

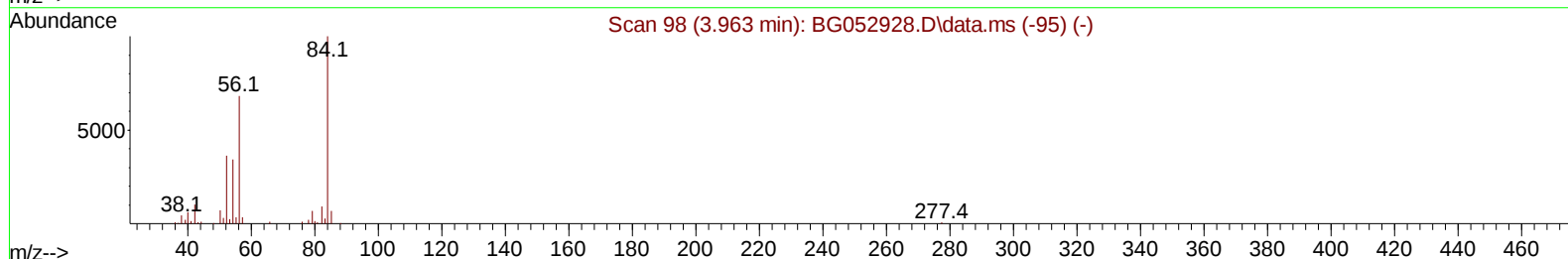
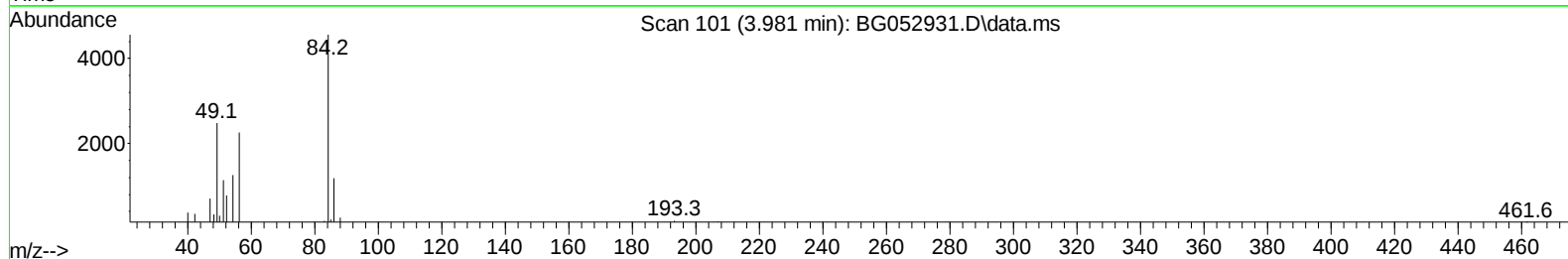
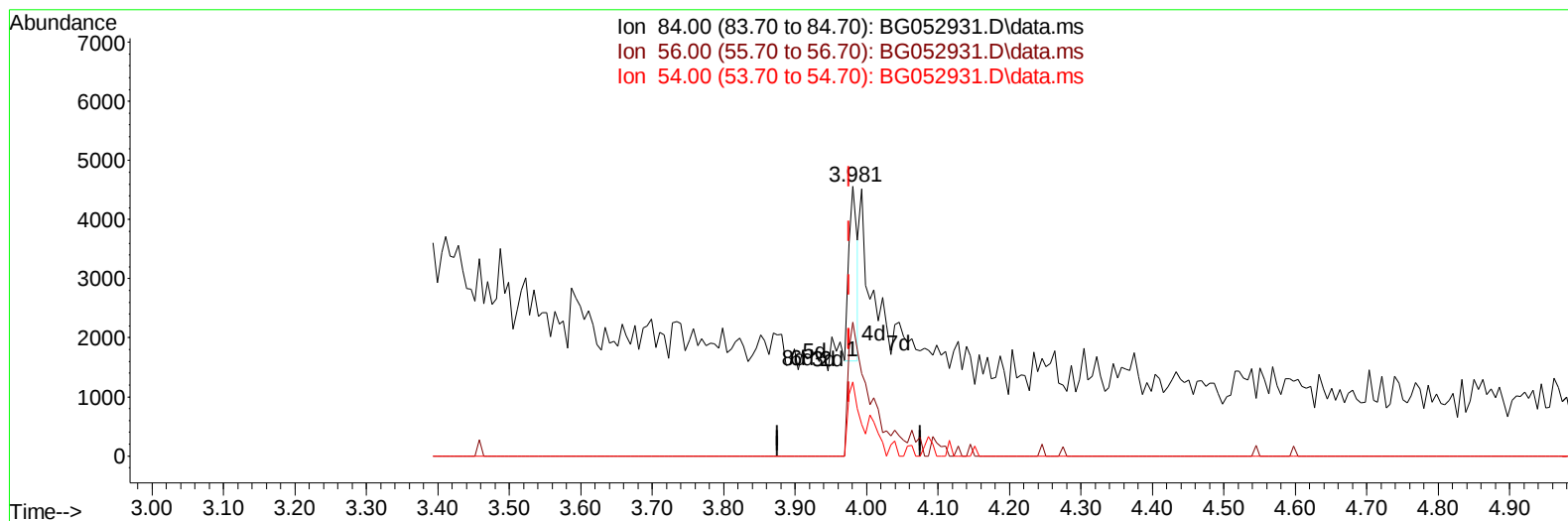
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032822\
 Data File : BG052931.D
 Acq On : 29 Mar 2022 17:44
 Operator : CG/JU
 Sample : N2097-18
 Misc :
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AE8

Manual IntegrationsAPPROVED

Quant Time: Mar 30 01:10:18 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG031822.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Mar 23 13:16:32 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/30/2022
 Supervised By :Yogesh Patel 04/02/2022



TIC: BG052931.D\data.ms

(4) Pyridine-d5 (S)

3.981min (+ 0.006) 1.35 ng/u1

response 2450

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	49.68#
54.00	28.70	27.57
0.00	0.00	0.00

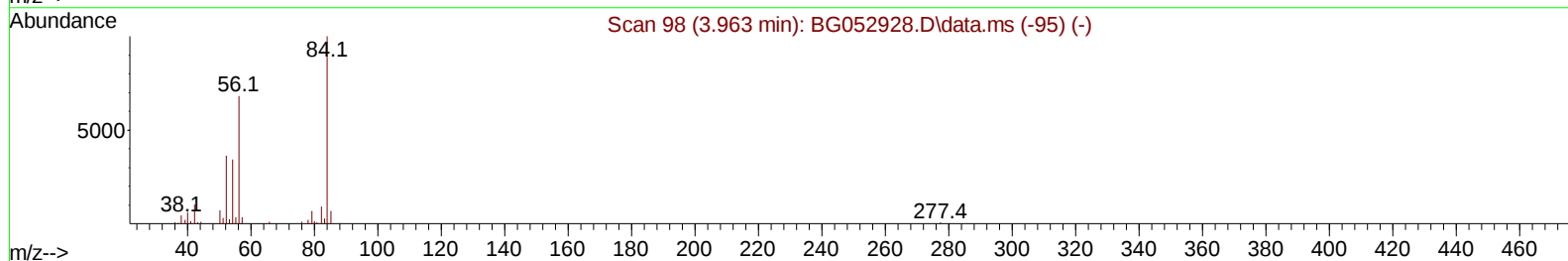
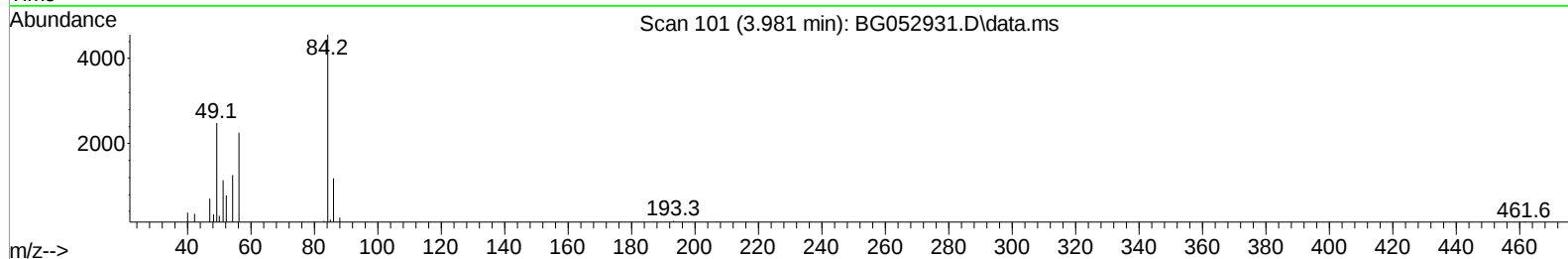
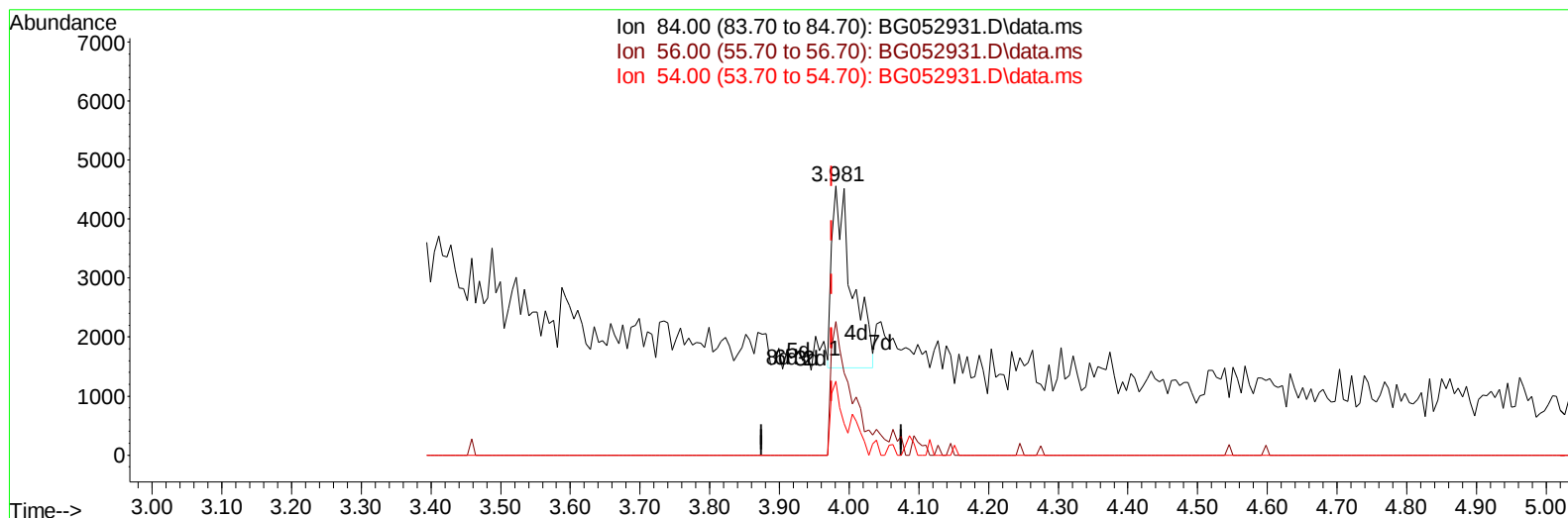
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TIC: BG052931.D\data.ms

(4) Pyridine-d5 (S)

3.981min (+ 0.006) 3.36 ng/u1 m

response 6084

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	49.68#
54.00	28.70	27.57
0.00	0.00	0.00

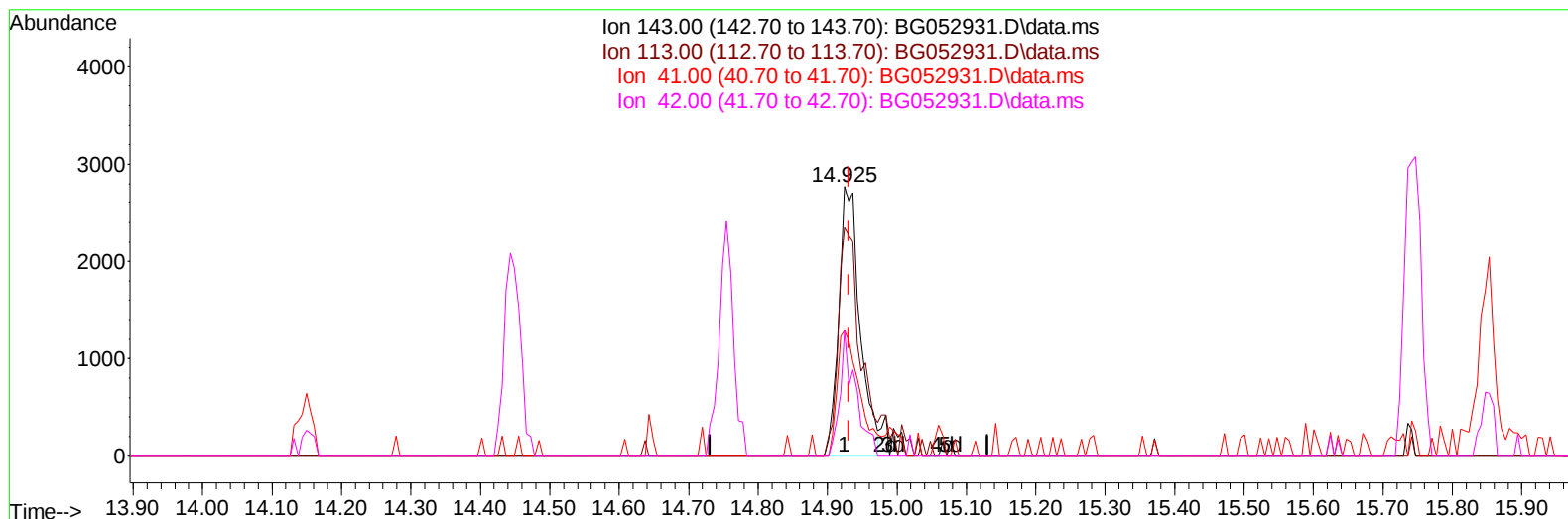
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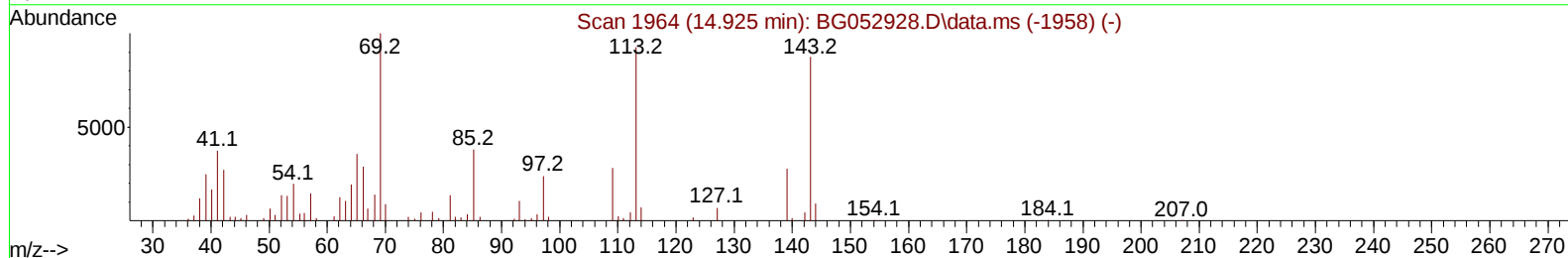
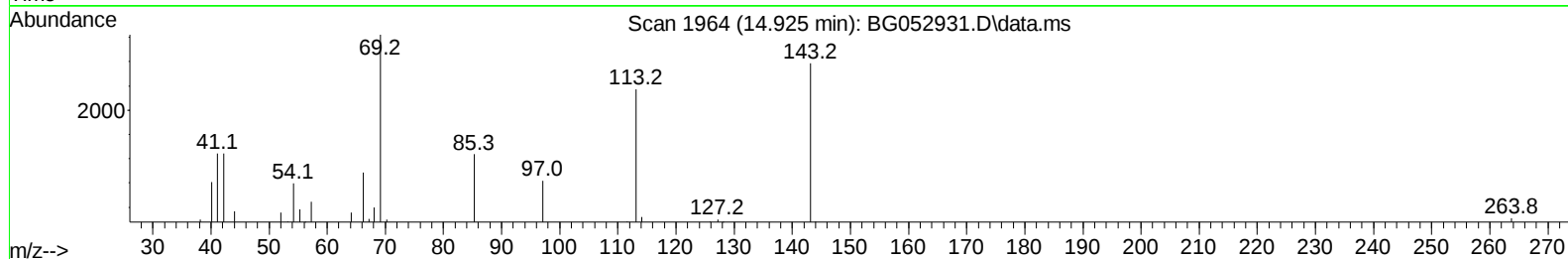
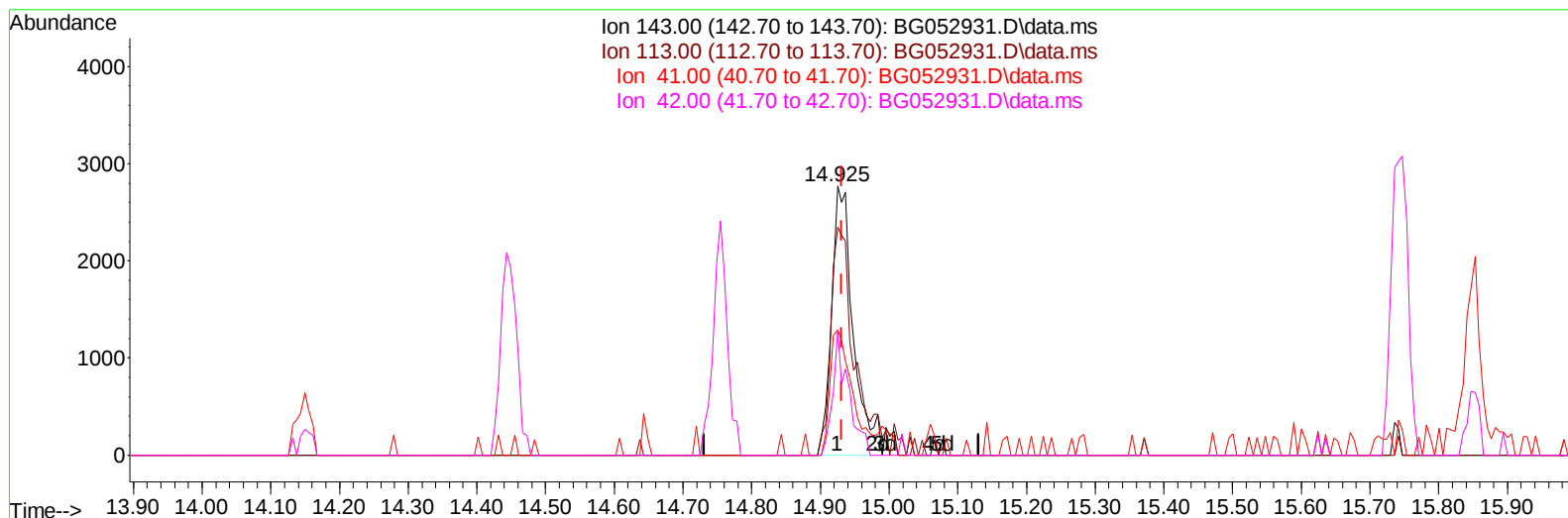
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TIC: BG052931.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.925min (-0.006) 4.85 ng/ul m

response 6058

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	84.80
41.00	47.20	46.51
42.00	29.40	46.48#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.140	152	25637	20.000	ng/ul	0.00
20) Naphthalene-d8	10.948	136	116909	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.755	164	96793	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.498	188	228084	20.000	ng/ul	0.00
79) Chrysene-d12	21.780	240	249854	20.000	ng/ul	#-0.01
88) Perylene-d12	25.093	264	251425	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.552	96	2472	3.406	ng/uL	0.00
4) Pyridine-d5	3.981	84	6084m	3.358	ng/ul	0.00
7) Phenol-d5	7.282	99	13992	6.094	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.453	67	43472	30.998	ng/ul	0.00
11) 2-Chlorophenol-d4	7.664	132	34179	21.922	ng/ul	-0.01
15) 4-Methylphenol-d8	8.827	113	24639	14.813	ng/ul	0.00
21) Nitrobenzene-d5	9.291	128	25581	30.397	ng/ul	0.00
24) 2-Nitrophenol-d4	10.020	143	26298	29.627	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.560	165	48678	25.208	ng/ul	0.00
31) 4-Chloroaniline-d4	11.071	131	61406	24.060	ng/ul	0.00
46) Dimethylphthalate-d6	14.149	166	217461	29.244	ng/ul	-0.01
49) Acenaphthylene-d8	14.449	160	241363	26.712	ng/ul	0.00
54) 4-Nitrophenol-d4	14.925	143	6058m	4.845	ng/ul	0.00
60) Fluorene-d10	15.741	176	197722	30.283	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.853	200	32437	27.946	ng/ul	0.00
73) Anthracene-d10	17.598	188	345689	32.548	ng/ul	0.00
81) Pyrene-d10	19.877	212	415270	29.710	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.864	264	419603	32.299	ng/ul	-0.01

Target Compounds Qvalue

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

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