

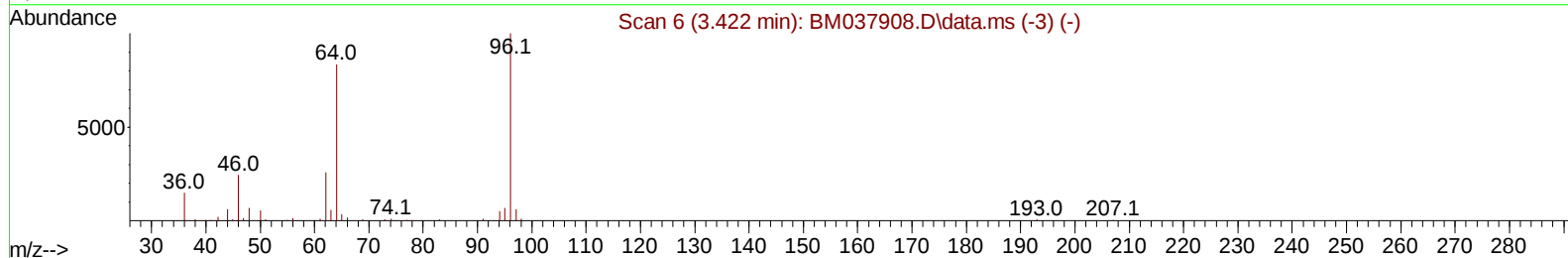
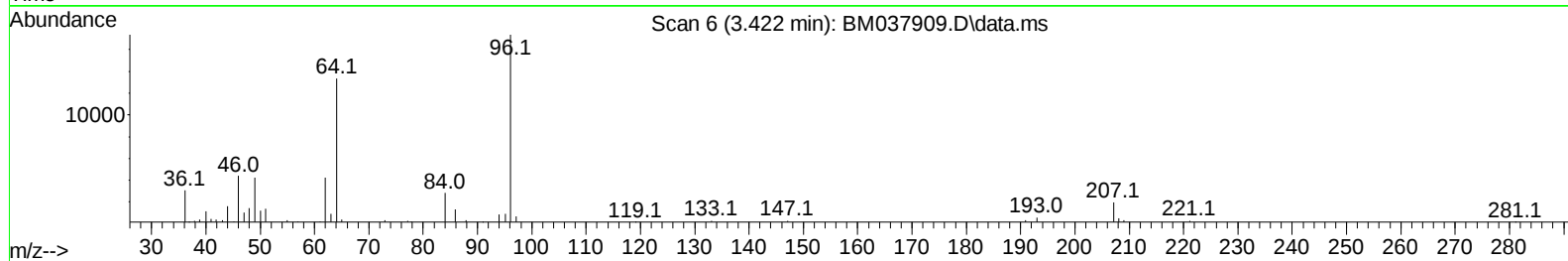
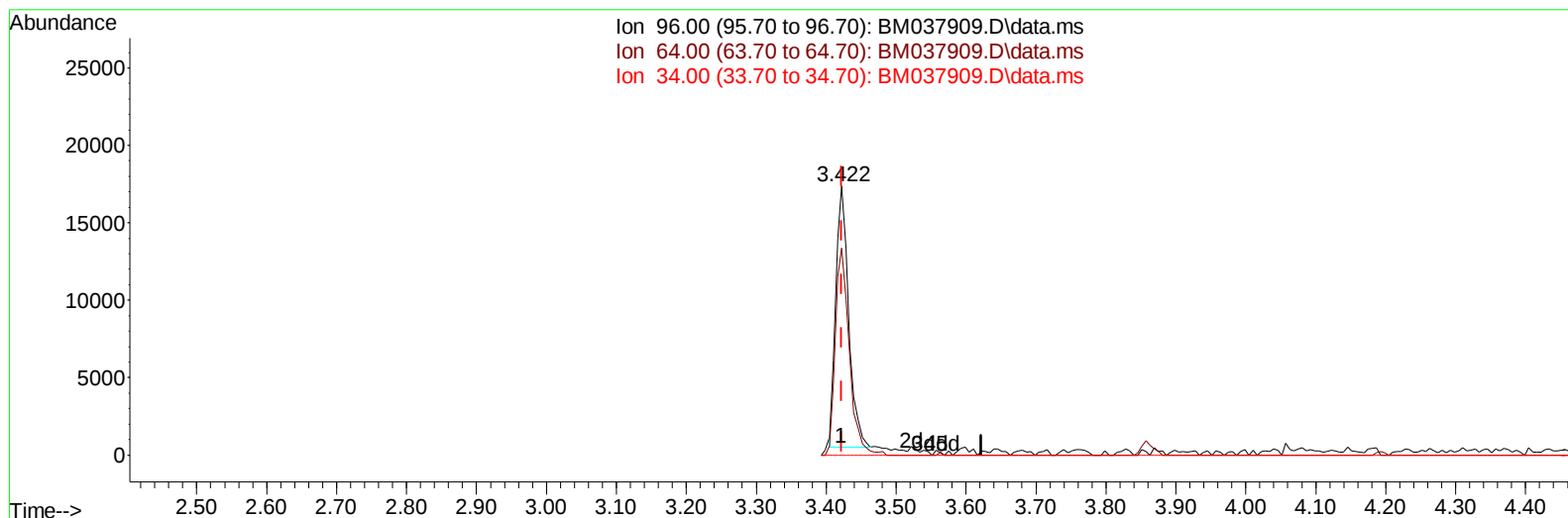
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120922\
 Data File : BM037909.D
 Acq On : 09 Dec 2022 10:22
 Operator : CG/JU
 Sample : PB149458BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK458

Manual IntegrationsAPPROVED

Quant Time: Dec 09 21:46:51 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 09 21:34:57 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/10/2022
 Supervised By :mohammad ahmed 12/10/2022



TIC: BM037909.D\data.ms

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

3.422min (-0.000) 4.17 ng/uL

response 21563

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	76.40	76.97
34.00	0.00	0.00
0.00	0.00	0.00

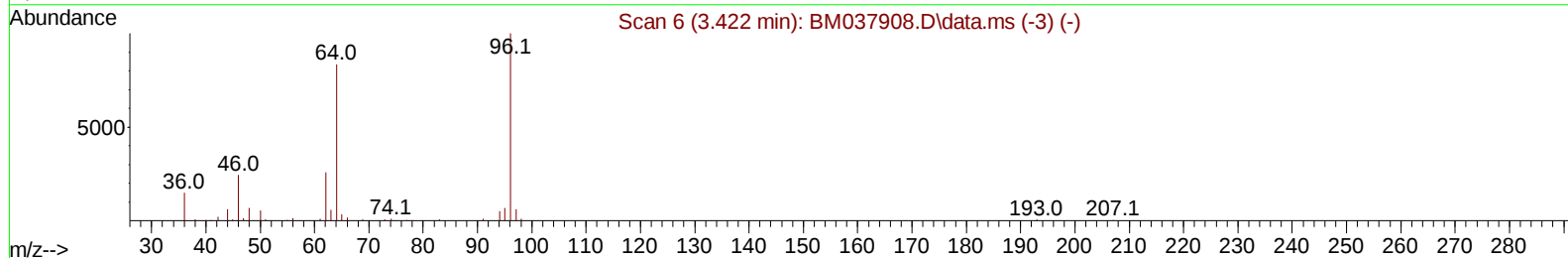
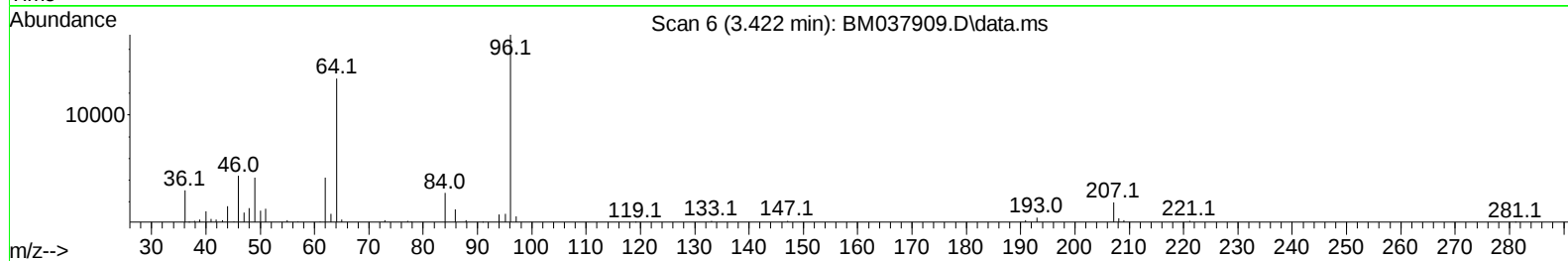
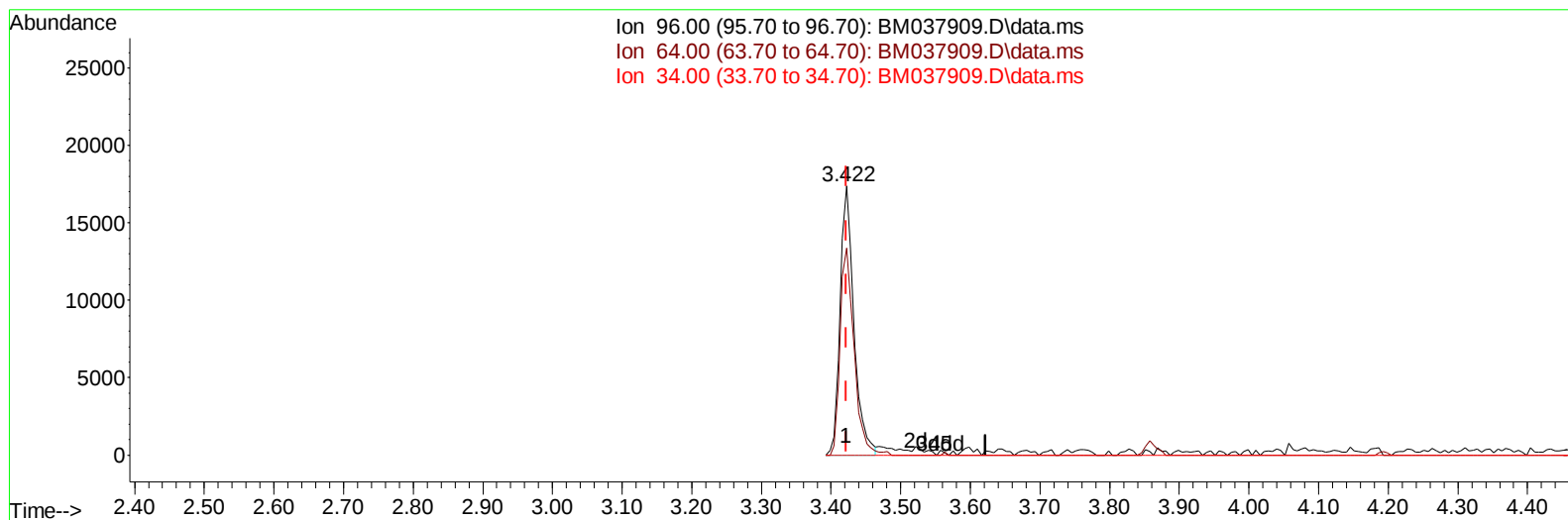
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(3) 1,4-Dioxane-d8 (S)

3.422min (-0.000) 4.62 ng/uL m

response 23888

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	76.40	76.97
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	8.081	152	234546	20.000	ng/ul	0.00
20) Naphthalene-d8	10.904	136	947158	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.710	164	513337	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.451	188	934941	20.000	ng/ul	0.00
79) Chrysene-d12	21.621	240	729062	20.000	ng/ul	0.00
88) Perylene-d12	24.103	264	778731	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.422	96	23888m	4.624	ng/uL	0.00
4) Pyridine-d5	3.857	84	358839	23.417	ng/ul	0.00
7) Phenol-d5	7.234	99	480166	25.193	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.404	67	290073	24.994	ng/ul	0.00
11) 2-Chlorophenol-d4	7.610	132	409334	26.181	ng/ul	0.00
15) 4-Methylphenol-d8	8.792	113	373161	25.082	ng/ul	0.00
21) Nitrobenzene-d5	9.257	128	202596	27.016	ng/ul	0.00
24) 2-Nitrophenol-d4	9.980	143	213819	27.207	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.522	165	366479	24.551	ng/ul	0.00
31) 4-Chloroaniline-d4	11.039	131	518367	25.157	ng/ul	0.00
46) Dimethylphthalate-d6	14.121	166	970143	26.509	ng/ul	0.00
49) Acenaphthylene-d8	14.410	160	1163338	26.009	ng/ul	0.00
54) 4-Nitrophenol-d4	14.904	143	112055	17.966	ng/ul	0.00
60) Fluorene-d10	15.698	176	837813	25.695	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.815	200	76952	15.395	ng/ul	0.00
73) Anthracene-d10	17.551	188	1183640	27.126	ng/ul	0.00
81) Pyrene-d10	19.833	212	1216244	27.771	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.938	264	1100015	28.281	ng/ul	0.00

Target Compounds Qvalue

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

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