

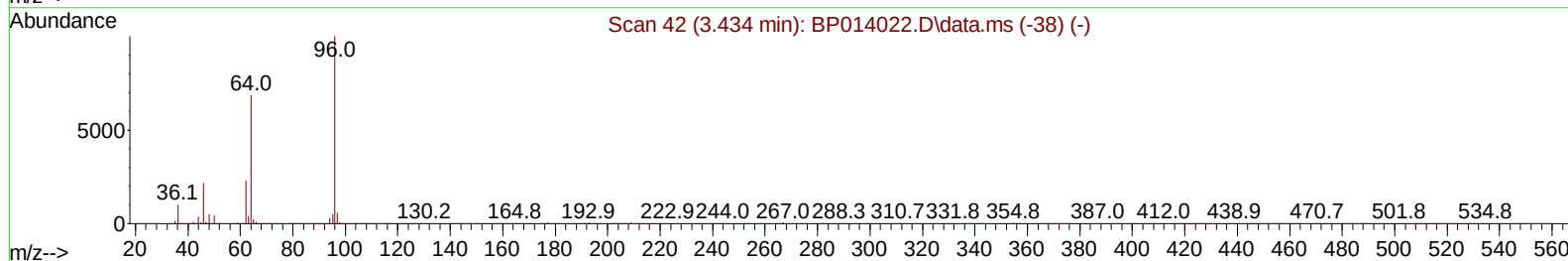
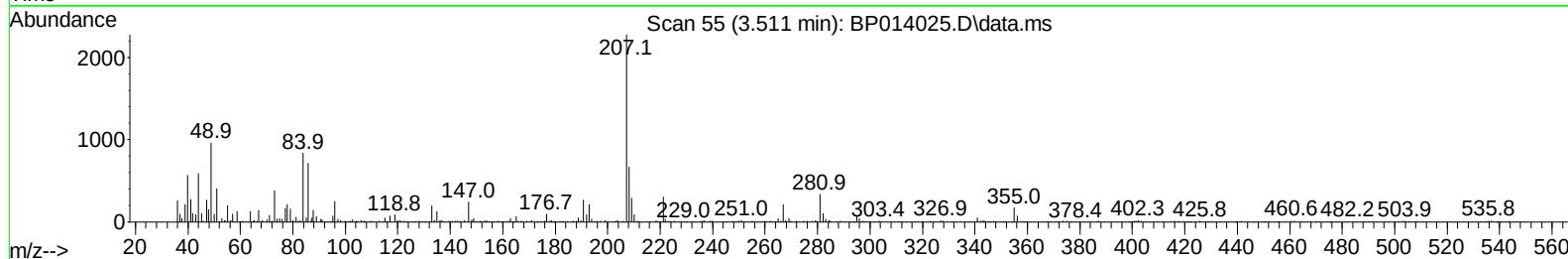
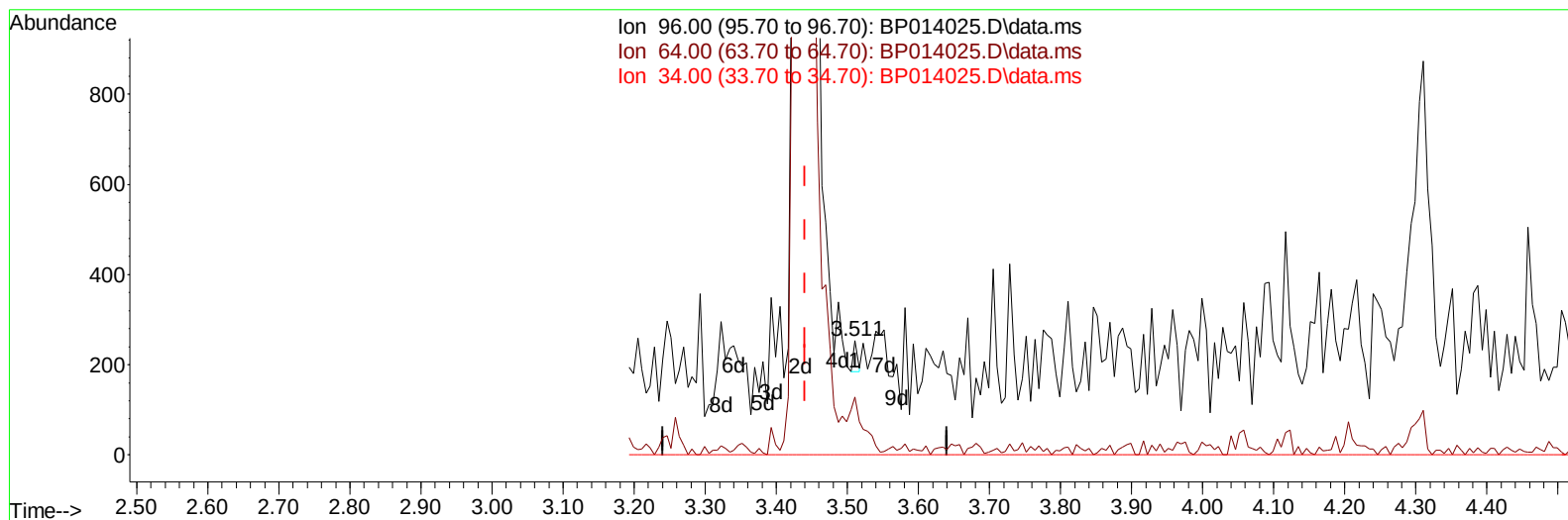
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031023\
Data File : BP014025.D
Acq On : 10 Mar 2023 12:04
Operator : CG/JU
Sample : 01713-13
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBZZ1

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 03/13/2023
Supervised By :Jagrut Upadhyay 03/13/2023

Quant Time: Mar 10 21:28:33 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Mar 09 01:10:12 2023
Response via : Initial Calibration



TIC: BP014025.D\data.ms

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

3.511min (+ 0.071) 0.01 ng/uL

response 27

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	66.60	58.10
34.00	0.00	0.00
0.00	0.00	0.00

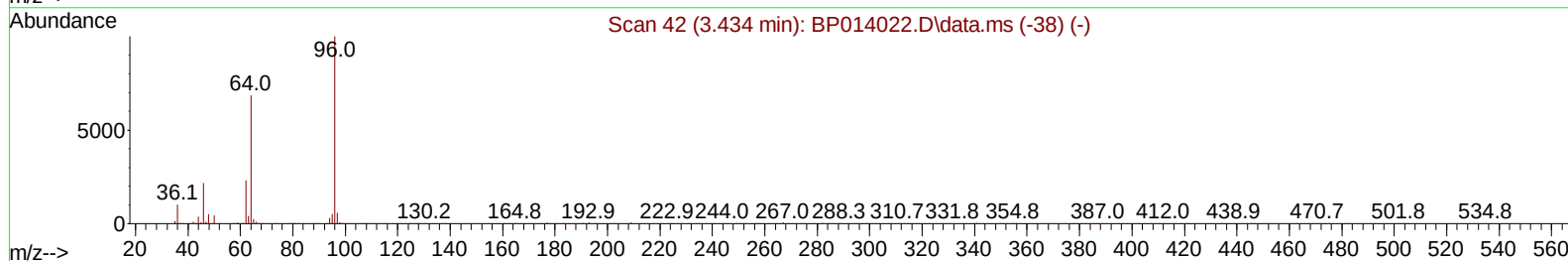
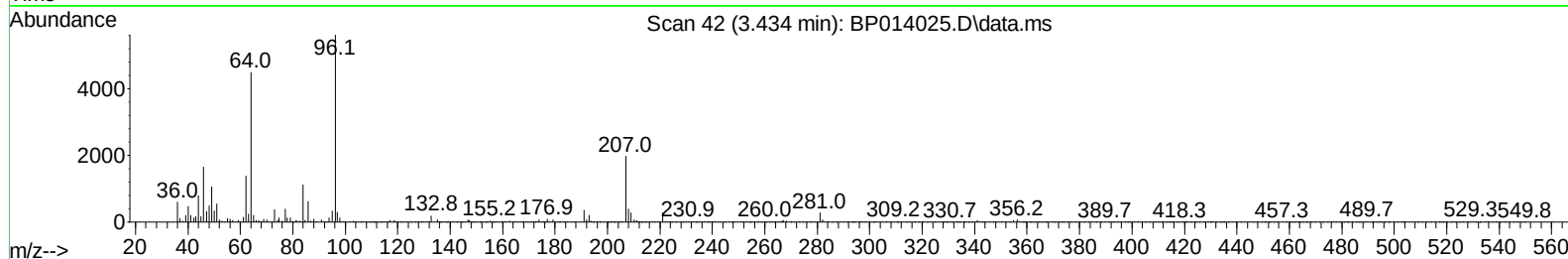
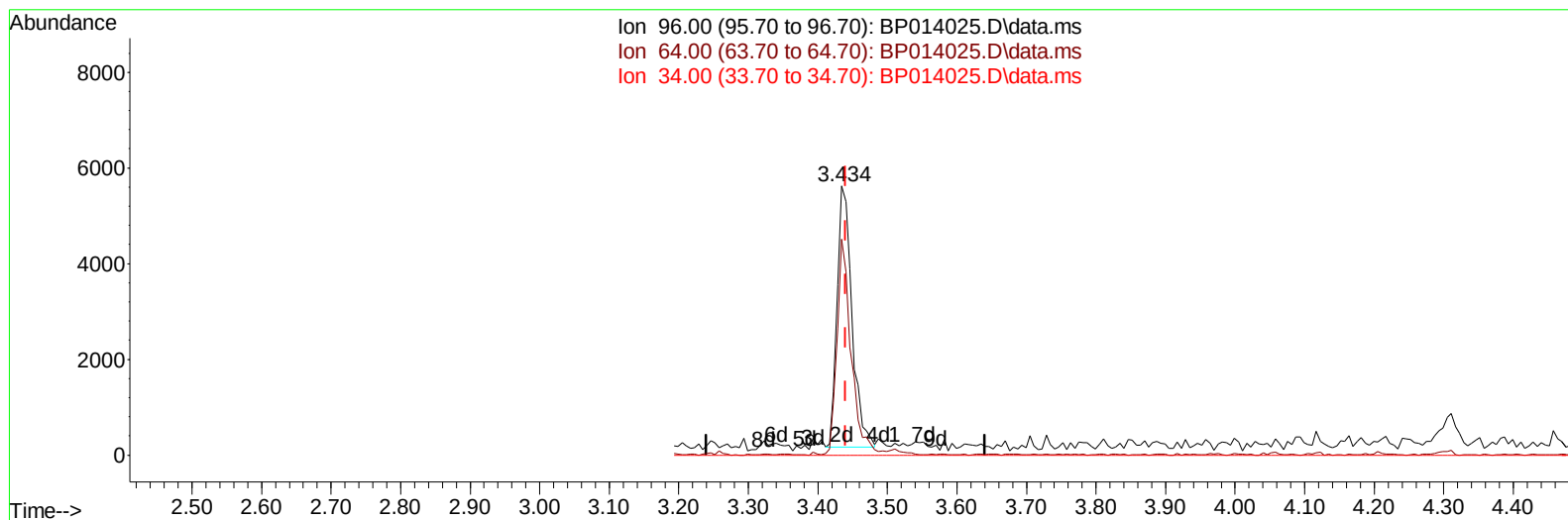
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TIC: BP014025.D\data.ms

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

3.434min (-0.006) 2.67 ng/uL m

response 8094

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	66.60	80.18#
34.00	0.00	0.00
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.075	152	114332	20.000	ng/ul	0.00
20) Naphthalene-d8	10.899	136	501079	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.716	164	360091	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.469	188	852811	20.000	ng/ul	0.00
79) Chrysene-d12	21.545	240	941869	20.000	ng/ul	-0.01
88) Perylene-d12	24.080	264	954953	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.434	96	8094m	2.675	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.228	99	143155	14.189	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.399	67	93695	15.923	ng/ul	0.00
11) 2-Chlorophenol-d4	7.605	132	113634	14.783	ng/ul	0.00
15) 4-Methylphenol-d8	8.787	113	93660	11.364	ng/ul	0.00
21) Nitrobenzene-d5	9.252	128	55412	13.860	ng/ul	0.00
24) 2-Nitrophenol-d4	9.975	143	65097	13.602	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.516	165	117521	13.050	ng/ul	0.00
31) 4-Chloroaniline-d4	11.040	131	129191	10.841	ng/ul	0.00
46) Dimethylphthalate-d6	14.116	166	479909	15.704	ng/ul	0.00
49) Acenaphthylene-d8	14.410	160	482971	14.919	ng/ul	0.00
54) 4-Nitrophenol-d4	14.910	143	54779	10.172	ng/ul	0.00
60) Fluorene-d10	15.704	176	408498	15.772	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.822	200	67999	10.550	ng/ul	-0.01
73) Anthracene-d10	17.569	188	646213	15.609	ng/ul	0.00
81) Pyrene-d10	19.786	212	916196	17.637	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.916	264	867033	17.069	ng/ul	-0.01

Target Compounds Qvalue

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

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