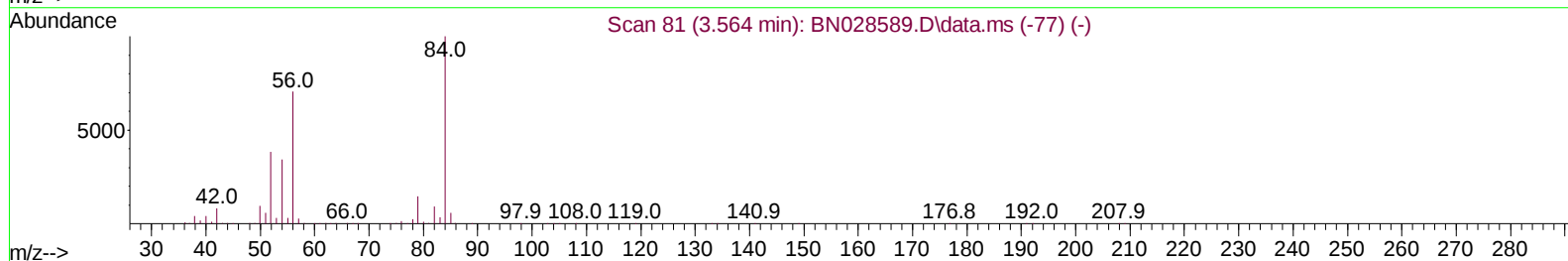
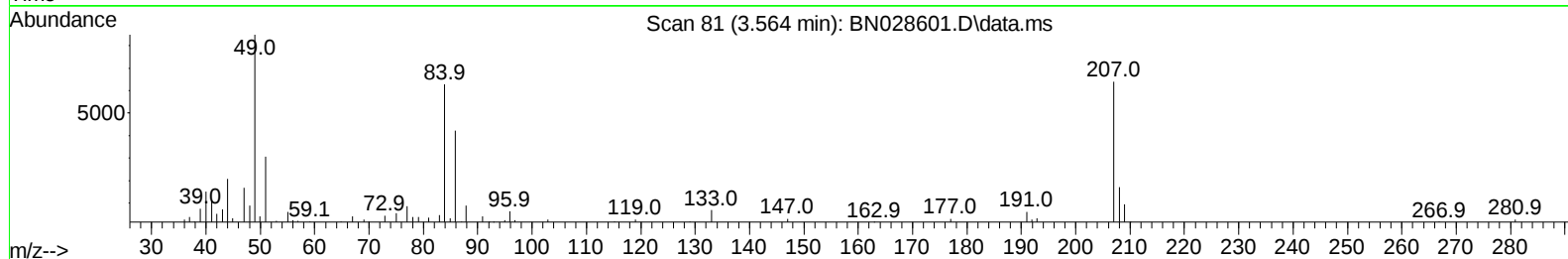
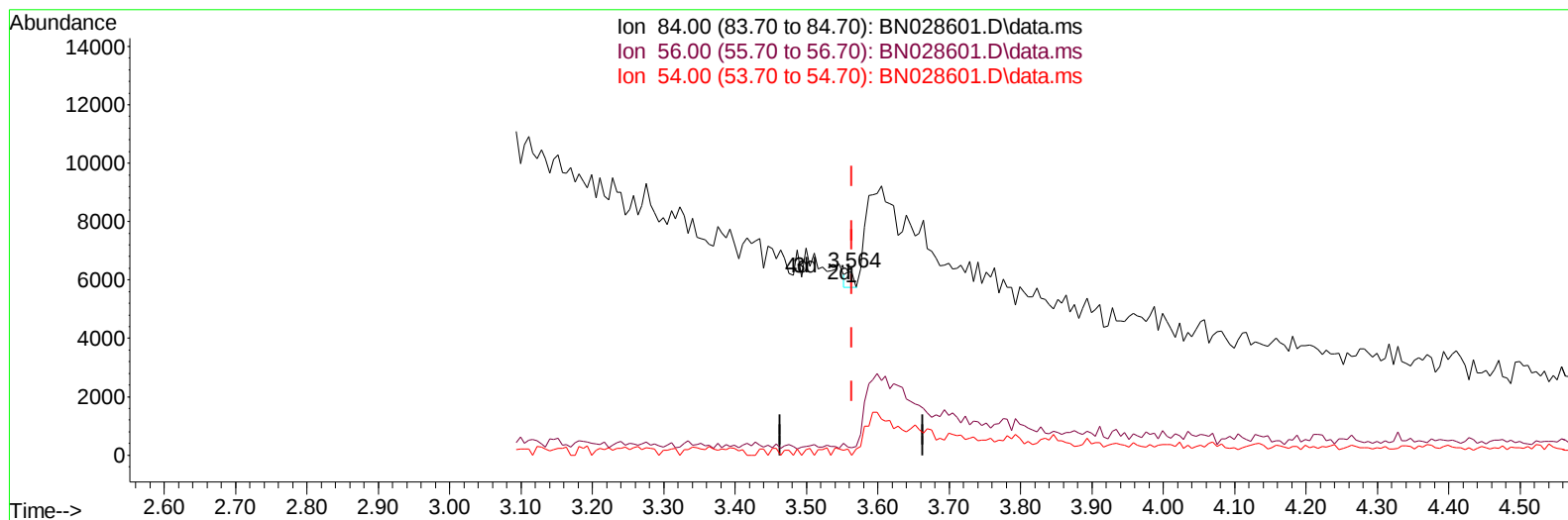


Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111423\
Data File : BN028601.D
Acq On : 14 Nov 2023 14:10
Operator : MA/JU
Sample : 05281-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Quant Time: Nov 14 15:25:26 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: BN028601.D\data.ms

(4) Pyridine-d5 (S)

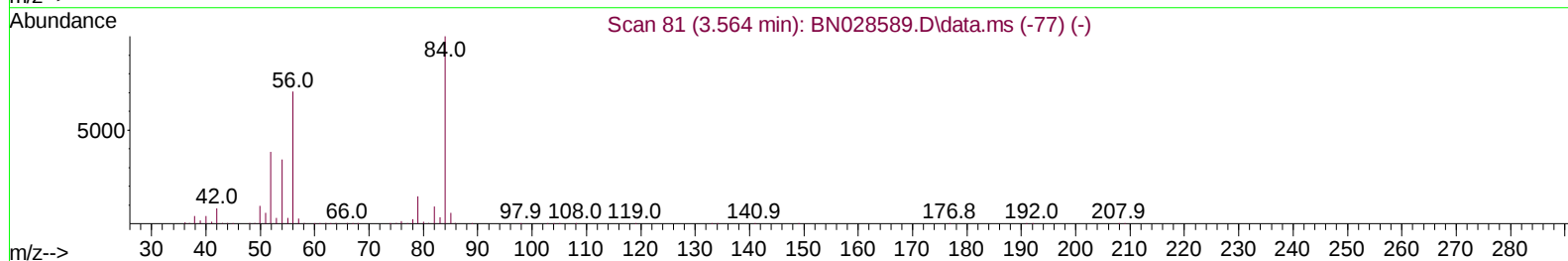
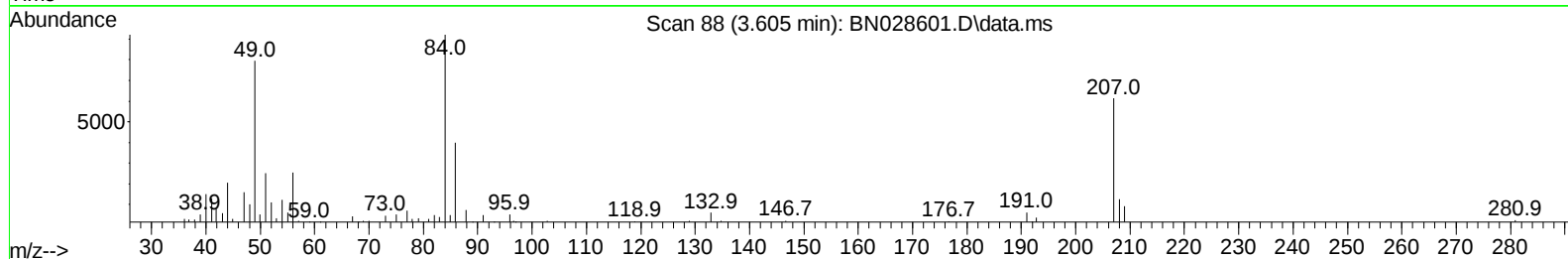
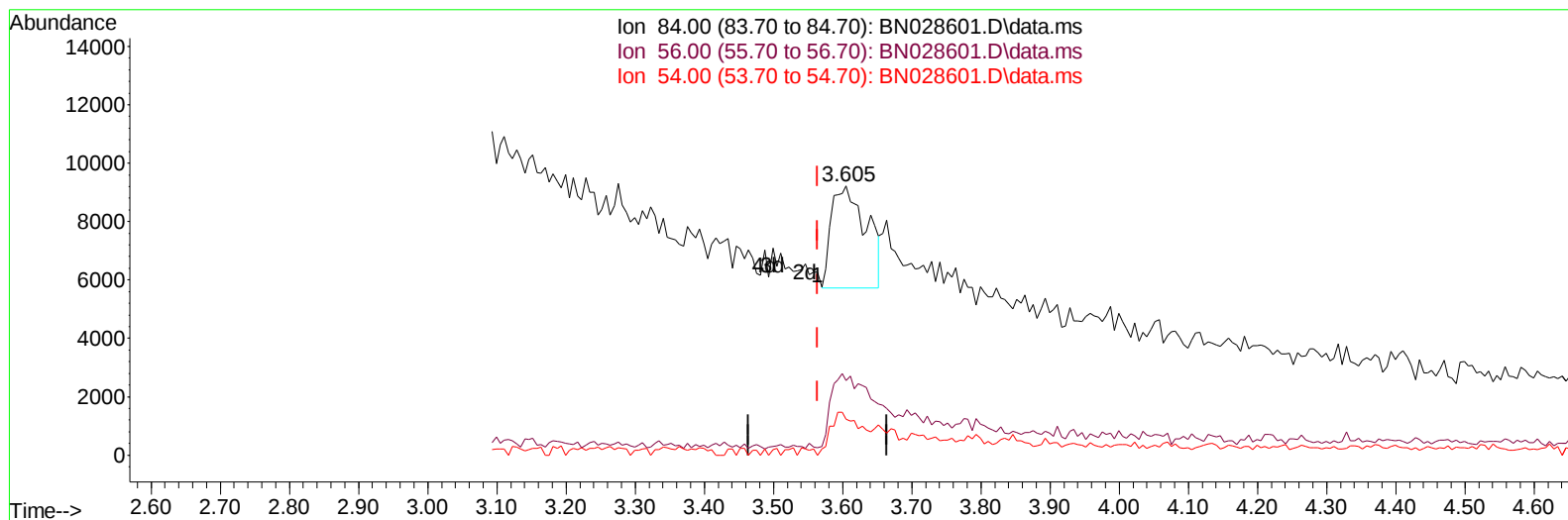
3.564min (+ 0.000) 0.03 ng/u1

response 367

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.50	3.94#
54.00	32.50	0.00#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111423\
Data File : BN028601.D
Acq On : 14 Nov 2023 14:10
Operator : MA/JU
Sample : 05281-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Quant Time: Nov 14 15:25:26 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: BN028601.D\data.ms

(4) Pyridine-d5 (S)

3.605min (+ 0.041) 1.07 ng/ul m

response 12185

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.50	27.68#
54.00	32.50	13.41#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111423\
 Data File : BN028601.D
 Acq On : 14 Nov 2023 14:10
 Operator : MA/JU
 Sample : 05281-04
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Quant Time: Nov 14 15:27:01 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	7.699	152	186448	20.000	ng/ul	0.00
20) Naphthalene-d8	10.487	136	1089753	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.345	164	804625	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.092	188	1710132	20.000	ng/ul	0.00
79) Chrysene-d12	21.304	240	1348516	20.000	ng/ul	0.00
88) Perylene-d12	23.604	264	1423839	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	23771	4.565	ng/uL	0.00
4) Pyridine-d5	3.605	84	12185m	1.067	ng/ul	0.04
7) Phenol-d5	6.934	99	50890	2.857	ng/ul	0.05
9) Bis-(2-Chloroethyl)eth...	7.040	67	285483	25.913	ng/ul	0.00
11) 2-Chlorophenol-d4	7.234	132	316896	20.358	ng/ul	0.00
15) 4-Methylphenol-d8	8.440	113	142224	9.697	ng/ul	0.01
21) Nitrobenzene-d5	8.863	128	206219	21.076	ng/ul	0.00
24) 2-Nitrophenol-d4	9.581	143	231912	22.643	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.122	165	307845	17.677	ng/ul	0.00
31) 4-Chloroaniline-d4	10.657	131	314814	11.928	ng/ul	0.01
46) Dimethylphthalate-d6	13.763	166	2161051	29.787	ng/ul	0.00
49) Acenaphthylene-d8	14.040	160	2319254	27.736	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.340	176	1811243	30.281	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.463	200	267511	28.160	ng/ul	0.00
73) Anthracene-d10	17.192	188	3388841	35.500	ng/ul	0.00
81) Pyrene-d10	19.498	212	4135398	40.433	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.463	264	3504055	39.311	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.181	88	9175	1.650	ng/uL	91

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

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