

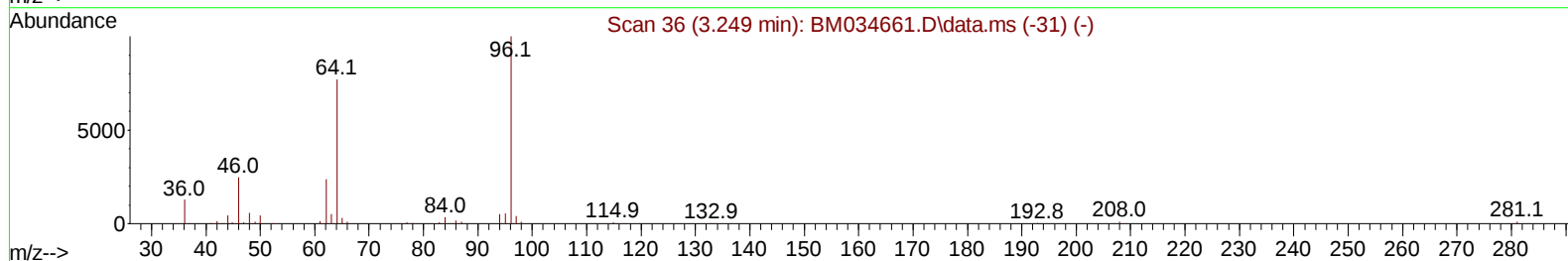
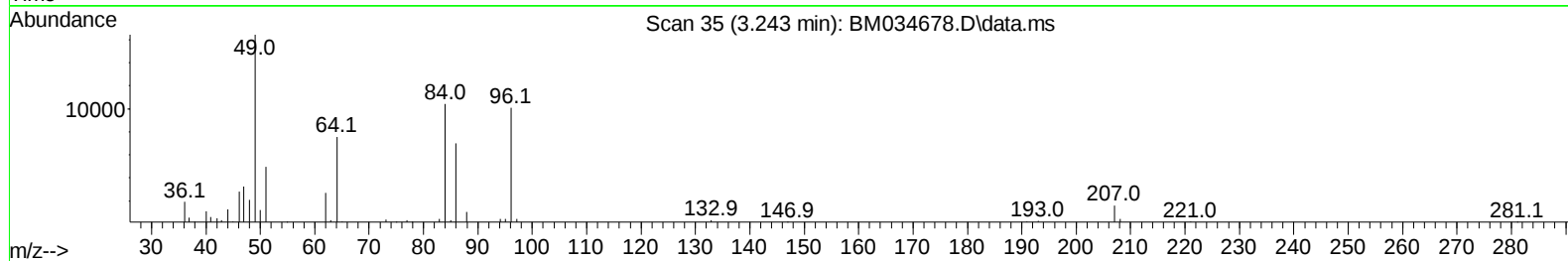
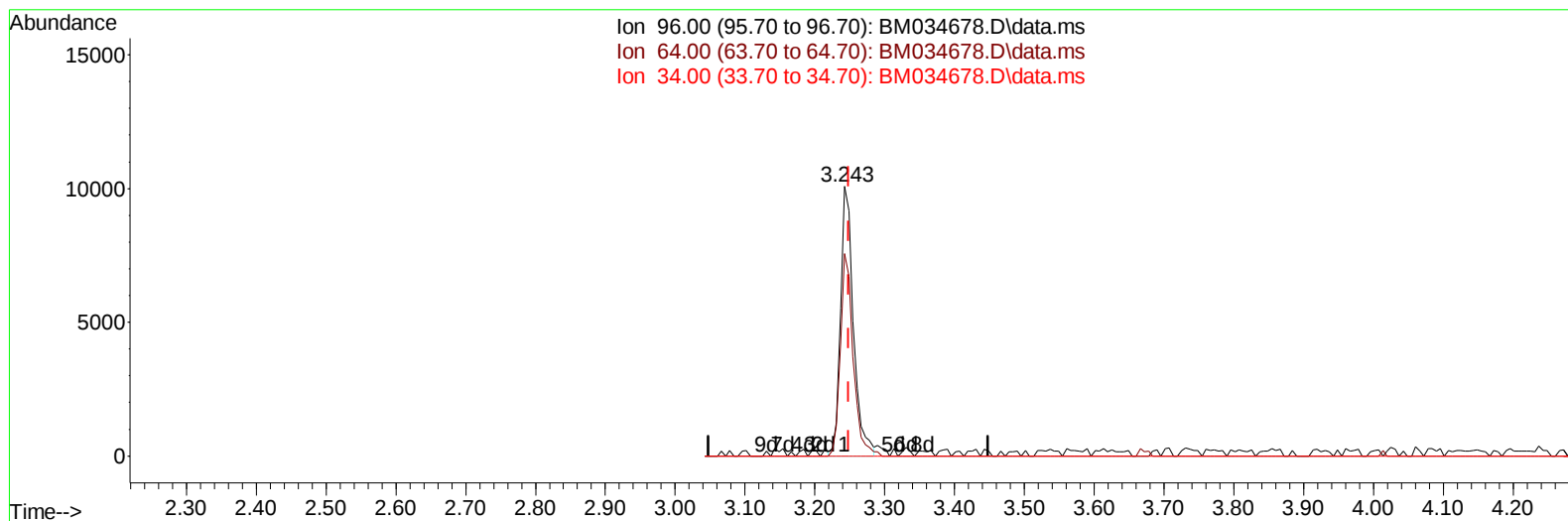
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
Data File : BM034678.D
Acq On : 14 Apr 2022 21:38
Operator : CG/JU
Sample : N2316-05
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETNQ3

Manual IntegrationsAPPROVED

Quant Time: Apr 14 23:45:05 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 14 23:42:00 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/15/2022
Supervised By :Yogesh Patel 04/23/2022



TIC: BM034678.D\data.ms

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

3.243min (-0.006) 4.08 ng/uL

response 12922

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	68.90	75.01
34.00	0.00	0.00
0.00	0.00	0.00

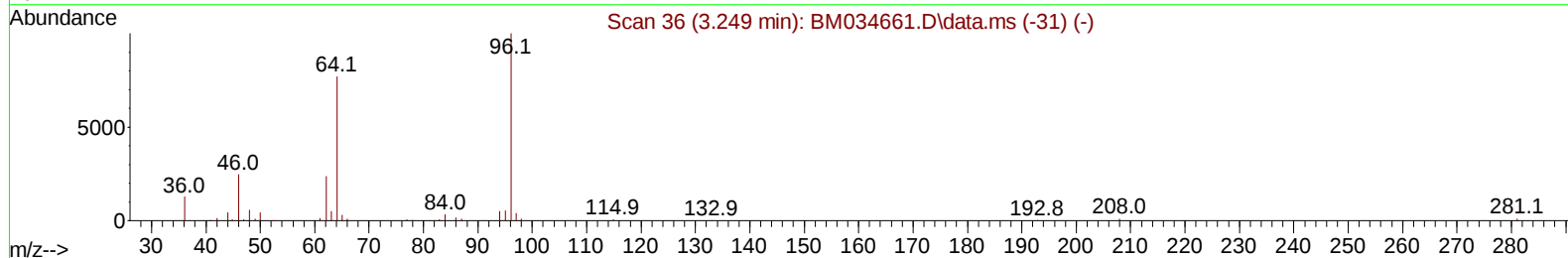
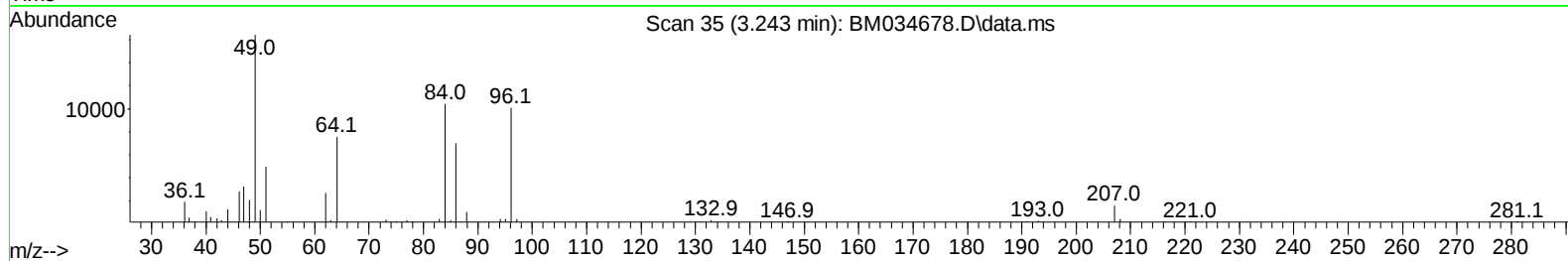
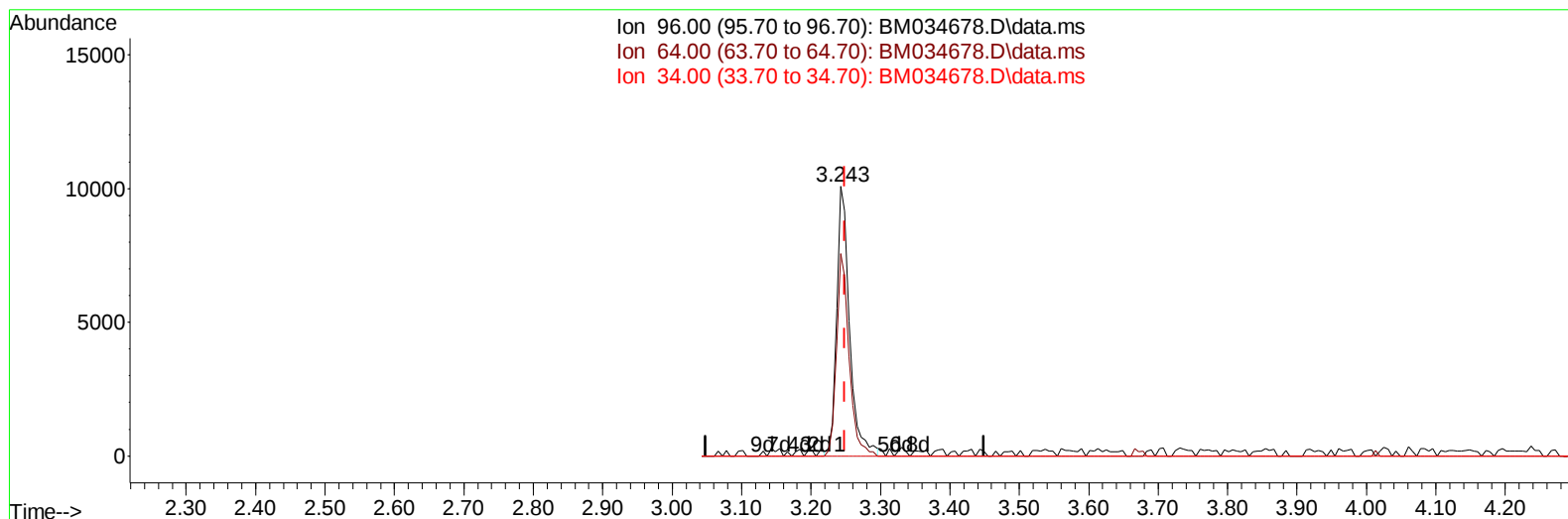
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TIC: BM034678.D\data.ms

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

3.243min (-0.006) 4.16 ng/uL m

response 13158

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	68.90	75.01
34.00	0.00	0.00
0.00	0.00	0.00

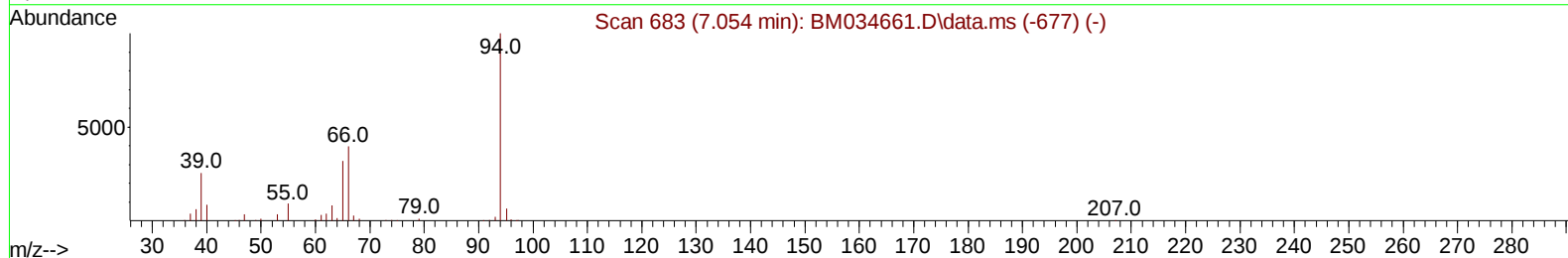
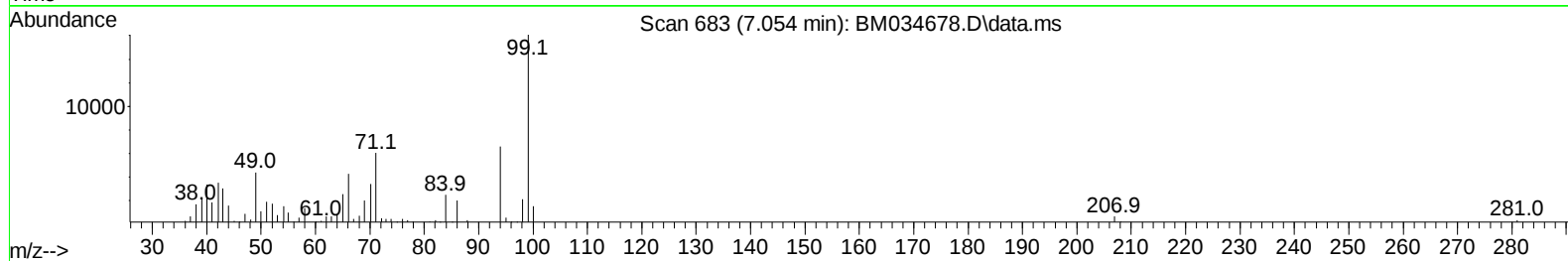
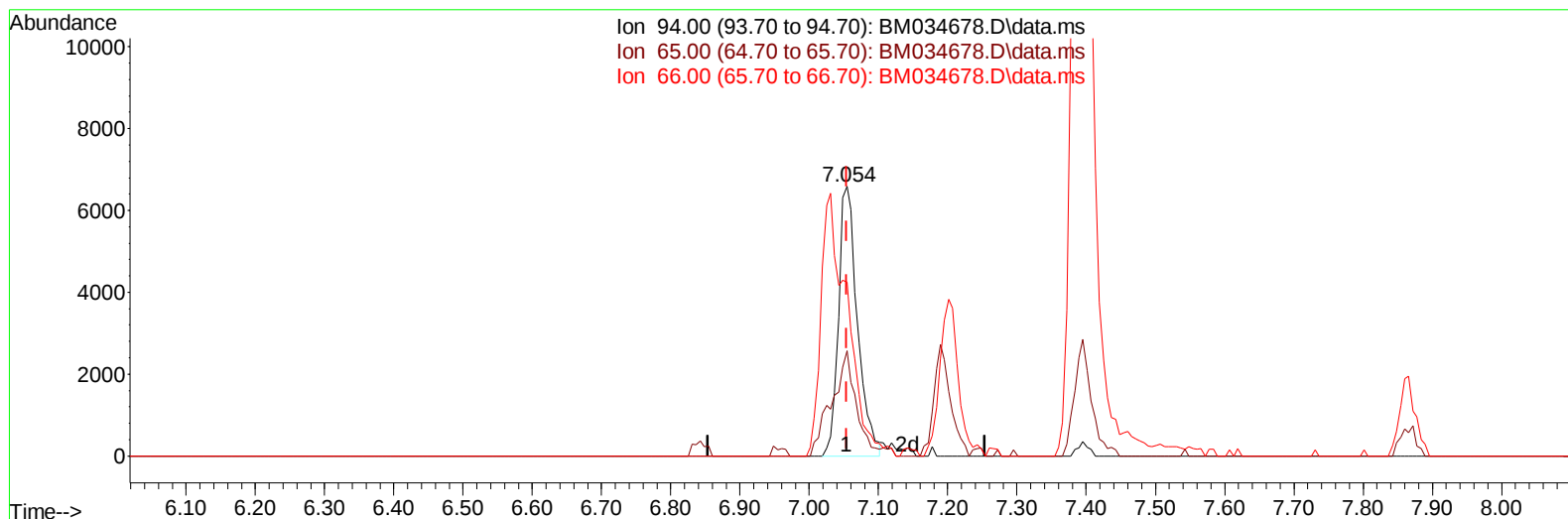
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TIC: BM034678.D\data.ms

(8) Phenol

7.054min (+ 0.000) 1.32 ng/u1

response 12525

Ion	Exp%	Act%
94.00	100.00	100.00
65.00	28.60	38.98#
66.00	42.10	64.51#
0.00	0.00	0.00

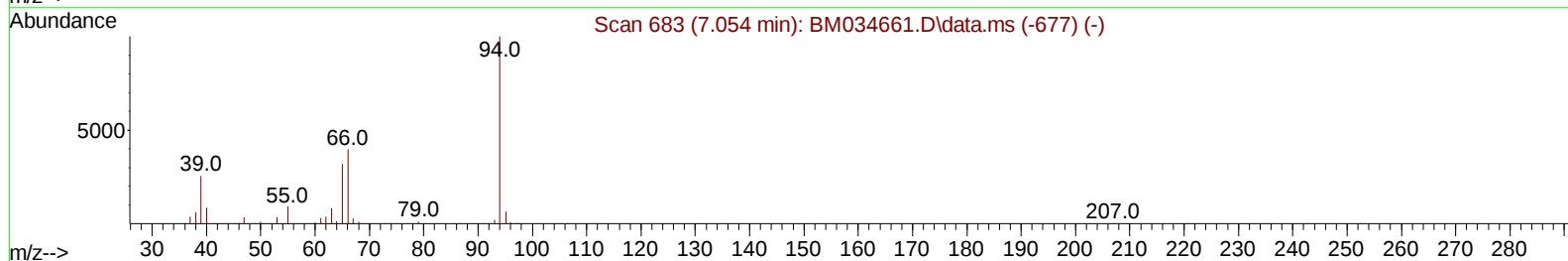
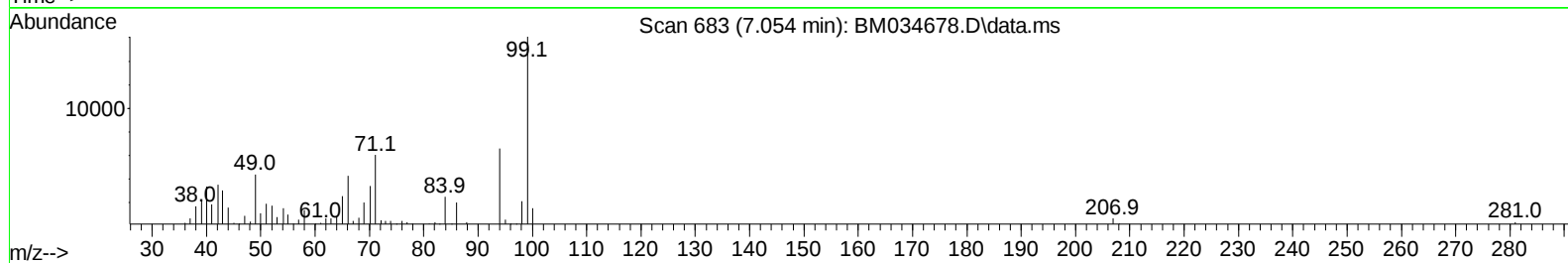
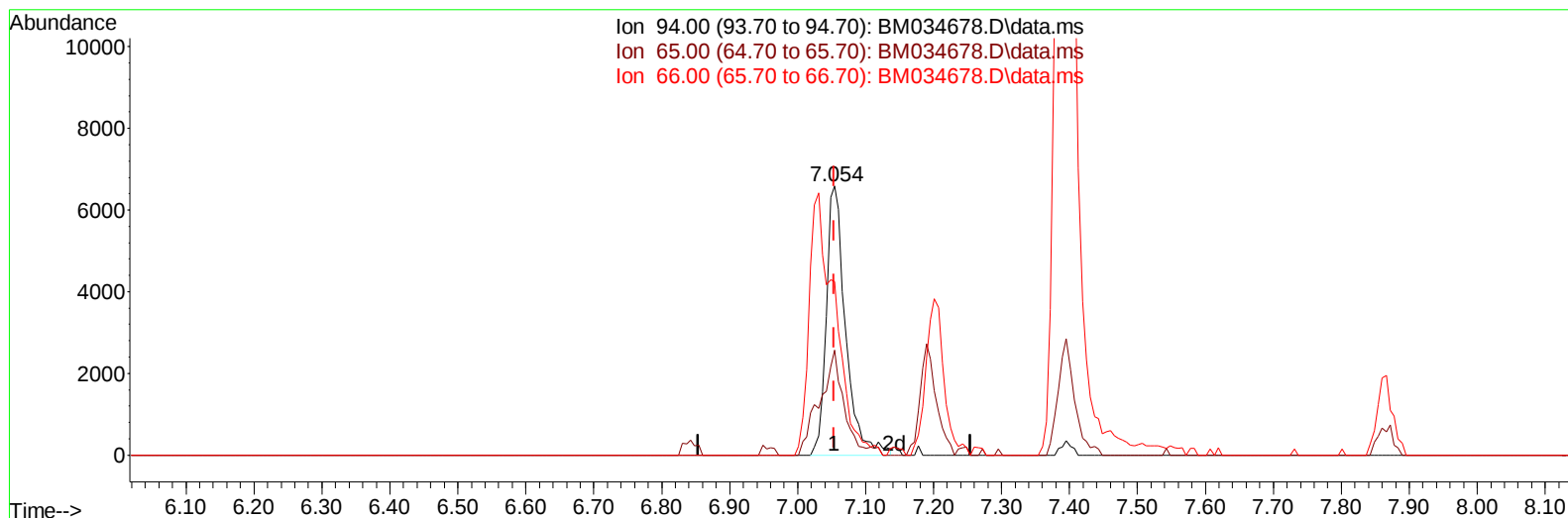
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Instrument :
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Manual IntegrationsAPPROVED

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TIC: BM034678.D\data.ms

(8) Phenol

7.054min (+ 0.000) 1.37 ng/ul m

response 12985

Ion	Exp%	Act%
94.00	100.00	100.00
65.00	28.60	38.98#
66.00	42.10	64.51#
0.00	0.00	0.00

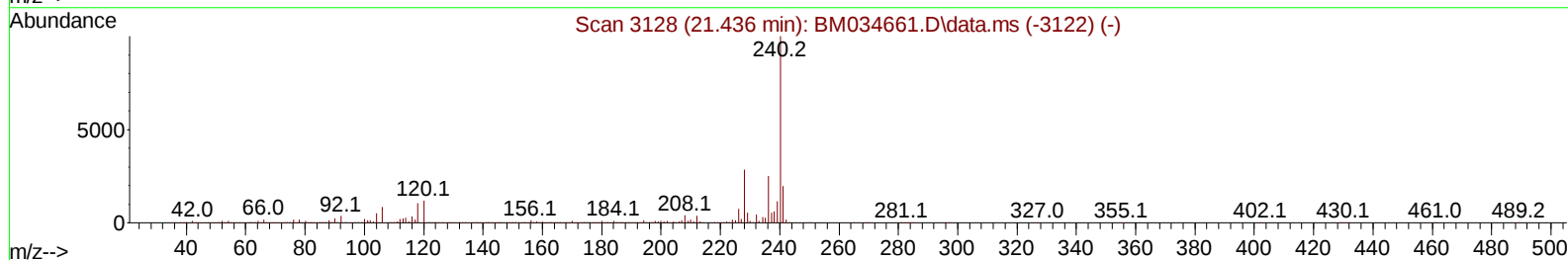
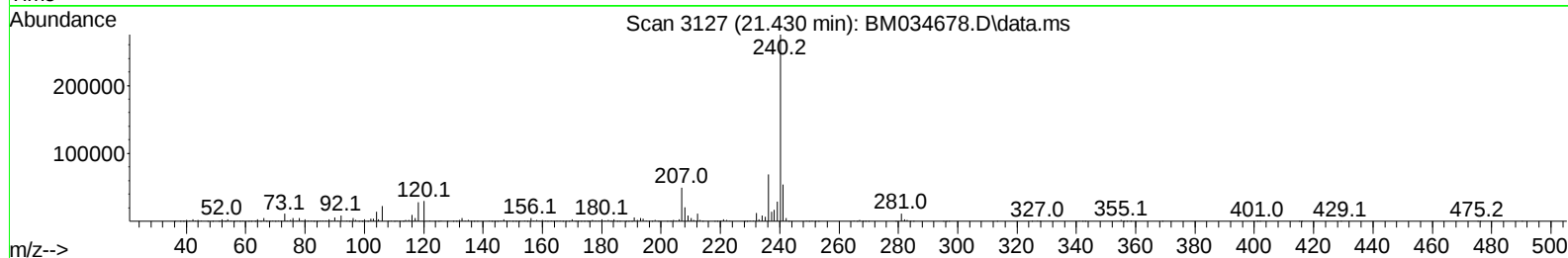
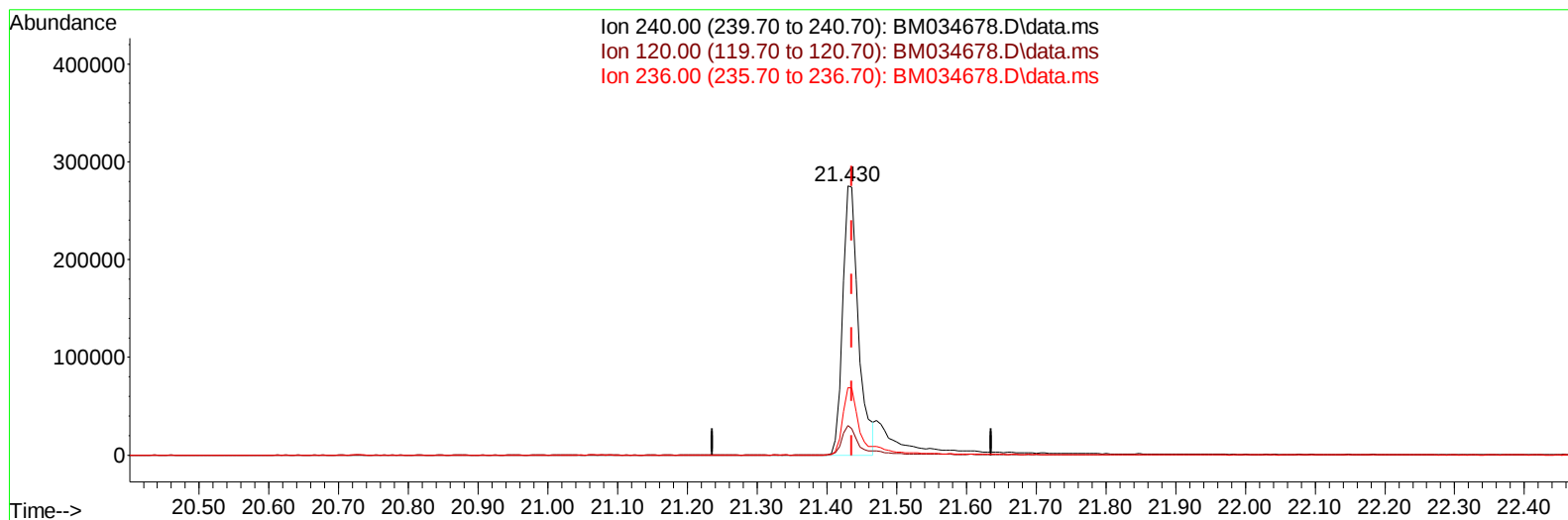
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TIC: BM034678.D\data.ms

(79) Chrysene-d12 (I)

21.430min (-0.006) 20.00 ng/u1

response 426173

Ion	Exp%	Act%
240.00	100.00	100.00
120.00	7.90	10.89#
236.00	24.80	25.13
0.00	0.00	0.00

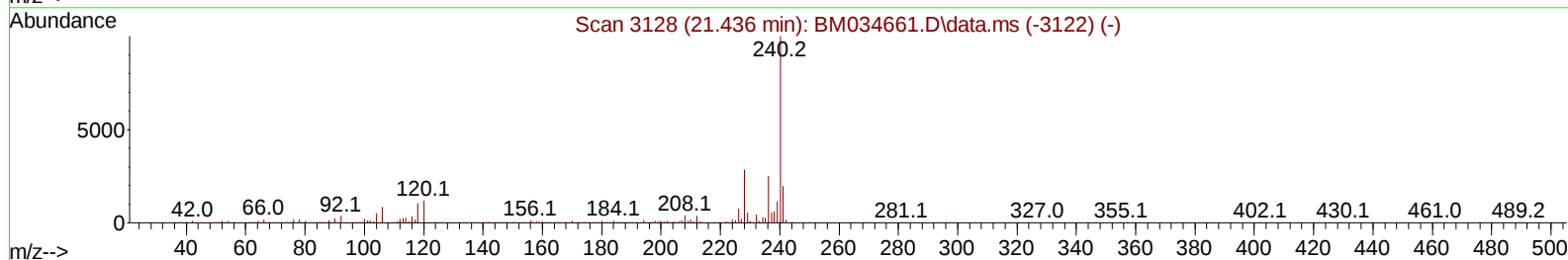
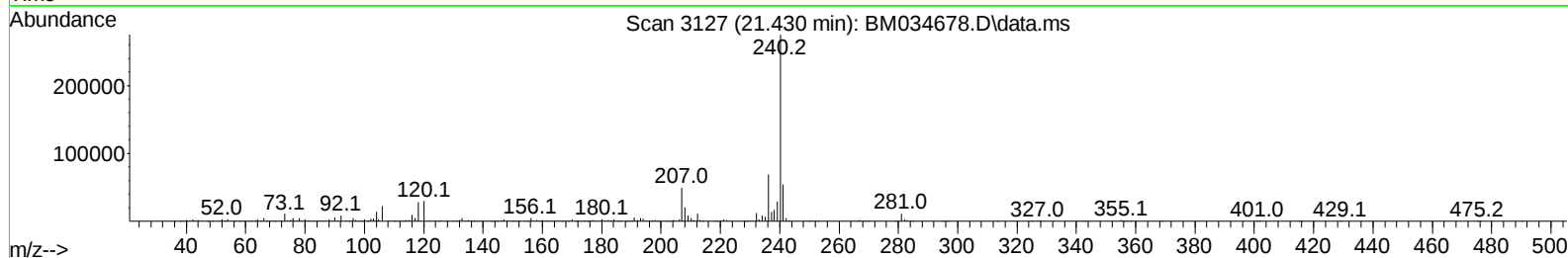
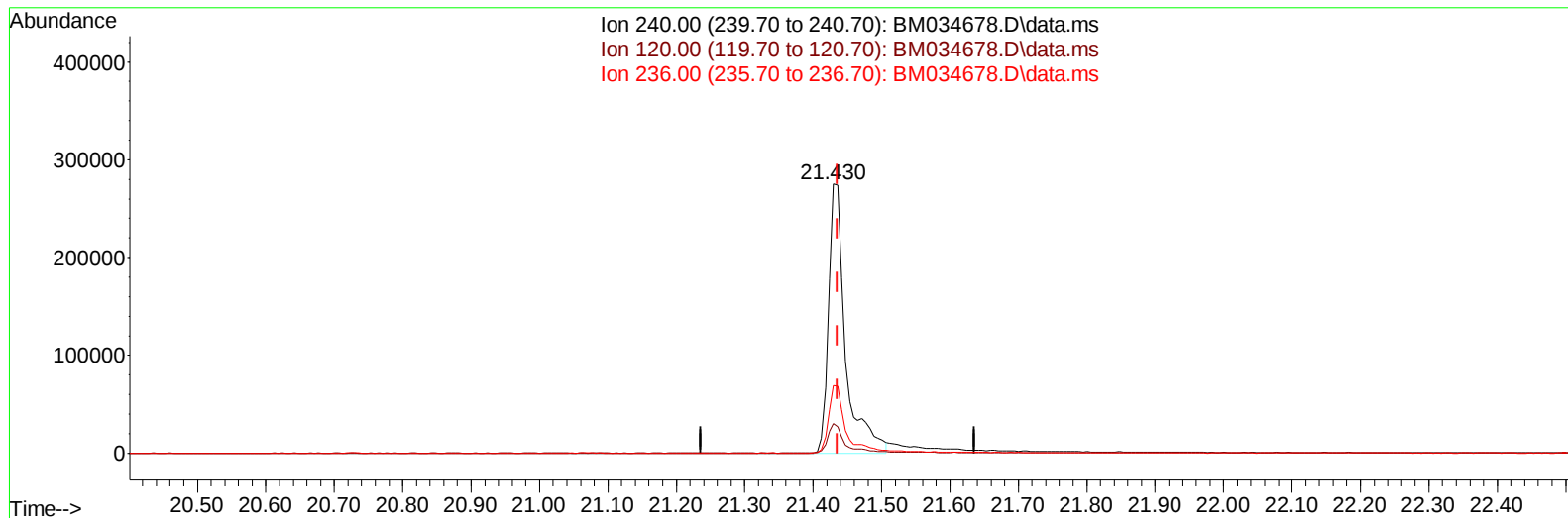
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Operator : CG/JU
Sample : N2316-05
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETNQ3

Manual IntegrationsAPPROVED

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TIC: BM034678.D\data.ms

(79) Chrysene-d12 (I)

21.430min (-0.006) 20.00 ng/ul m

response 478354

Ion	Exp%	Act%
240.00	100.00	100.00
120.00	7.90	10.89#
236.00	24.80	25.13
0.00	0.00	0.00

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 ALS Vial : 20 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	7.866	152		112890	20.000	ng/ul	0.00
20) Naphthalene-d8	10.672	136		462026	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.513	164		272201	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.254	188		538077	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.430	240		478354m	20.000	ng/ul	0.00
88) Perylene-d12	23.806	264		534635	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.243	96		13158m	4.156	ng/uL	0.00
4) Pyridine-d5	3.666	84		150445	19.692	ng/ul	0.00
7) Phenol-d5	7.025	99		186601	19.860	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.190	67		144947	23.150	ng/ul	0.00
11) 2-Chlorophenol-d4	7.395	132		161700	22.262	ng/ul	0.00
15) 4-Methylphenol-d8	8.578	113		160611	21.883	ng/ul	0.00
21) Nitrobenzene-d5	9.025	128		76457	22.516	ng/ul	0.00
24) 2-Nitrophenol-d4	9.754	143		77136	21.819	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.295	165		144183	21.659	ng/ul	0.00
31) 4-Chloroaniline-d4	10.813	131		184276	20.065	ng/ul	0.00
46) Dimethylphthalate-d6	13.924	166		485463	24.682	ng/ul	0.00
49) Acenaphthylene-d8	14.207	160		570687	22.575	ng/ul	0.00
54) 4-Nitrophenol-d4	14.718	143		33053	11.021	ng/ul	0.00
60) Fluorene-d10	15.501	176		418251	24.607	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.624	200		59521	18.120	ng/ul	0.00
73) Anthracene-d10	17.354	188		616996	24.051	ng/ul	0.00
81) Pyrene-d10	19.642	212		728983	27.921	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.653	264		638490	23.742	ng/ul	0.00
Target Compounds							
8) Phenol	7.054	94		12985m	1.368	ng/ul	Qvalue

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

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