

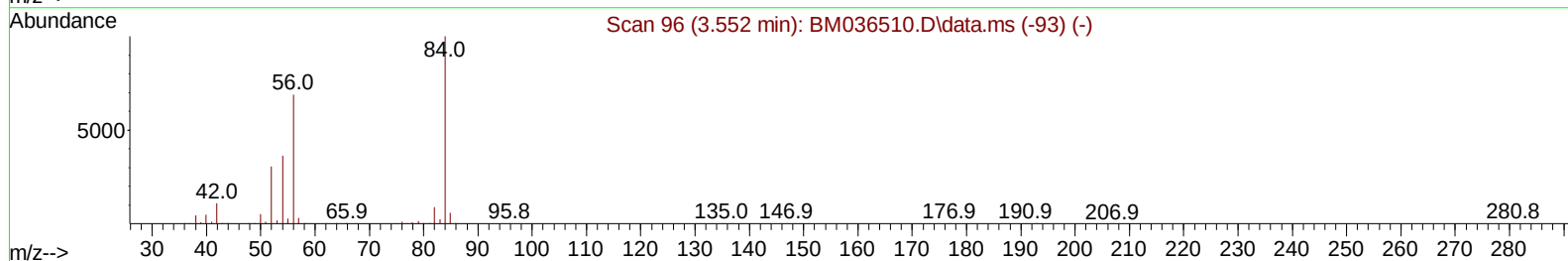
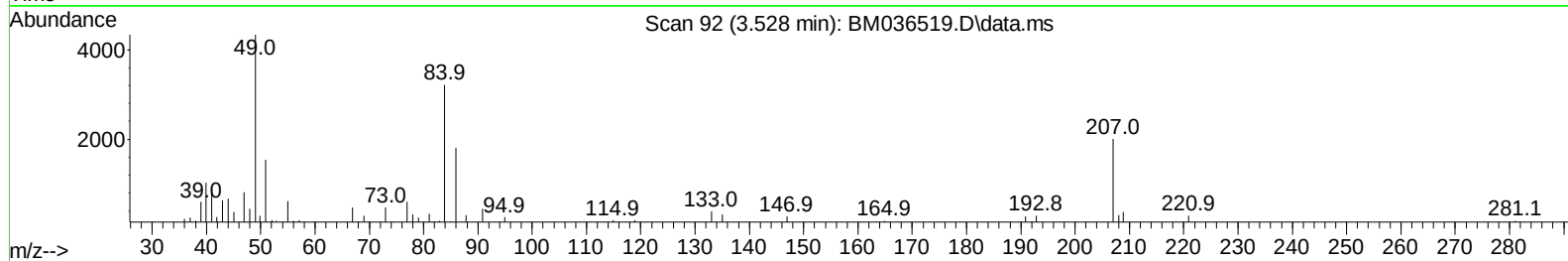
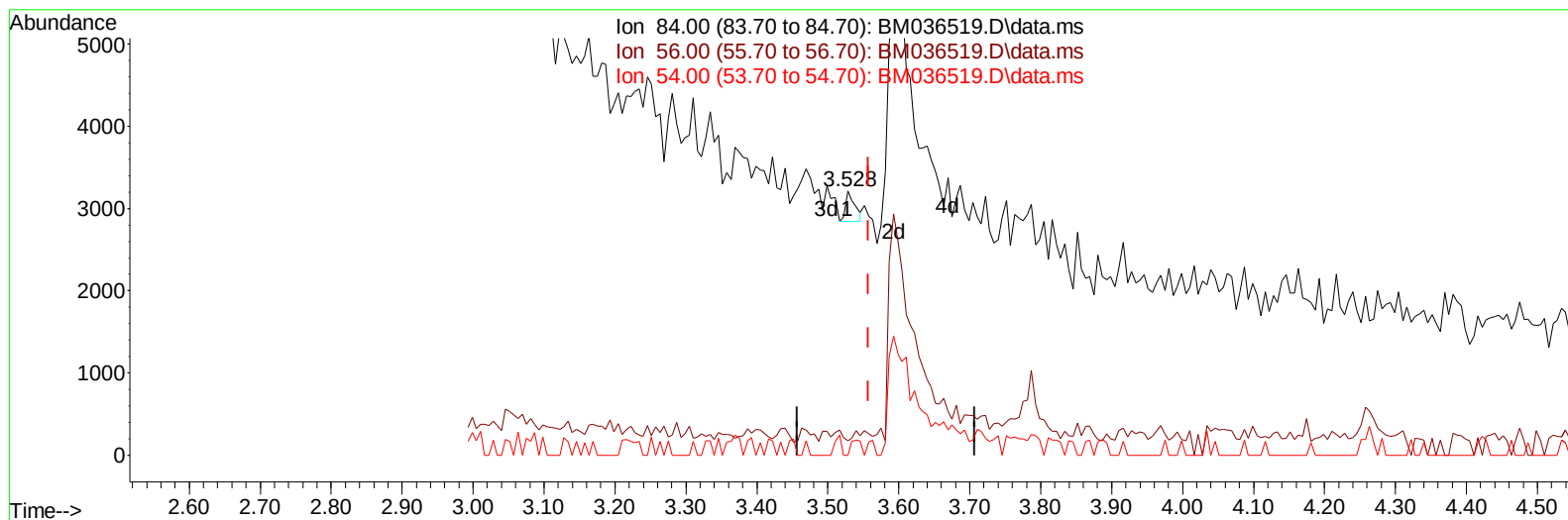
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090822\
 Data File : BM036519.D
 Acq On : 07 Sep 2022 21:13
 Operator : CG/JU
 Sample : N4483-12
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW360

Manual IntegrationsAPPROVED

Quant Time: Sep 08 00:54:44 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM081622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 02 01:33:23 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 09/08/2022
 Supervised By :mohammad ahmed 09/19/2022



TIC: BM036519.D\data.ms

(4) Pyridine-d5 (S)

3.528min (-0.030) 0.02 ng/ul

response 336

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	70.90	5.47#
54.00	32.70	0.00#
0.00	0.00	0.00

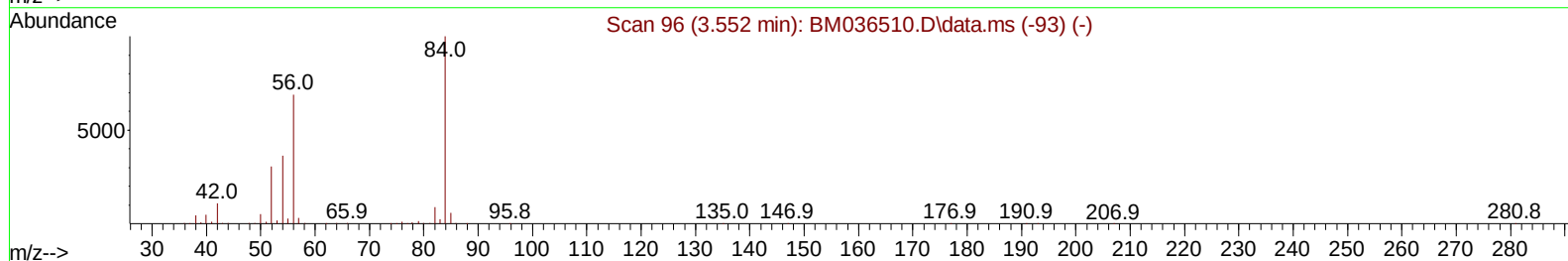
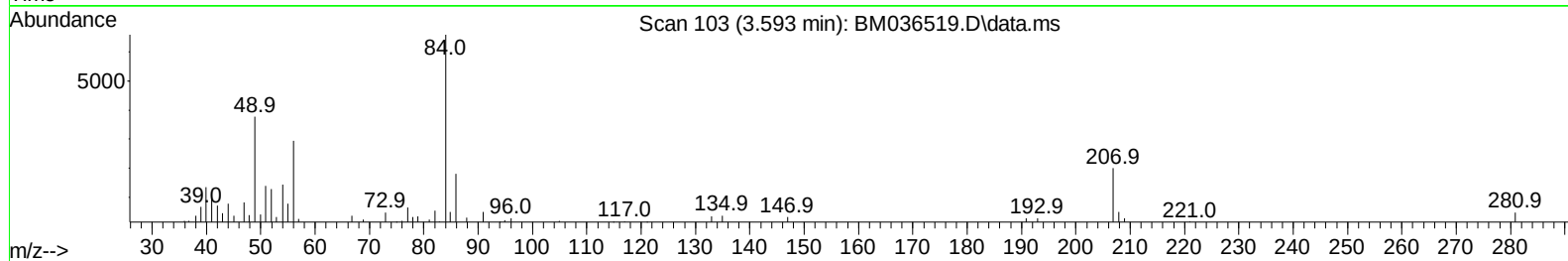
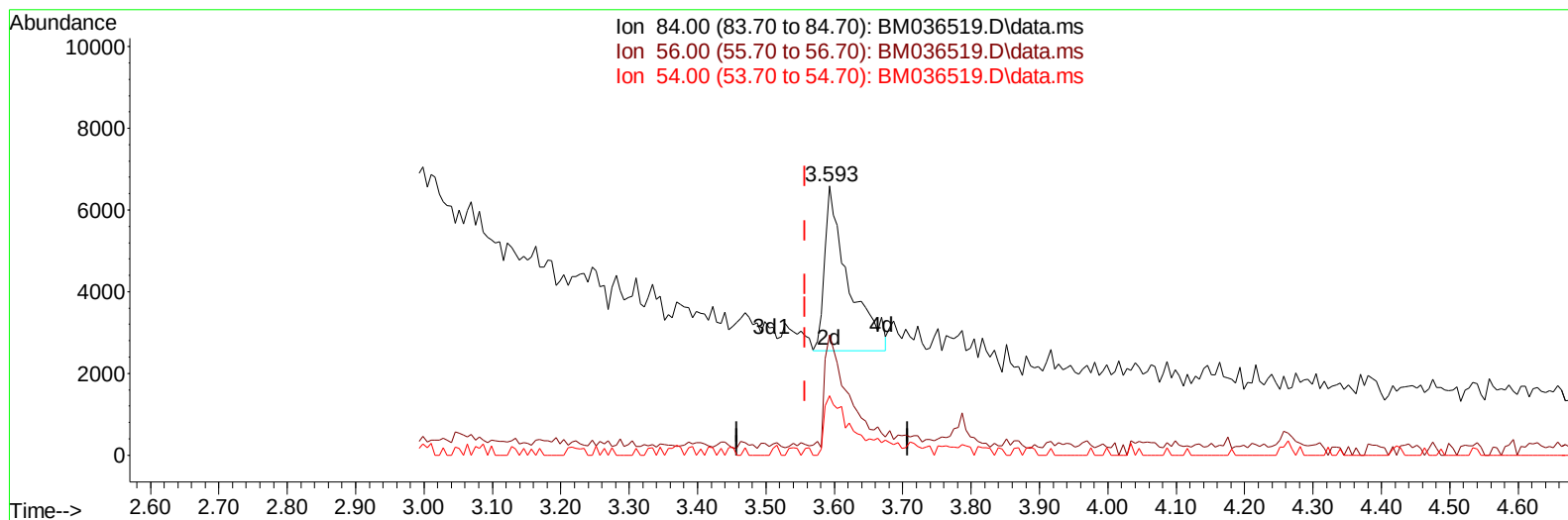
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(4) Pyridine-d5 (S)

3.593min (+ 0.035) 0.66 ng/ul m

response 9722

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	70.90	44.61#
54.00	32.70	22.00#
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	7.722	152	203364	20.000	ng/ul	0.00
20) Naphthalene-d8	10.522	136	957263	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.374	164	609907	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.115	188	1236571	20.000	ng/ul	0.00
79) Chrysene-d12	21.309	240	995011	20.000	ng/ul	0.00
88) Perylene-d12	23.597	264	817082	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.128	96	22581	4.232	ng/uL	0.00
4) Pyridine-d5	3.593	84	9722m	0.661	ng/ul	0.04
7) Phenol-d5	6.904	99	460251	26.514	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.063	67	309199	28.298	ng/ul	0.00
11) 2-Chlorophenol-d4	7.251	132	393893	28.880	ng/ul	0.00
15) 4-Methylphenol-d8	8.445	113	390132	30.100	ng/ul	0.00
21) Nitrobenzene-d5	8.892	128	207884	28.643	ng/ul	0.00
24) 2-Nitrophenol-d4	9.610	143	221211	30.048	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.151	165	386912	29.413	ng/ul	0.00
31) 4-Chloroaniline-d4	10.669	131	494955	24.271	ng/ul	0.00
46) Dimethylphthalate-d6	13.798	166	1243422	32.225	ng/ul	0.00
49) Acenaphthylene-d8	14.068	160	1518511	31.553	ng/ul	0.00
54) 4-Nitrophenol-d4	14.604	143	137540	18.812	ng/ul	0.00
60) Fluorene-d10	15.368	176	1109529	31.822	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.504	200	130654	20.552	ng/ul	0.00
73) Anthracene-d10	17.215	188	1682186	30.594	ng/ul	0.00
81) Pyrene-d10	19.515	212	1918942	37.178	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.450	264	1319438	32.936	ng/ul	0.00
Target Compounds						
					Qvalue	
8) Phenol	6.928	94	24323	1.348	ng/ul#	79
80) Fluoranthene	19.180	202	152588	2.446	ng/ul	98
82) Pyrene	19.545	202	135129	2.076	ng/ul	96
86) Bis(2-ethylhexyl)phtha...	21.227	149	50650	1.506	ng/ul	100
87) Chrysene	21.344	228	62054	1.056	ng/ul	97
90) Benzo(b)fluoranthene	22.903	252	77333	1.552	ng/ul#	86

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

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