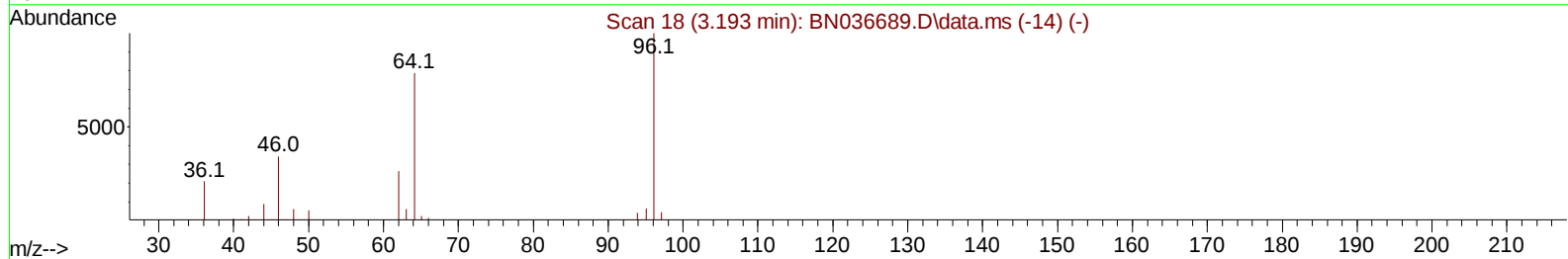
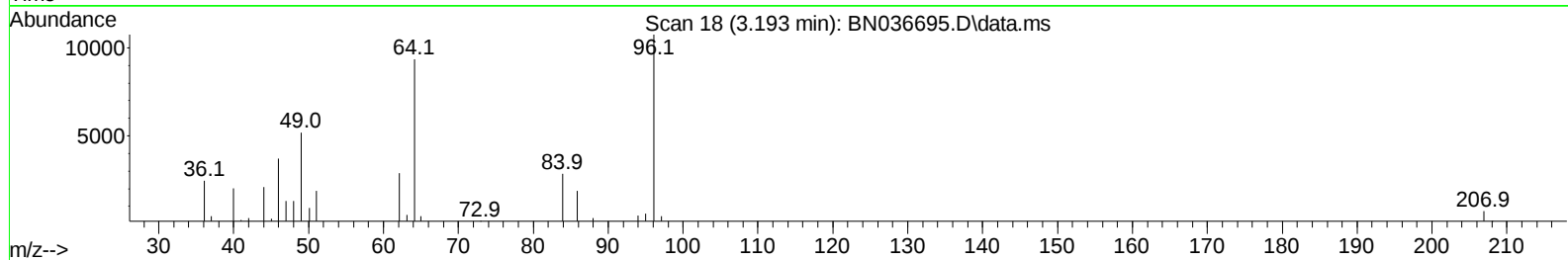
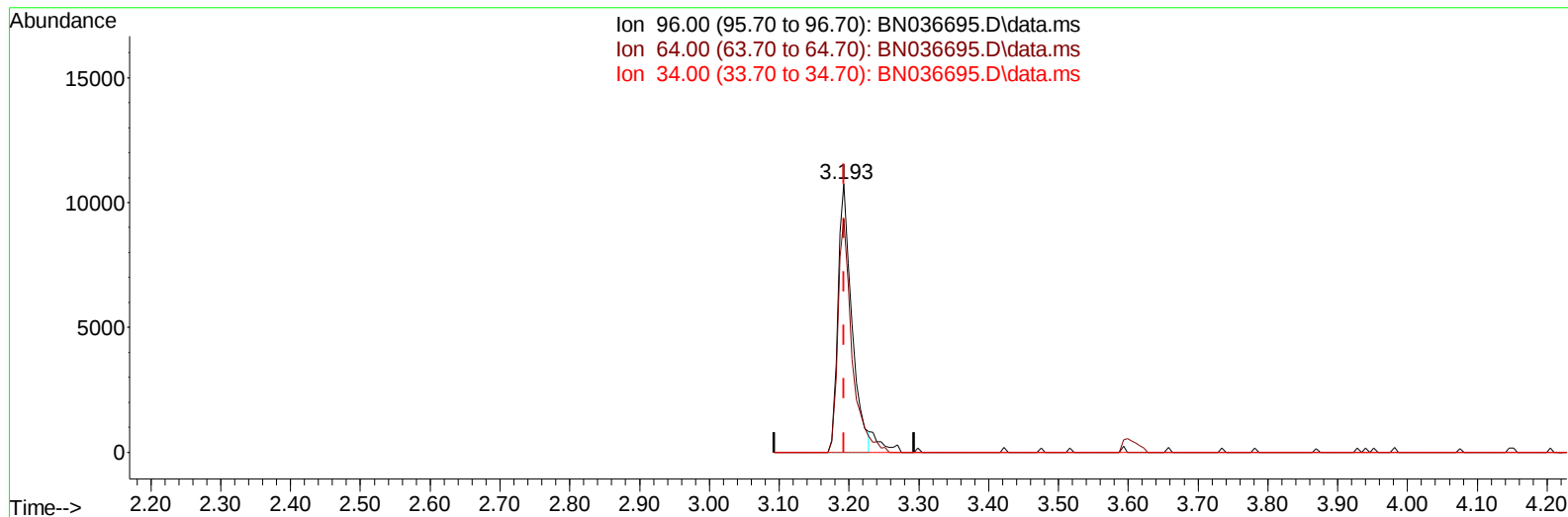


Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN032025\
Data File : BN036695.D
Acq On : 20 Mar 2025 14:11
Operator : RC/JU
Sample : PB167182BL
Misc :
ALS Vial : 8 Sample Multiplier: 1

Quant Time: Mar 20 14:45:09 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN022825.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Mar 20 13:16:23 2025
Response via : Initial Calibration



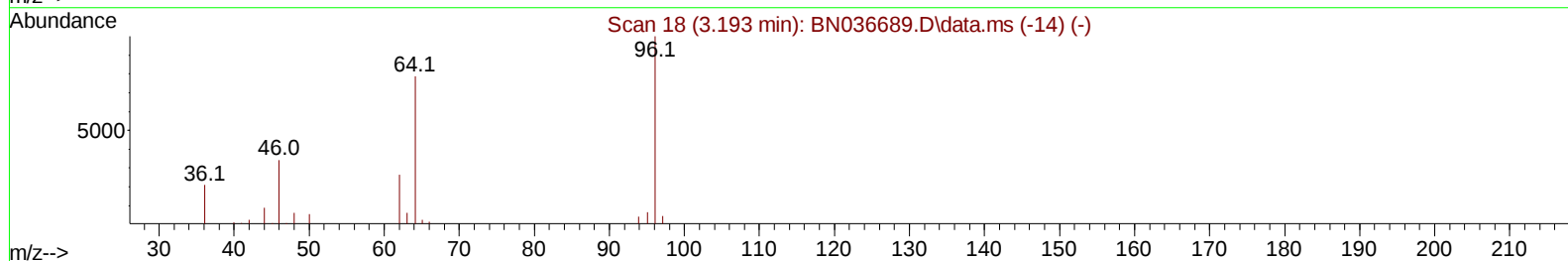
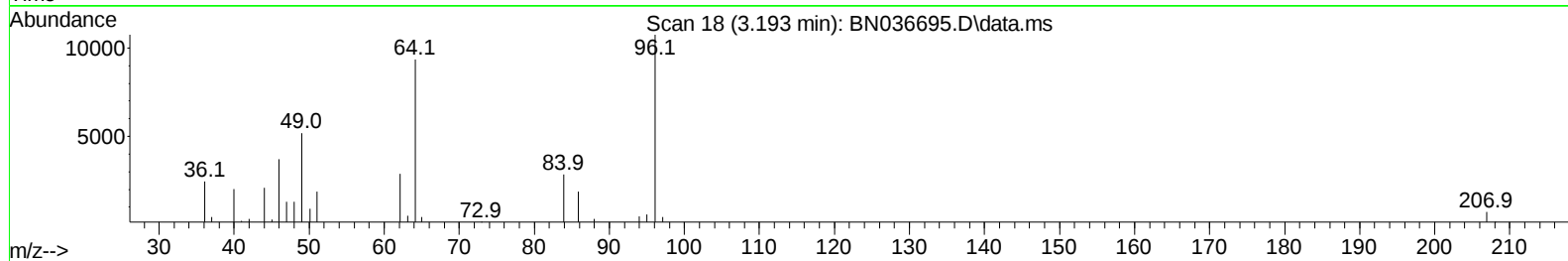
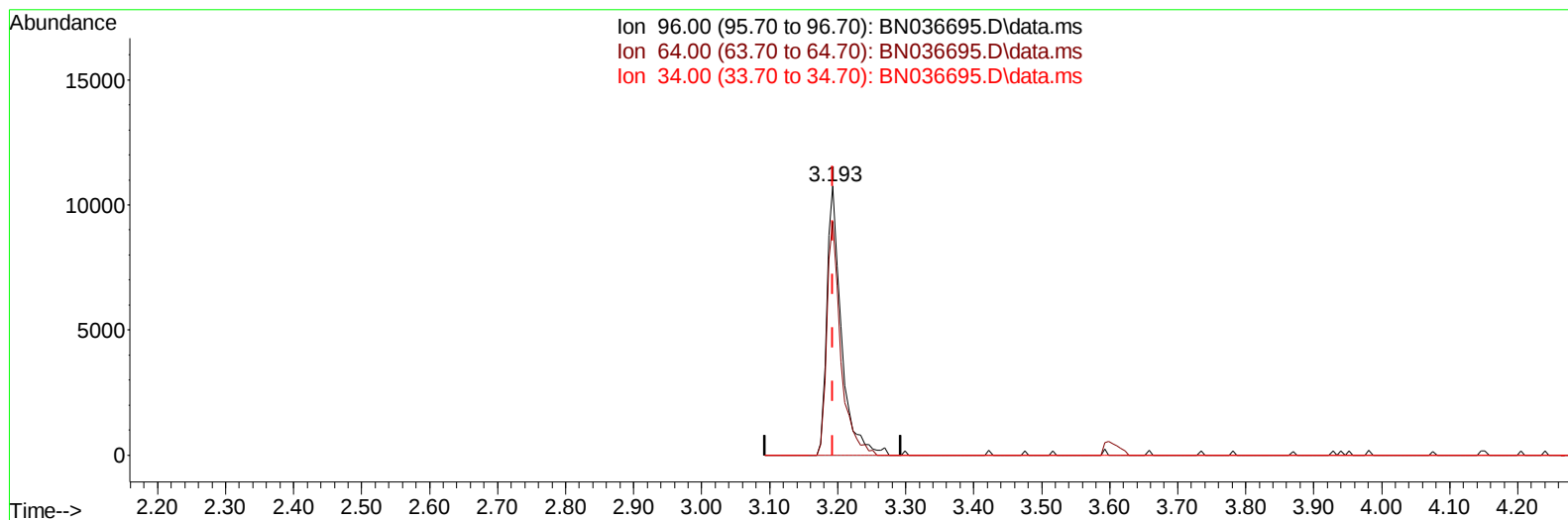
TIC: BN036695.D\data.ms

(3) 1,4-Dioxane-d8 (S)**3.193min (-0.000) 6.82 ng/uL****response 15001**

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	87.00	86.83
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN032025\
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Misc :
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QLast Update : Thu Mar 20 13:16:23 2025
Response via : Initial Calibration



TIC: BN036695.D\data.ms

(3) 1,4-Dioxane-d8 (S)**3.193min (-0.000) 7.23 ng/uL m****response 15909**

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	87.00	86.83
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN032025\
 Data File : BN036695.D
 Acq On : 20 Mar 2025 14:11
 Operator : RC/JU
 Sample : PB167182BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Quant Time: Mar 20 14:45:38 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN022825.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Mar 20 13:16:23 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.704	152	85537	20.000	ng/ul	0.00
20) Naphthalene-d8	10.493	136	362865	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.345	164	243044	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.092	188	509002	20.000	ng/ul	0.00
79) Chrysene-d12	21.327	240	516039	20.000	ng/ul	0.00
88) Perylene-d12	24.280	264	532816	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.193	96	15909m	7.231	ng/uL	0.00
4) Pyridine-d5	3.599	84	233824	34.769	ng/ul	0.00
7) Phenol-d5	6.875	99	283350	34.683	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.040	67	190559	32.785	ng/ul	0.00
11) 2-Chlorophenol-d4	7.234	132	208362	36.271	ng/ul	0.00
15) 4-Methylphenol-d8	8.410	113	228150	33.822	ng/ul	0.00
21) Nitrobenzene-d5	8.857	128	97388	33.953	ng/ul	0.00
24) 2-Nitrophenol-d4	9.575	143	105861	34.933	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.110	165	194492	35.385	ng/ul	0.00
31) 4-Chloroaniline-d4	10.628	131	275164	32.418	ng/ul	0.00
46) Dimethylphthalate-d6	13.757	166	661578	36.195	ng/ul	0.00
49) Acenaphthylene-d8	14.039	160	709904	34.769	ng/ul	0.00
54) 4-Nitrophenol-d4	14.545	143	108308	29.915	ng/ul	0.00
60) Fluorene-d10	15.345	176	542142	35.166	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.463	200	103944	29.842	ng/ul	0.00
73) Anthracene-d10	17.192	188	885461	35.351	ng/ul	0.00
81) Pyrene-d10	19.492	212	1028138	36.470	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.080	264	984082	37.067	ng/ul	0.00

Target Compounds	Qvalue

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

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