

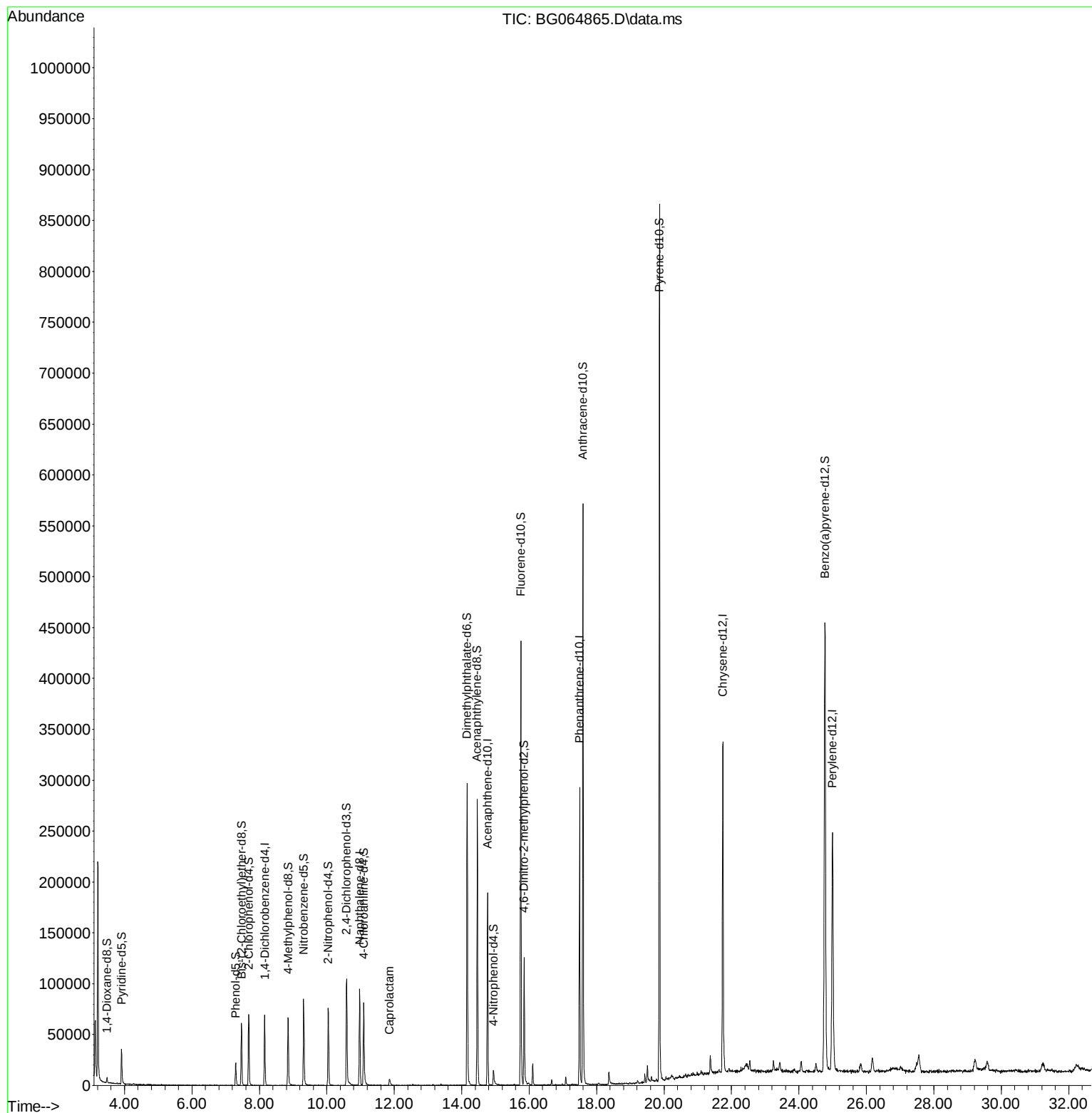
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112525\
Data File : BG064865.D
Acq On : 25 Nov 2025 17:58
Operator : RC/JU
Sample : Q3688-10
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
M-32S

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/26/2025
Supervised By :mohammad ahmed 12/03/2025

Quant Time: Nov 26 02:05:24 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112425.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 25 10:23:09 2025
Response via : Initial Calibration



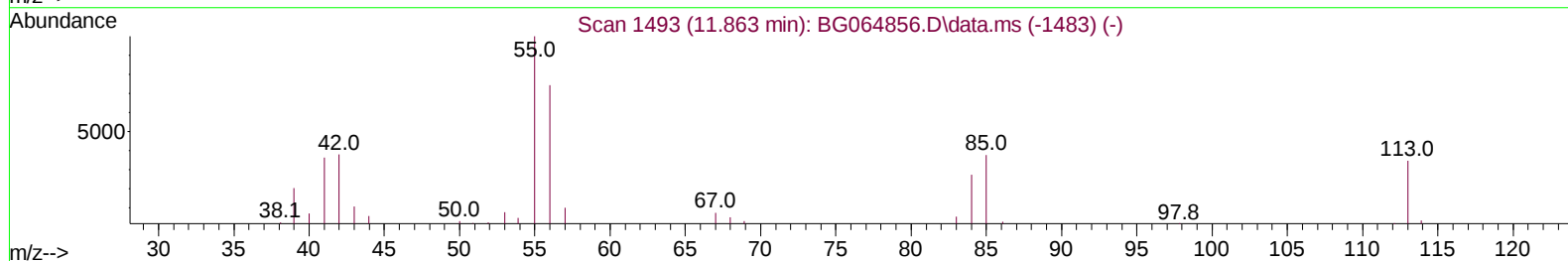
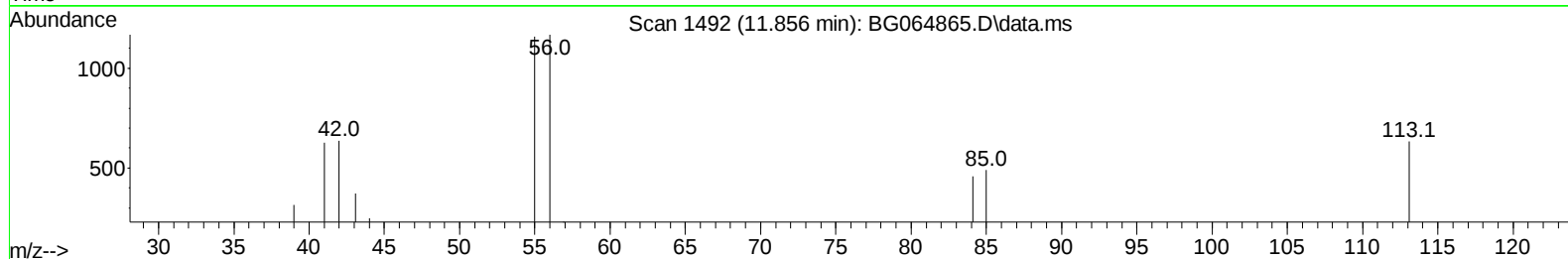
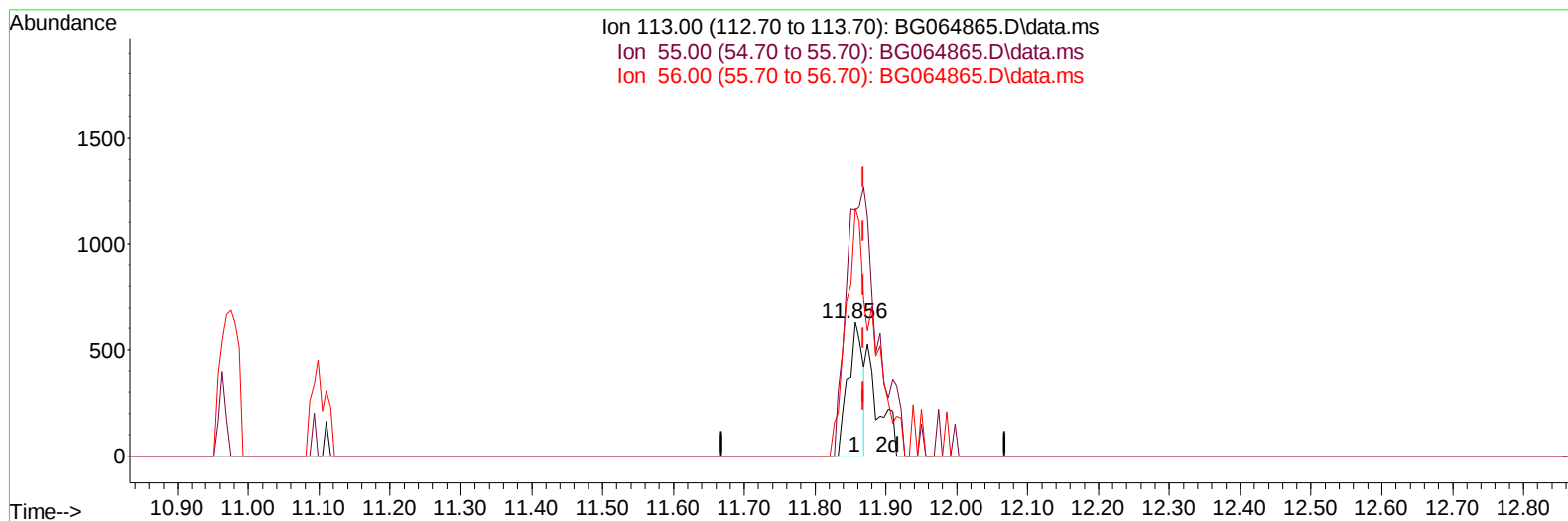
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Quant Time: Nov 26 07:06:26 2025
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TIC: BG064865.D\data.ms

(34) Caprolactam

11.856min (-0.012) 1.75 ng/u1

response 894

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	234.90	183.10#
56.00	159.40	184.52
0.00	0.00	0.00

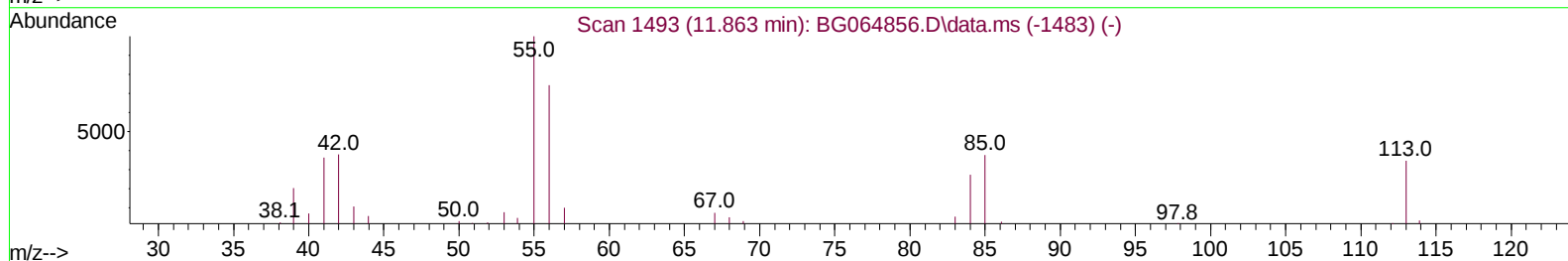
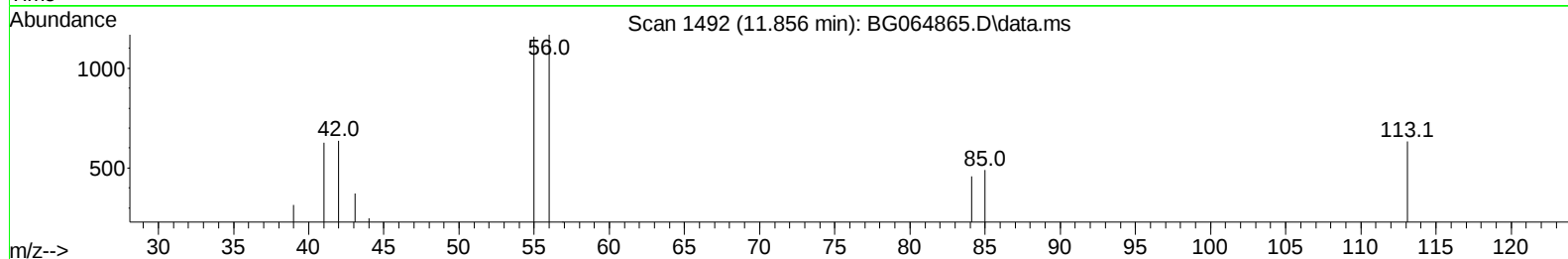
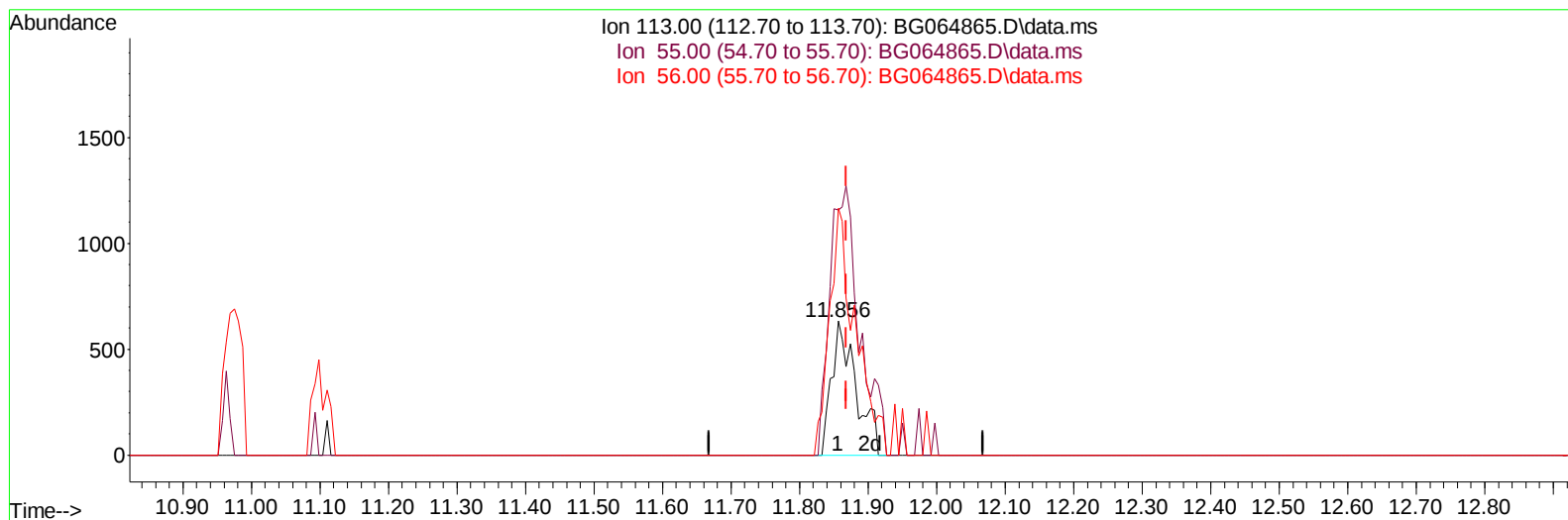
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TIC: BG064865.D\data.ms

(34) Caprolactam

11.856min (-0.012) 3.06 ng/ul m

response 1562

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	234.90	183.10#
56.00	159.40	184.52
0.00	0.00	0.00

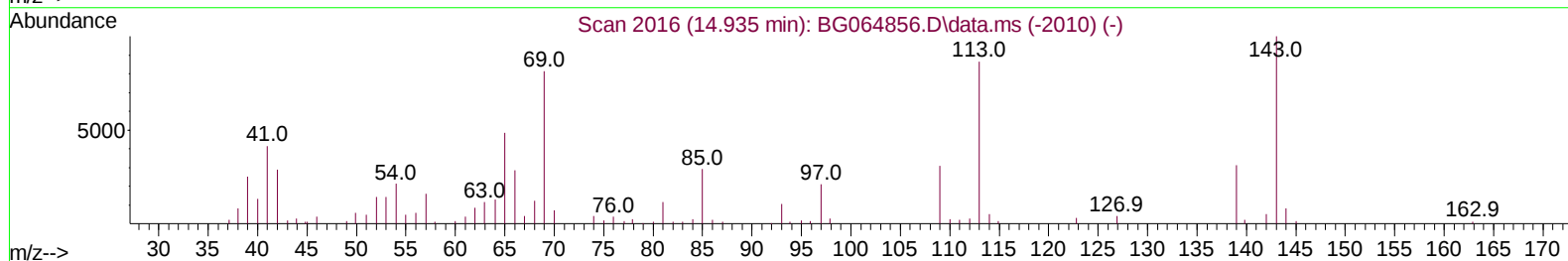
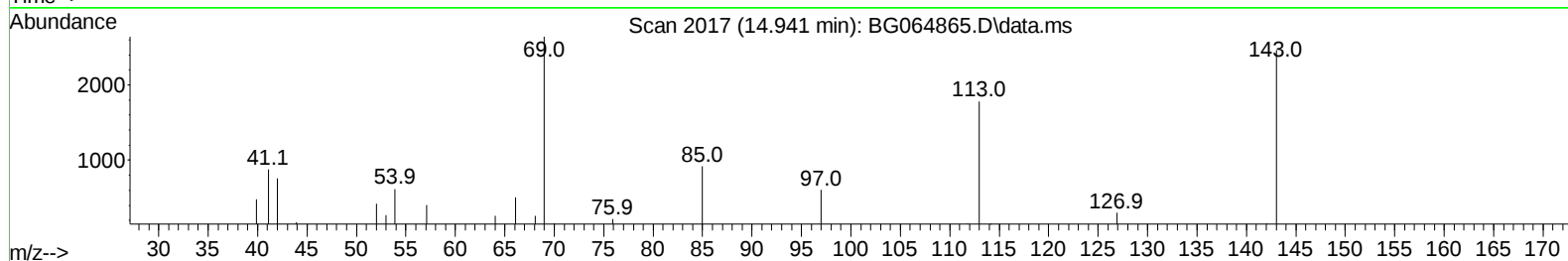
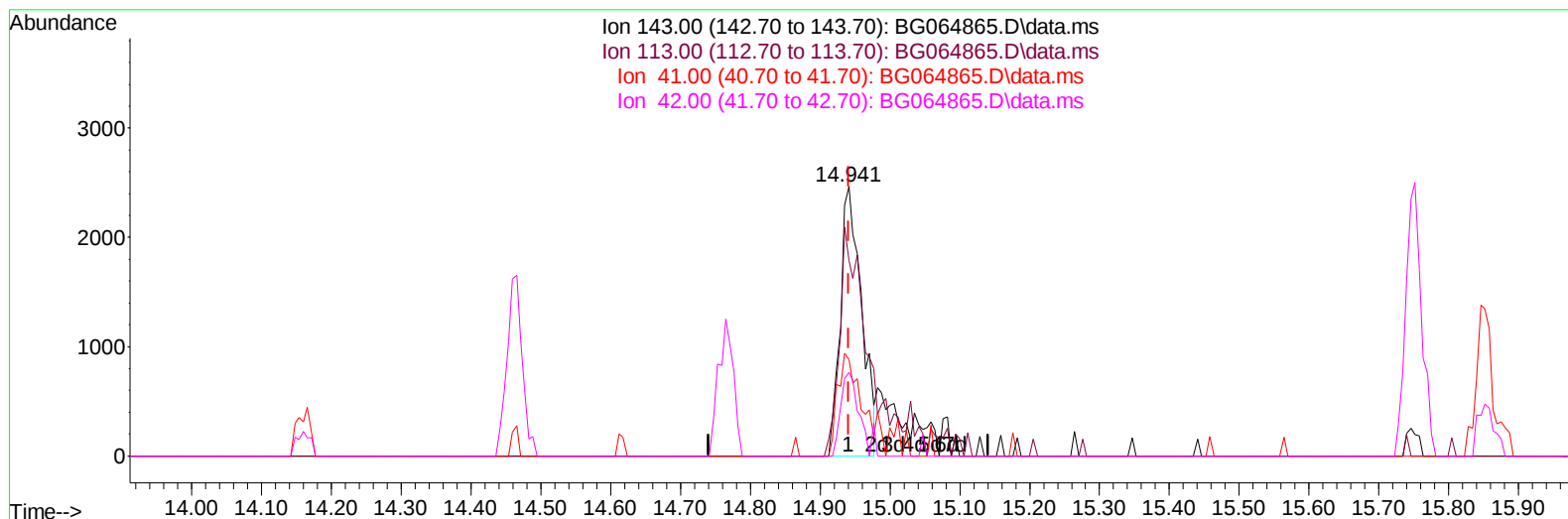
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TIC: BG064865.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.941min (+ 0.000) 5.69 ng/u1

response 5160

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	90.10	72.38
41.00	40.40	35.64
42.00	23.60	30.94#

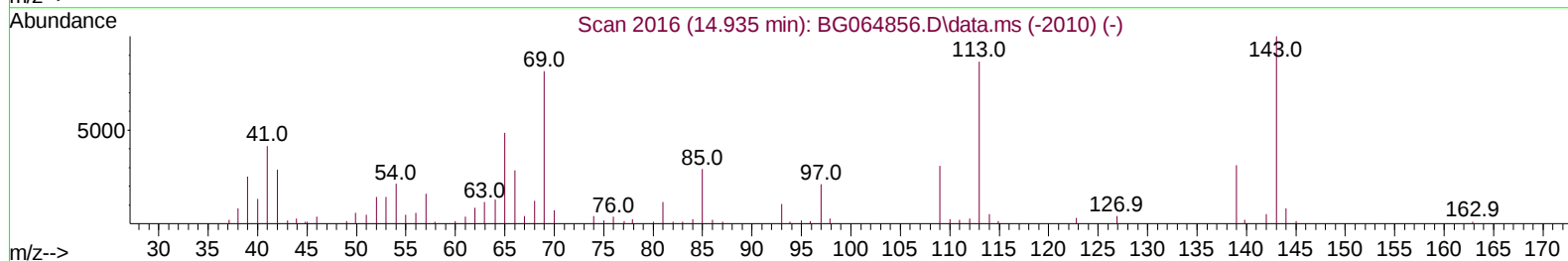
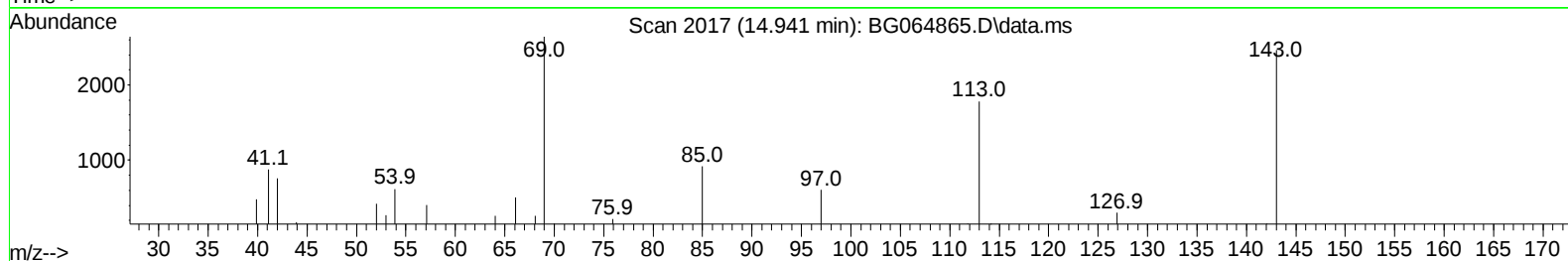
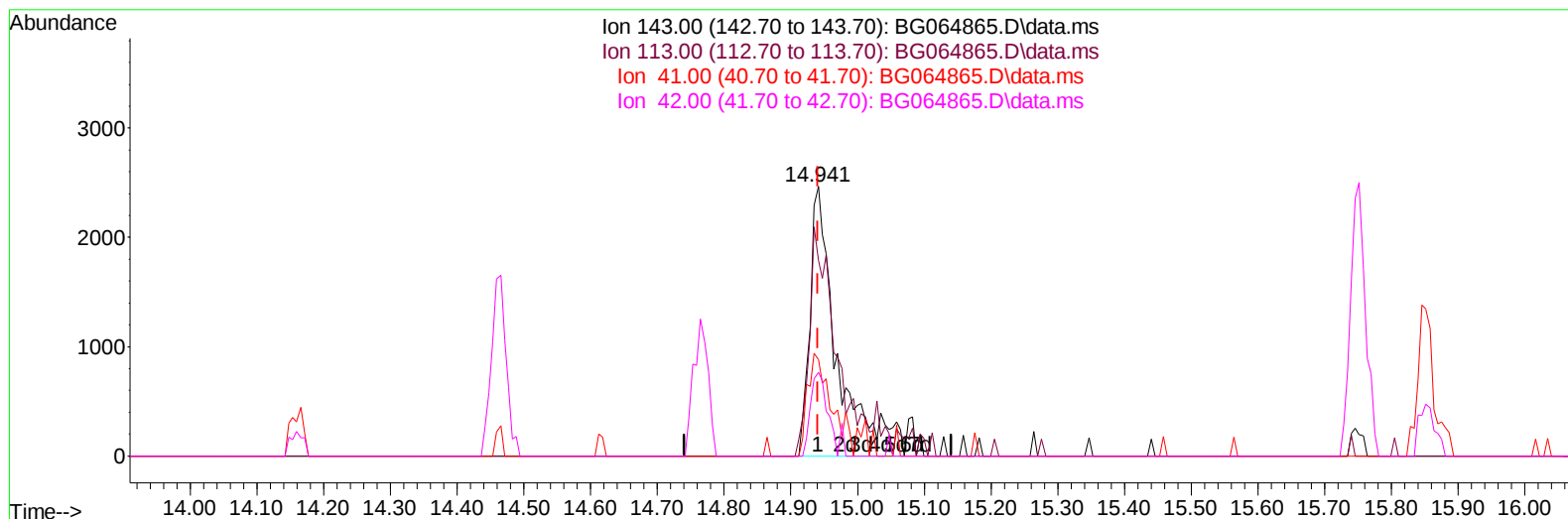
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Instrument :
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TIC: BG064865.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.941min (+ 0.000) 7.76 ng/ul m

response 7043

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	90.10	72.38
41.00	40.40	35.64
42.00	23.60	30.94#

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 ALS Vial : 11 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.161	152	20582	20.000	ng/ul	0.00
20) Naphthalene-d8	10.975	136	84782	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.765	164	76060	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.497	188	198464	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.745	240	248527	20.000	ng/ul	# 0.00
88) Perylene-d12	24.994	264	274864	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.484	96	2415	4.651	ng/uL	0.00
4) Pyridine-d5	3.913	84	16676	11.655	ng/ul	0.00
7) Phenol-d5	7.297	99	13441	8.101	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.467	67	30685	28.917	ng/ul	0.00
11) 2-Chlorophenol-d4	7.685	132	33957	26.996	ng/ul	0.00
15) 4-Methylphenol-d8	8.854	113	27171	19.001	ng/ul	0.00
21) Nitrobenzene-d5	9.312	128	20859	32.619	ng/ul	0.00
24) 2-Nitrophenol-d4	10.041	143	25211	33.193	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.587	165	46744	29.850	ng/ul	0.00
31) 4-Chloroaniline-d4	11.098	131	47122	23.542	ng/ul	0.00
46) Dimethylphthalate-d6	14.160	166	231100	38.626	ng/ul	0.00
49) Acenaphthylene-d8	14.459	160	212050	33.094	ng/ul	0.00
54) 4-Nitrophenol-d4	14.941	143	7043m	7.765	ng/ul	0.00
60) Fluorene-d10	15.752	176	199631	37.167	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.852	200	39190	30.843	ng/ul	0.00
73) Anthracene-d10	17.597	188	398890	39.776	ng/ul	0.00
81) Pyrene-d10	19.865	212	529796	41.501	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.771	264	581966	40.131	ng/ul	0.00
Target Compounds						
34) Caprolactam	11.856	113	1562m	3.061	ng/ul	Qvalue

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