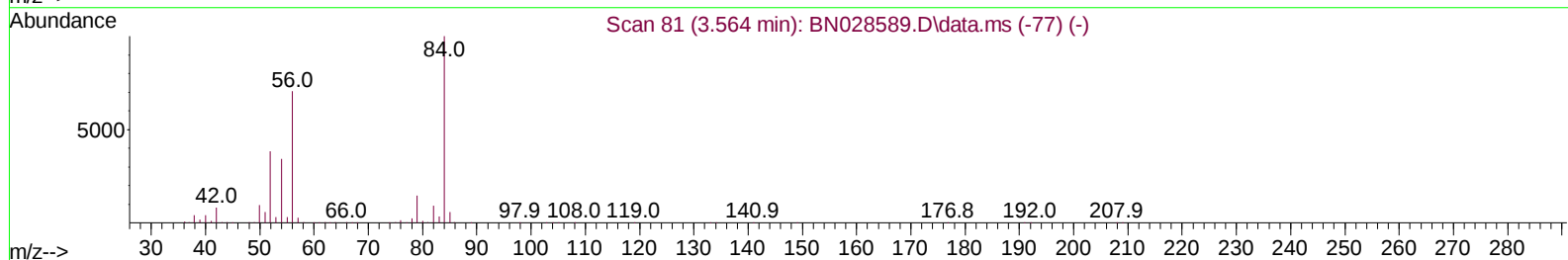
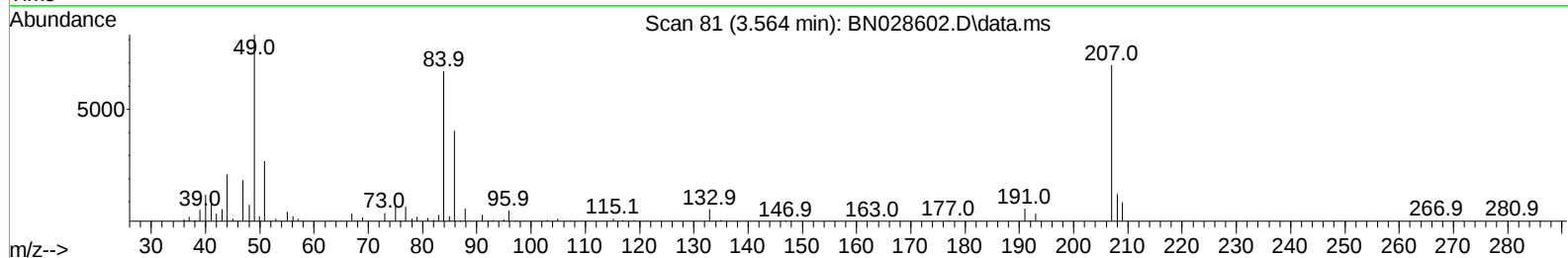
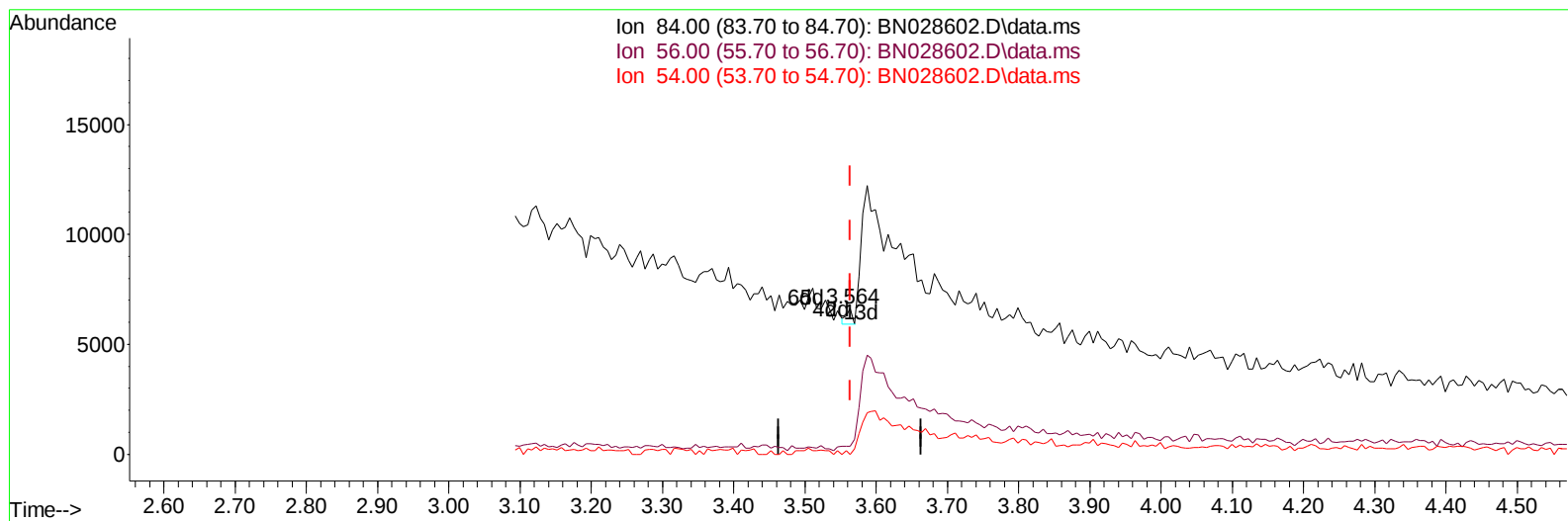


Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111423\
Data File : BN028602.D
Acq On : 14 Nov 2023 14:46
Operator : MA/JU
Sample : 05281-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

Quant Time: Nov 14 15:27:52 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN111423.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 14 04:50:19 2023
Response via : Initial Calibration



TIC: BN028602.D\data.ms

(4) Pyridine-d5 (S)

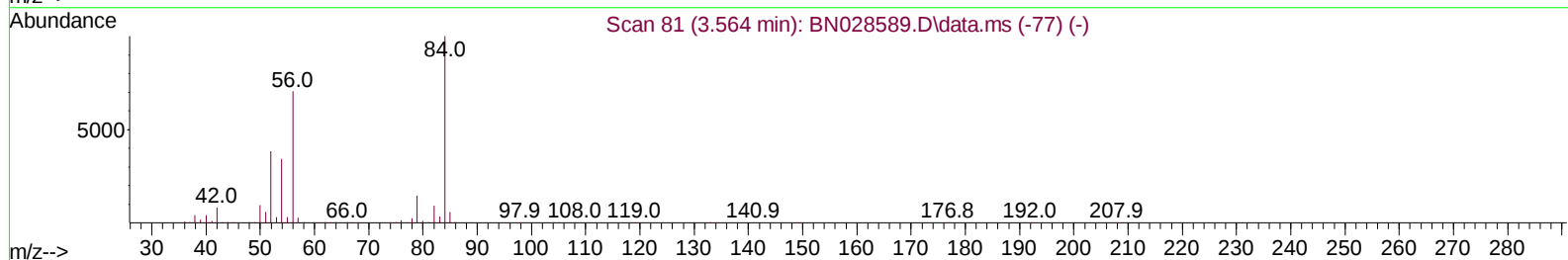
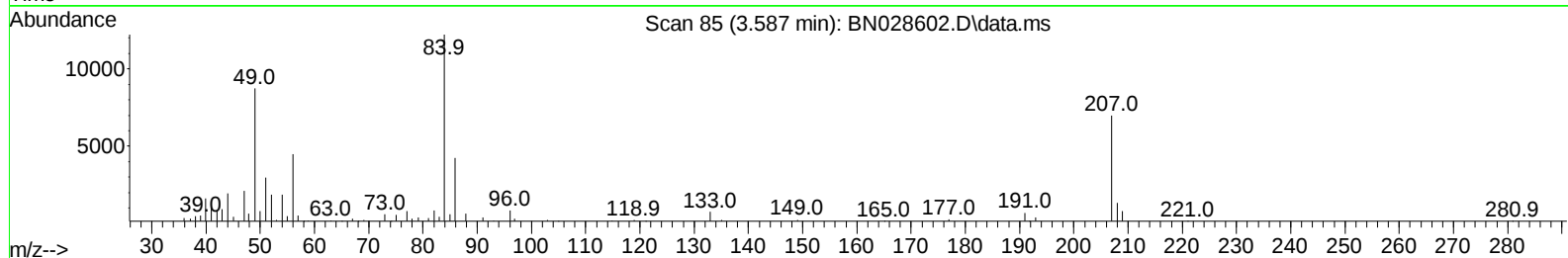
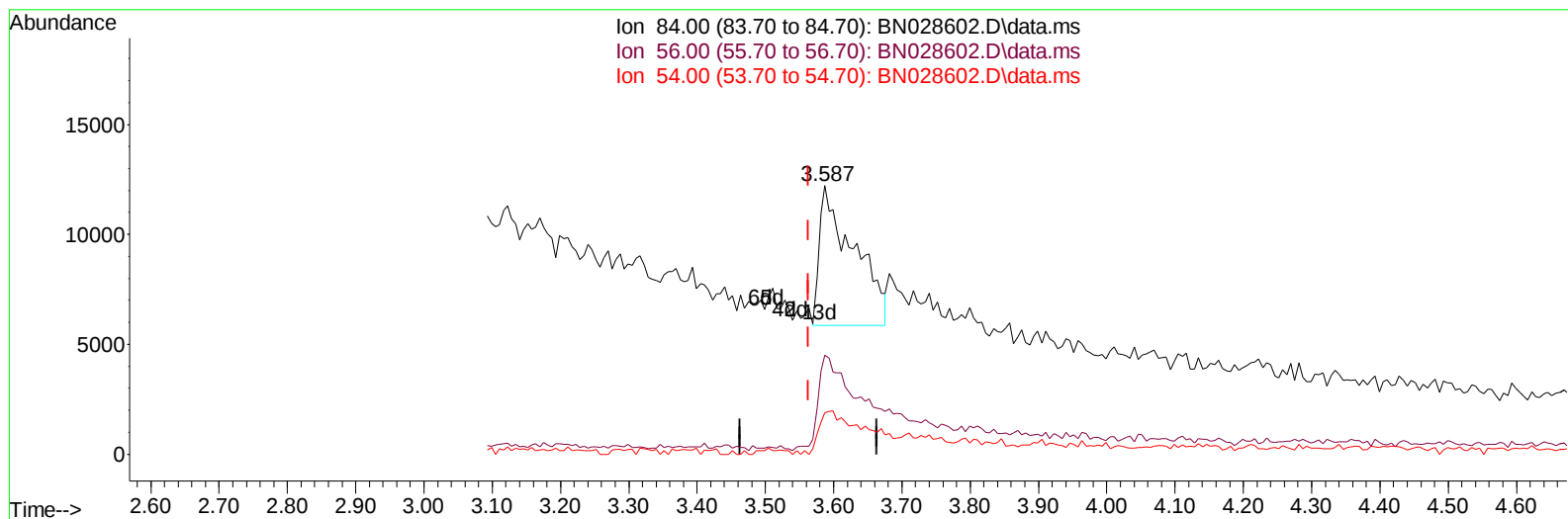
3.564min (-0.000) 0.05 ng/u1

response 396

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.50	5.67#
54.00	32.50	0.00#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111423\
Data File : BN028602.D
Acq On : 14 Nov 2023 14:46
Operator : MA/JU
Sample : 05281-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

Quant Time: Nov 14 15:27:52 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN111423.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 14 04:50:19 2023
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TIC: BN028602.D\data.ms

(4) Pyridine-d5 (S)

3.587min (+ 0.023) 2.95 ng/u1 m

response 22340

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.50	36.85#
54.00	32.50	15.44#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111423\
 Data File : BN028602.D
 Acq On : 14 Nov 2023 14:46
 Operator : MA/JU
 Sample : 05281-05
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Quant Time: Nov 14 15:28:25 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN111423.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 14 04:50:19 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.705	152	123790	20.000	ng/ul	0.00
20) Naphthalene-d8	10.487	136	723455	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.345	164	536164	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.092	188	1120918	20.000	ng/ul	0.00
79) Chrysene-d12	21.304	240	941157	20.000	ng/ul	0.00
88) Perylene-d12	23.604	264	993950	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.146	96	24242	7.012	ng/uL	0.00
4) Pyridine-d5	3.587	84	22340m	2.948	ng/ul	0.02
7) Phenol-d5	6.928	99	63766	5.392	ng/ul	0.04
9) Bis-(2-Chloroethyl)eth...	7.040	67	346122	47.319	ng/ul	0.00
11) 2-Chlorophenol-d4	7.234	132	385065	37.258	ng/ul	0.00
15) 4-Methylphenol-d8	8.434	113	170579	17.517	ng/ul	0.00
21) Nitrobenzene-d5	8.863	128	233449	35.940	ng/ul	0.00
24) 2-Nitrophenol-d4	9.581	143	251253	36.951	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.122	165	362149	31.325	ng/ul	0.00
31) 4-Chloroaniline-d4	10.657	131	294121	16.787	ng/ul	0.01
46) Dimethylphthalate-d6	13.763	166	2507412	51.866	ng/ul	0.00
49) Acenaphthylene-d8	14.034	160	2610532	46.851	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0	0.000	ng/ul	
60) Fluorene-d10	15.339	176	2068348	51.894	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.463	200	255145	40.977	ng/ul	0.00
73) Anthracene-d10	17.192	188	3378790	54.000	ng/ul	0.00
81) Pyrene-d10	19.498	212	4005770	56.118	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.463	264	3496816	56.197	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.181	88	9438	2.556	ng/uL#	94

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

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