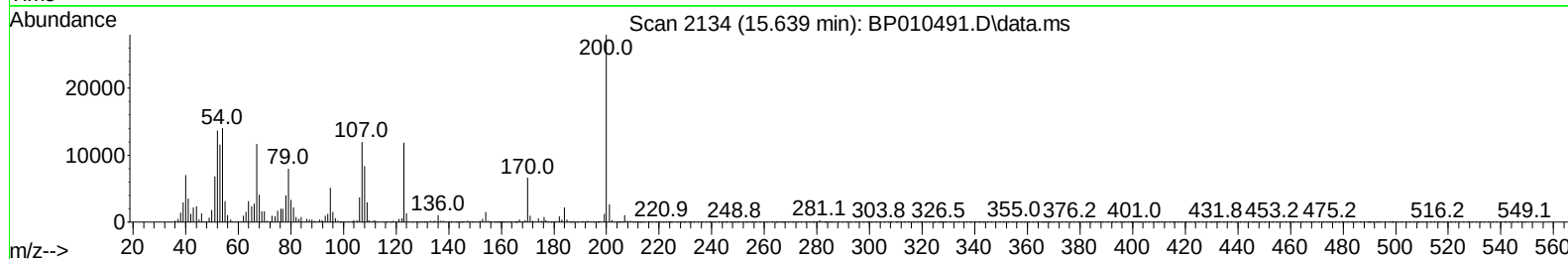
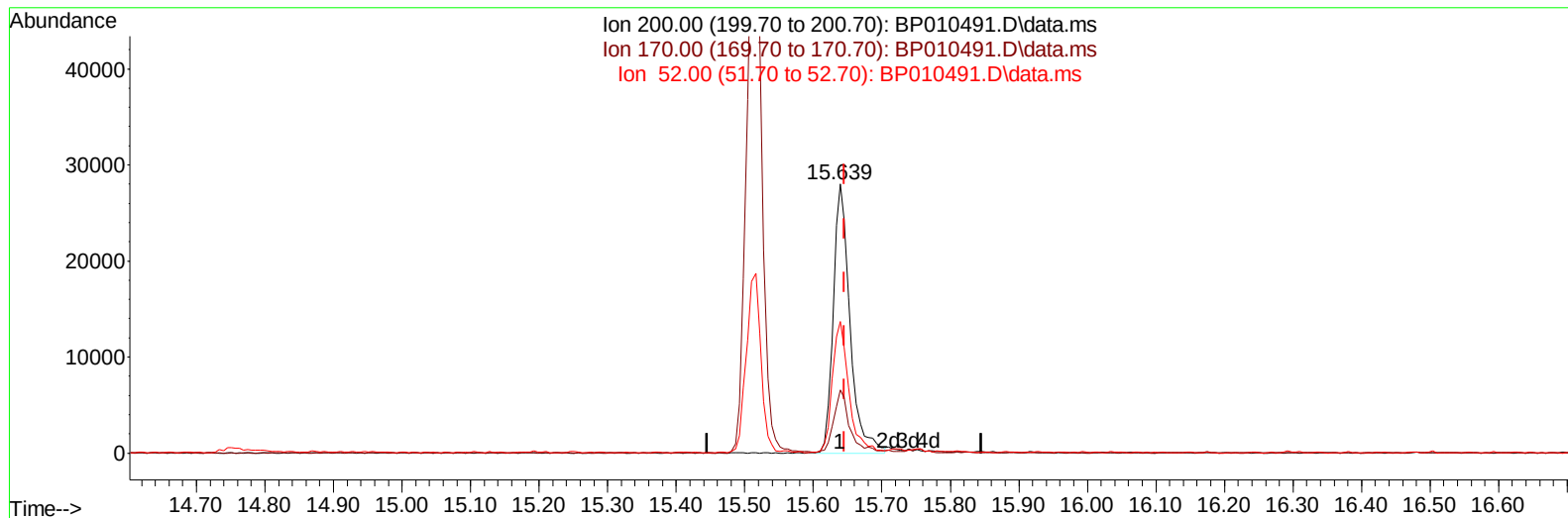
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Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AA2

Manual IntegrationsAPPROVED

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TIC: BP010491.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.639min (-0.006) 8.71 ng/ul m

response 47423

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.00	23.57
52.00	21.20	48.91#
0.00	0.00	0.00

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 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
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Manual IntegrationsAPPROVED

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	142057	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.687	136	644850	20.000	ng/ul	#-0.02
38) Acenaphthene-d10	14.522	164	441451	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.269	188	930513	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.351	240	771378	20.000	ng/ul	#-0.01
88) Perylene-d12	23.774	264	603701	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	14376m	4.314	ng/uL	0.00
4) Pyridine-d5	3.769	84	65439	6.915	ng/ul	0.00
7) Phenol-d5	7.057	99	56262	4.719	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.216	67	208675	26.049	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.422	132	179675	19.694	ng/ul	-0.01
15) 4-Methylphenol-d8	8.599	113	119846	11.828	ng/ul	-0.01
21) Nitrobenzene-d5	9.046	128	131993	26.448	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.769	143	134395	25.410	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.310	165	222615	22.260	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.828	131	323330	21.744	ng/ul	-0.01
46) Dimethylphthalate-d6	13.928	166	877996	26.995	ng/ul	-0.01
49) Acenaphthylene-d8	14.210	160	1019340	27.171	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.751	143	10045m	1.735	ng/ul	0.02
60) Fluorene-d10	15.516	176	779274	27.540	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.639	200	47423m	8.714	ng/ul	0.00
73) Anthracene-d10	17.369	188	1322319	31.383	ng/ul	-0.01
81) Pyrene-d10	19.604	212	1551199	37.866	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.615	264	985667	32.515	ng/ul	-0.02

Target Compounds Qvalue

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

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