

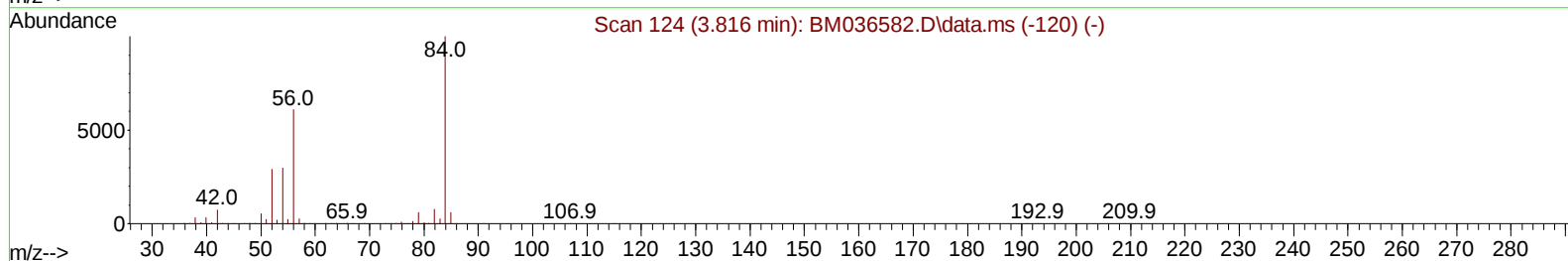
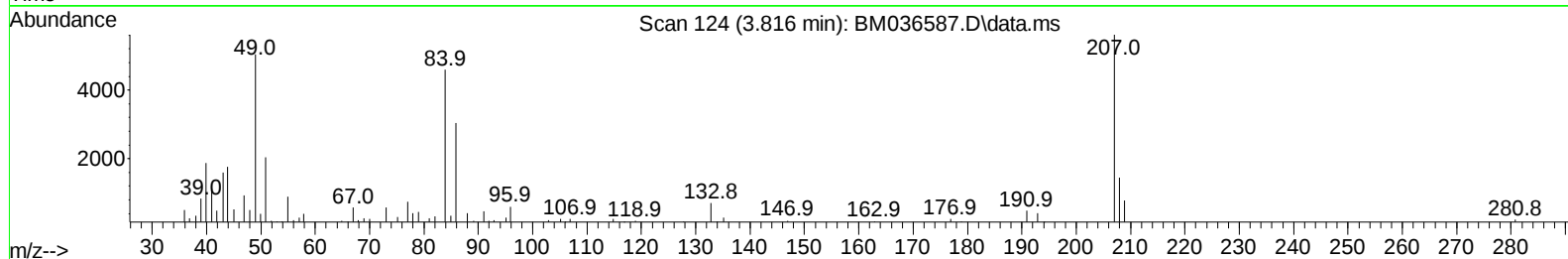
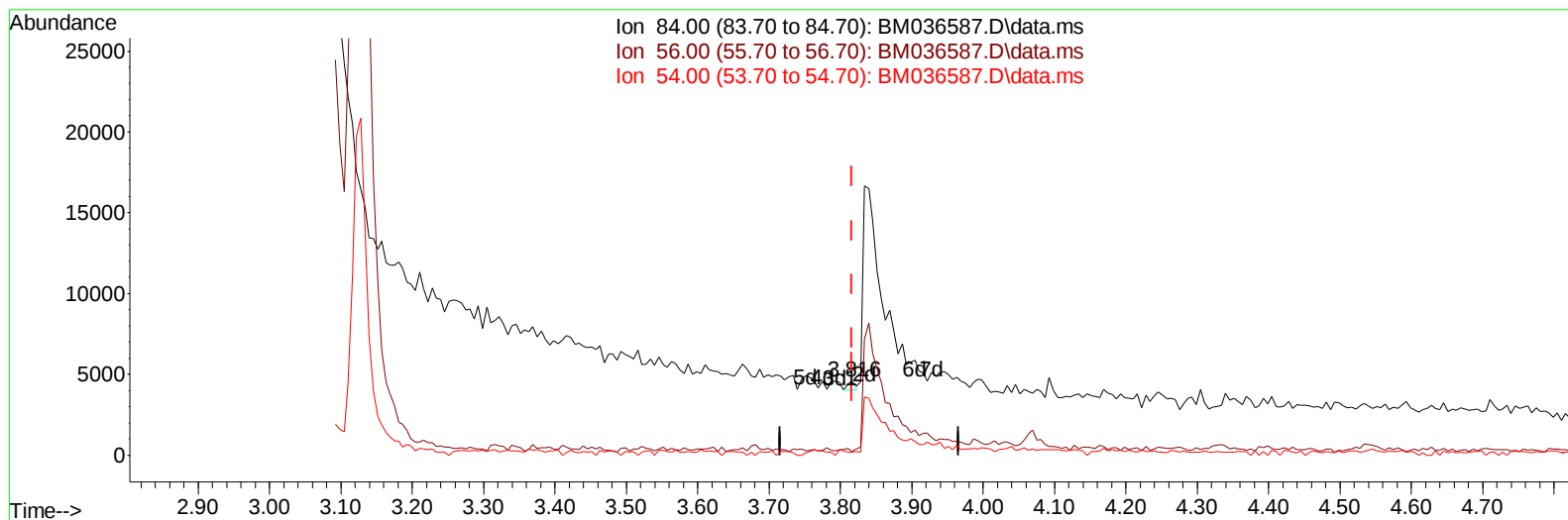
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091922\
 Data File : BM036587.D
 Acq On : 19 Sep 2022 19:23
 Operator : CG/JU
 Sample : N4604-15
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A5775

Manual IntegrationsAPPROVED

Quant Time: Sep 19 22:34:27 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 16 02:39:54 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 09/20/2022
 Supervised By :mohammad ahmed 09/21/2022



TIC: BM036587.D\data.ms

(4) Pyridine-d5 (S)

3.816min (+ 0.000) 0.01 ng/u1

response 403

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	61.90	4.71#
54.00	30.90	3.57#
0.00	0.00	0.00

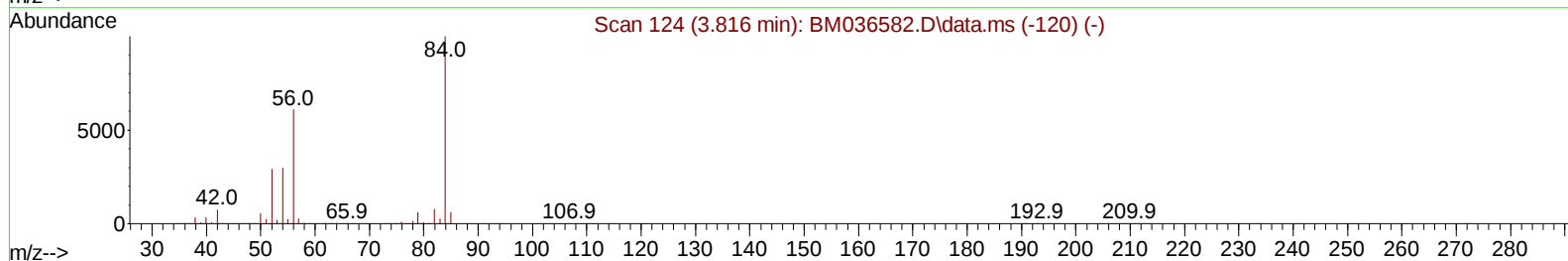
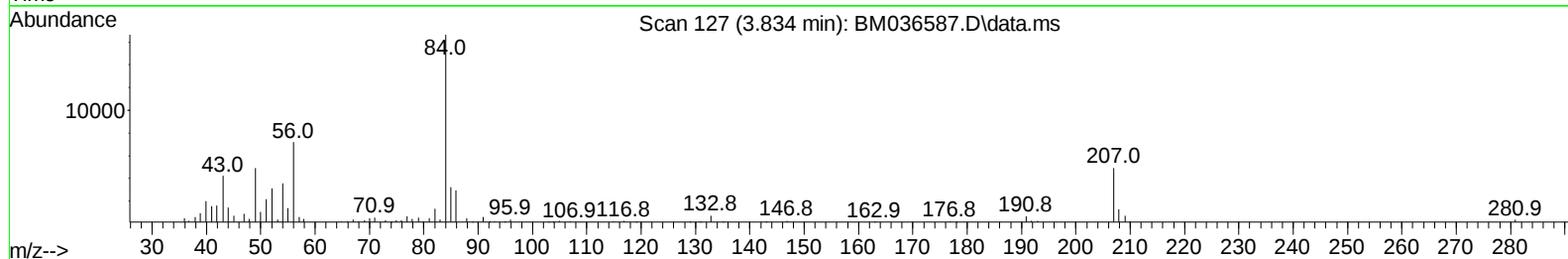
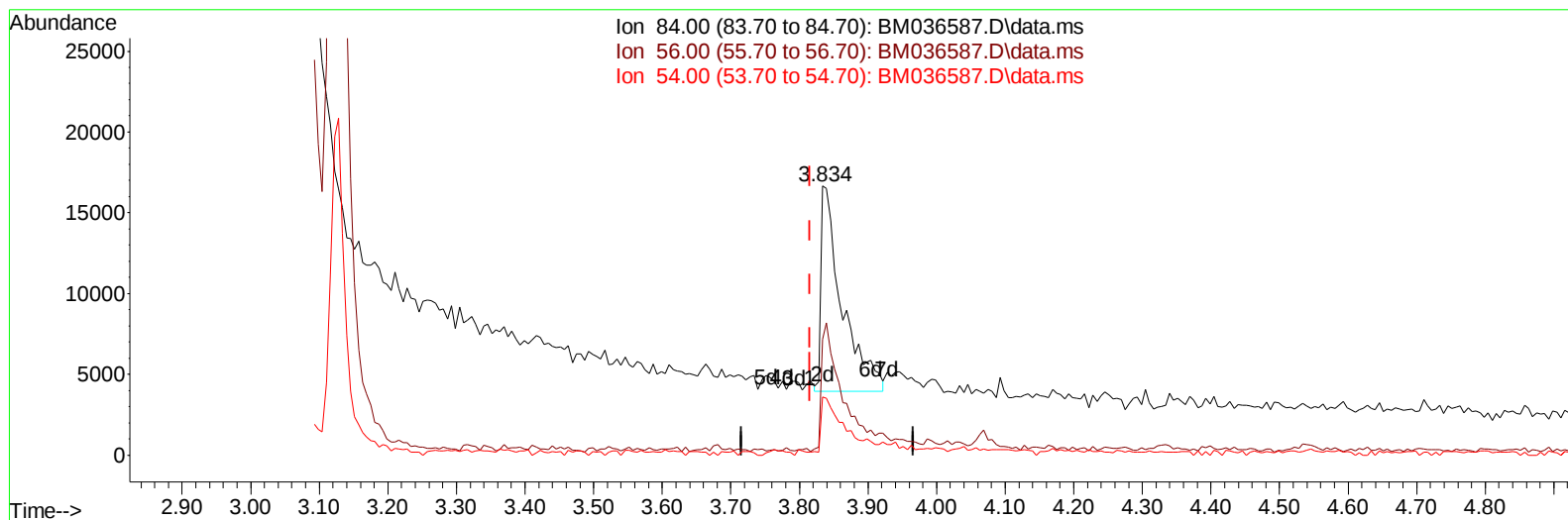
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TIC: BM036587.D\data.ms

(4) Pyridine-d5 (S)

3.834min (+ 0.018) 0.85 ng/u1 m

response 26916

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	61.90	43.15#
54.00	30.90	21.61#
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	8.028	152	407503	20.000	ng/ul	0.00
20) Naphthalene-d8	10.839	136	1844489	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.651	164	1222298	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.386	188	2502815	20.000	ng/ul	0.00
79) Chrysene-d12	21.550	240	2380248	20.000	ng/ul	0.00
88) Perylene-d12	23.991	264	2138795	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	36711	3.377	ng/uL	0.00
4) Pyridine-d5	3.834	84	26916m	0.849	ng/ul	0.02
7) Phenol-d5	7.175	99	729990	20.046	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.351	67	442520	20.010	ng/ul	0.00
11) 2-Chlorophenol-d4	7.551	132	565228	21.559	ng/ul	0.00
15) 4-Methylphenol-d8	8.728	113	569519	21.165	ng/ul	0.00
21) Nitrobenzene-d5	9.192	128	268011	21.468	ng/ul	0.00
24) 2-Nitrophenol-d4	9.916	143	254816	22.787	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.451	165	529530	20.433	ng/ul	0.00
31) 4-Chloroaniline-d4	10.975	131	566559	13.149	ng/ul	0.00
46) Dimethylphthalate-d6	14.063	166	1832286	22.226	ng/ul	0.00
49) Acenaphthylene-d8	14.345	160	2112005	21.599	ng/ul	0.00
54) 4-Nitrophenol-d4	14.827	143	290453	18.165	ng/ul	0.00
60) Fluorene-d10	15.639	176	1637180	22.161	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	160583	16.217	ng/ul	0.00
73) Anthracene-d10	17.486	188	2417121	21.122	ng/ul	0.00
81) Pyrene-d10	19.762	212	2959190	25.271	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.832	264	2512053	24.041	ng/ul	0.00

Target Compounds Qvalue

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

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