

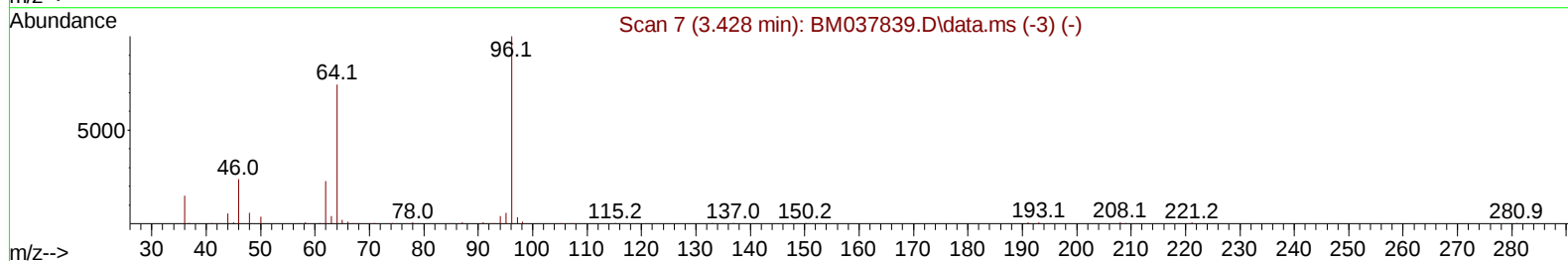
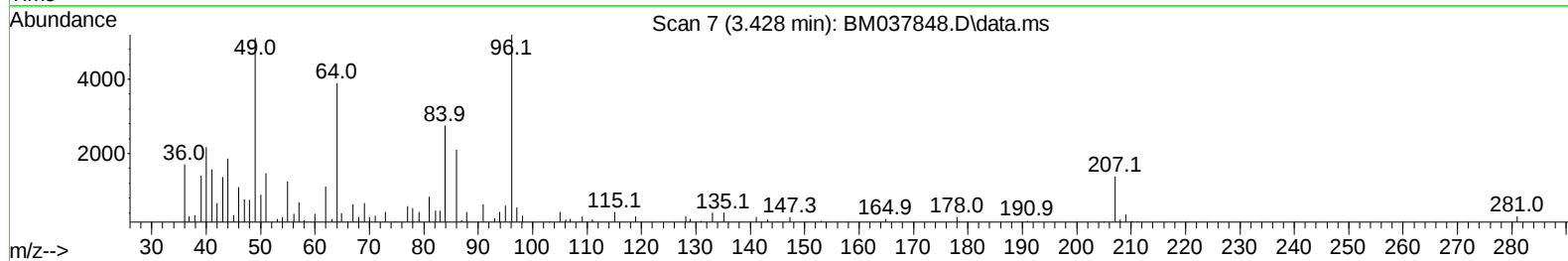
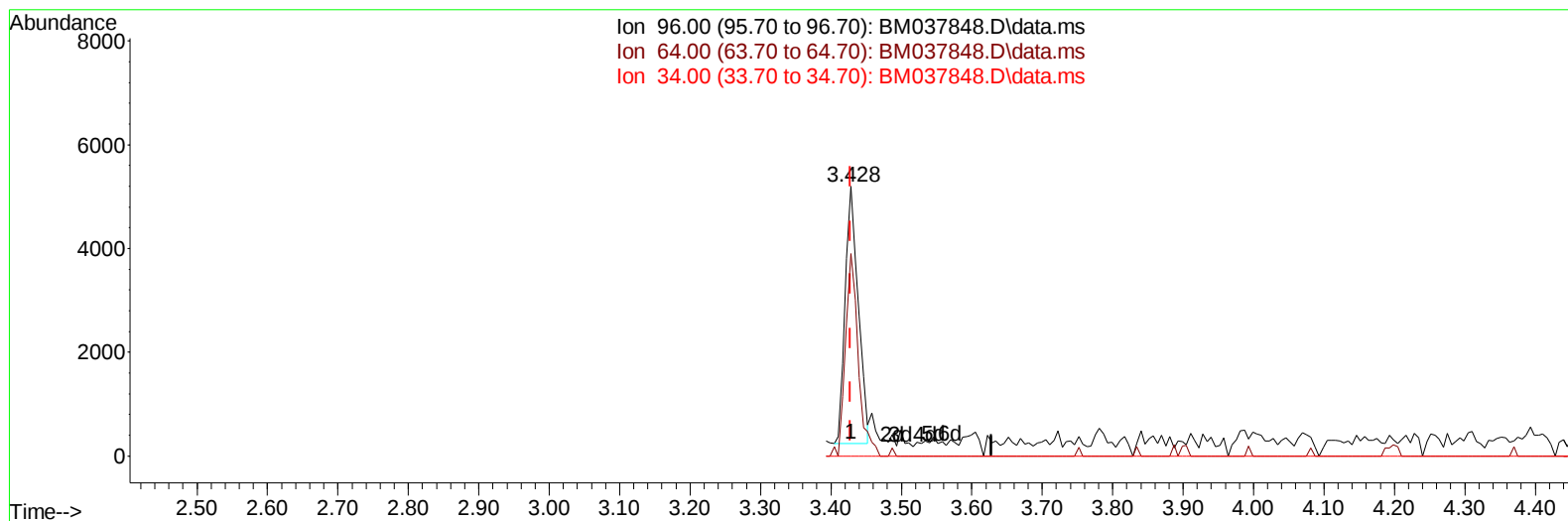
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120522\
 Data File : BM037848.D
 Acq On : 05 Dec 2022 17:30
 Operator : CG/JU
 Sample : N5738-01DL 5X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 GBNH2DL

Manual IntegrationsAPPROVED

Quant Time: Dec 05 23:27:09 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 02 01:14:59 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/06/2022
 Supervised By :mohammad ahmed 12/06/2022



TIC: BM037848.D\data.ms

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

3.428min (+ 0.000) 0.93 ng/uL

response 6286

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	82.40	74.98
34.00	0.00	0.00
0.00	0.00	0.00

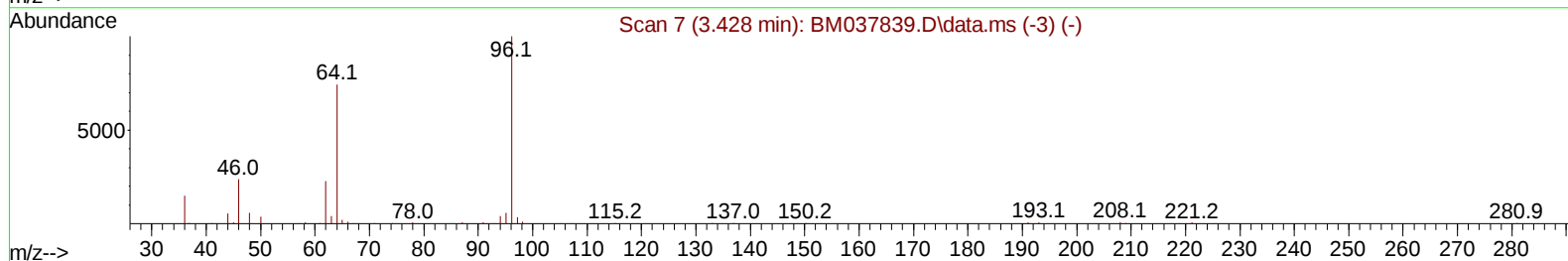
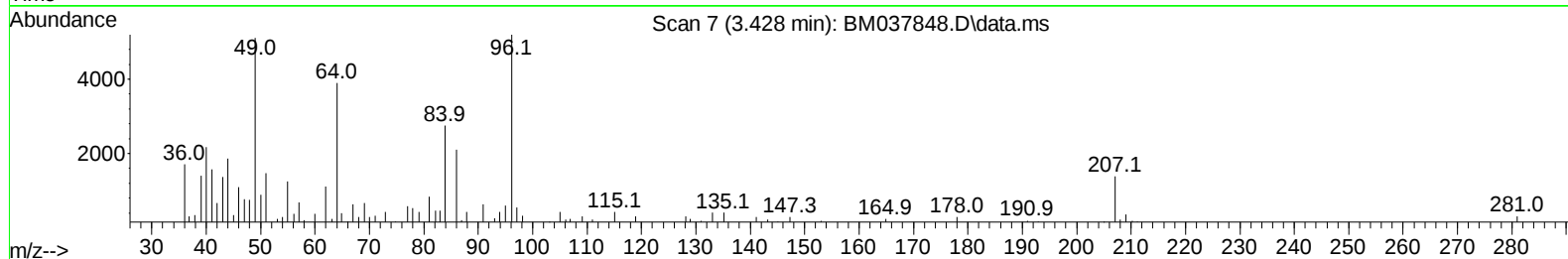
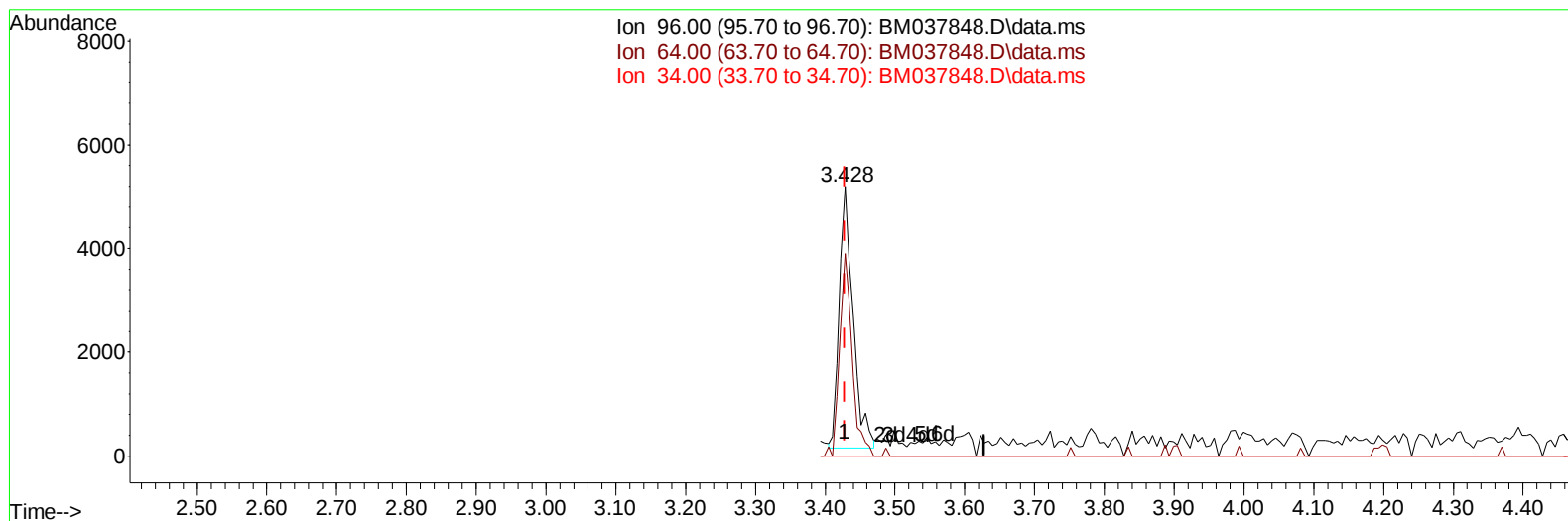
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TIC: BM037848.D\data.ms

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

3.428min (+ 0.000) 1.02 ng/uL m

response 6915

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	82.40	74.98
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	8.093	152	318605	20.000	ng/ul	0.00
20) Naphthalene-d8	10.916	136	1360787	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.722	164	802636	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.463	188	1597271	20.000	ng/ul	0.00
79) Chrysene-d12	21.627	240	992631	20.000	ng/ul	0.00
88) Perylene-d12	24.115	264	1062402	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.428	96	6915m	1.018	ng/uL	0.00
4) Pyridine-d5	3.887	84	36900	1.680	ng/ul	0.02
7) Phenol-d5	7.246	99	141505	4.930	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.416	67	88498	5.285	ng/ul	0.00
11) 2-Chlorophenol-d4	7.622	132	115426	5.318	ng/ul	0.00
15) 4-Methylphenol-d8	8.804	113	87774	3.845	ng/ul	0.00
21) Nitrobenzene-d5	9.263	128	60256	5.574	ng/ul	0.00
24) 2-Nitrophenol-d4	9.993	143	63885	5.711	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.528	165	106632	5.140	ng/ul	0.00
31) 4-Chloroaniline-d4	11.057	131	33008	1.037	ng/ul	0.00
46) Dimethylphthalate-d6	14.128	166	316297	5.739	ng/ul	0.00
49) Acenaphthylene-d8	14.416	160	380258	5.673	ng/ul	0.00
54) 4-Nitrophenol-d4	14.916	143	38823	3.394	ng/ul	0.00
60) Fluorene-d10	15.710	176	290158	5.803	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.822	200	28137	3.344	ng/ul	0.00
73) Anthracene-d10	17.557	188	402050	5.473	ng/ul	0.00
81) Pyrene-d10	19.839	212	395725	7.696	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.951	264	278165	5.226	ng/ul	0.00
Target Compounds						
					Qvalue	
71) Pentachlorophenol	17.110	266	576452	48.966	ng/ul	99

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

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