

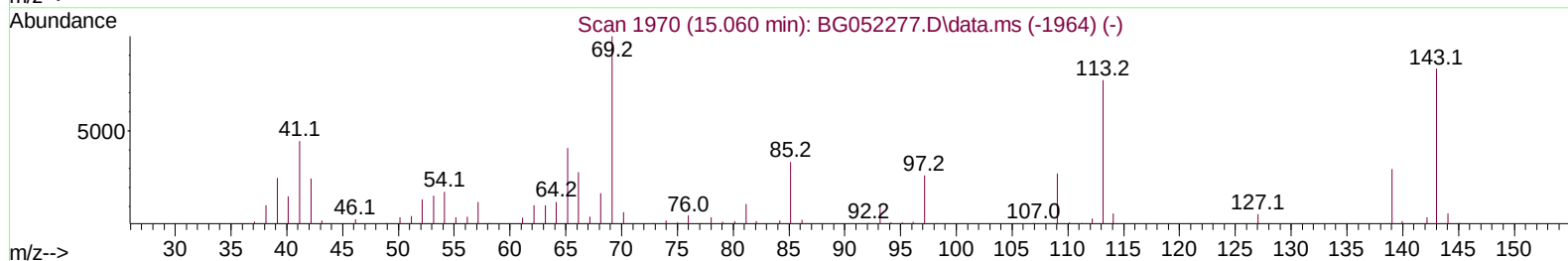
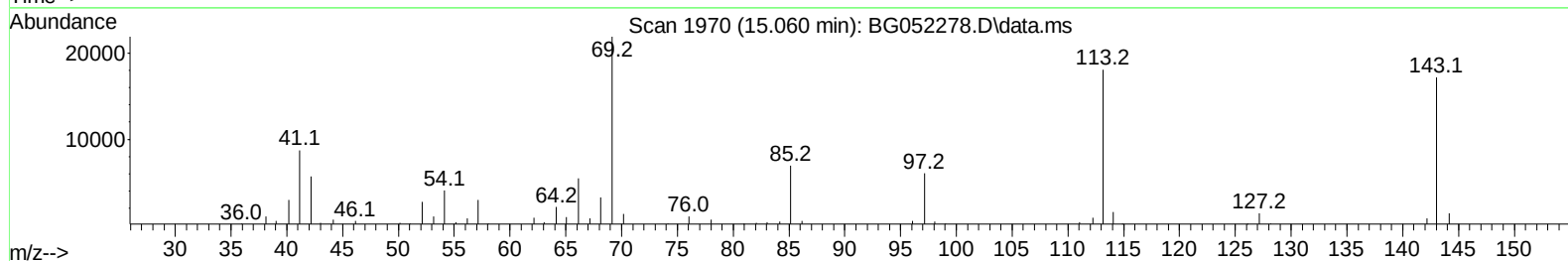
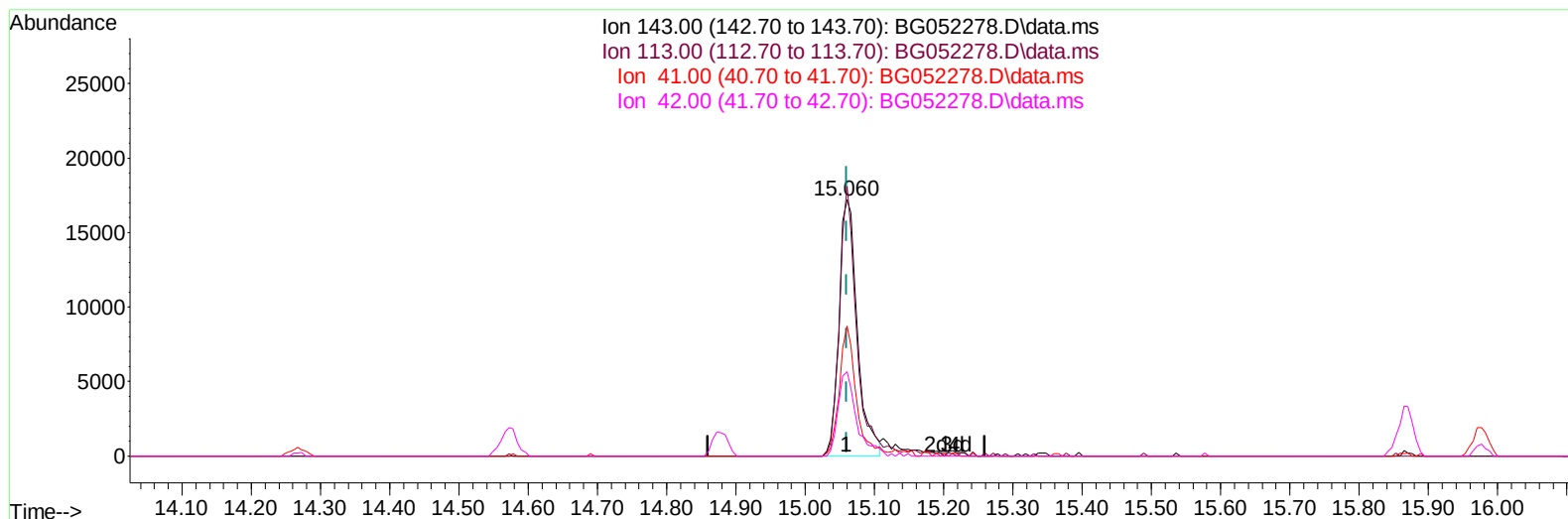
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG020322\
 Data File : BG052278.D
 Acq On : 3 Feb 2022 21:35
 Operator : CG/JU
 Sample : PB142431BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK431

Manual IntegrationsAPPROVED

Quant Time: Feb 04 00:02:55 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG013122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jan 31 18:07:10 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 02/04/2022
 Supervised By :Yogesh Patel 02/08/2022



TIC: BG052278.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.060min (0.000) 33.01 ng/ul

response 31747

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	105.21#
41.00	44.40	50.82
42.00	29.70	33.00

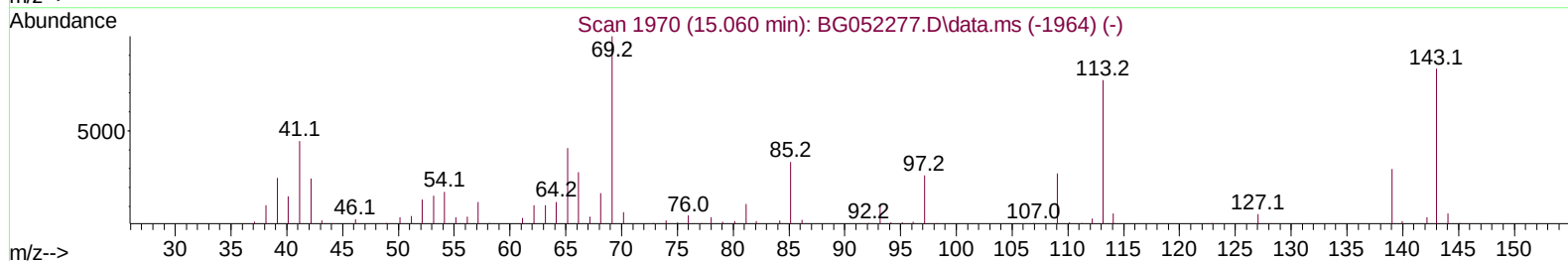
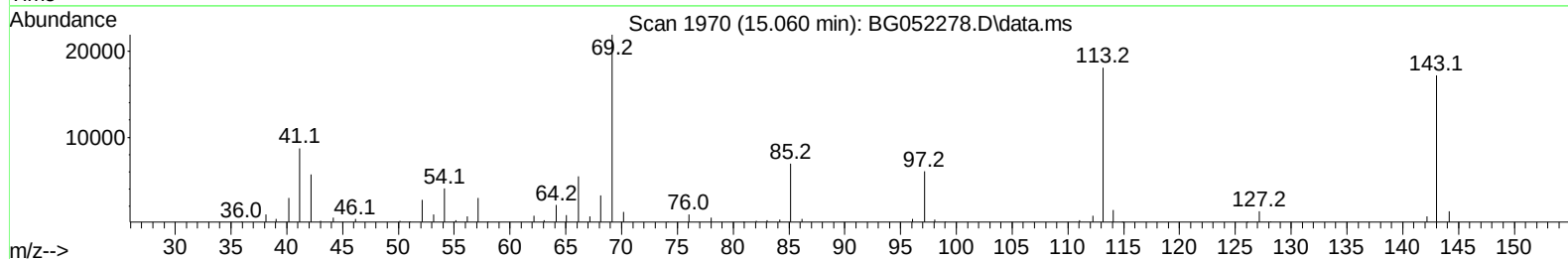
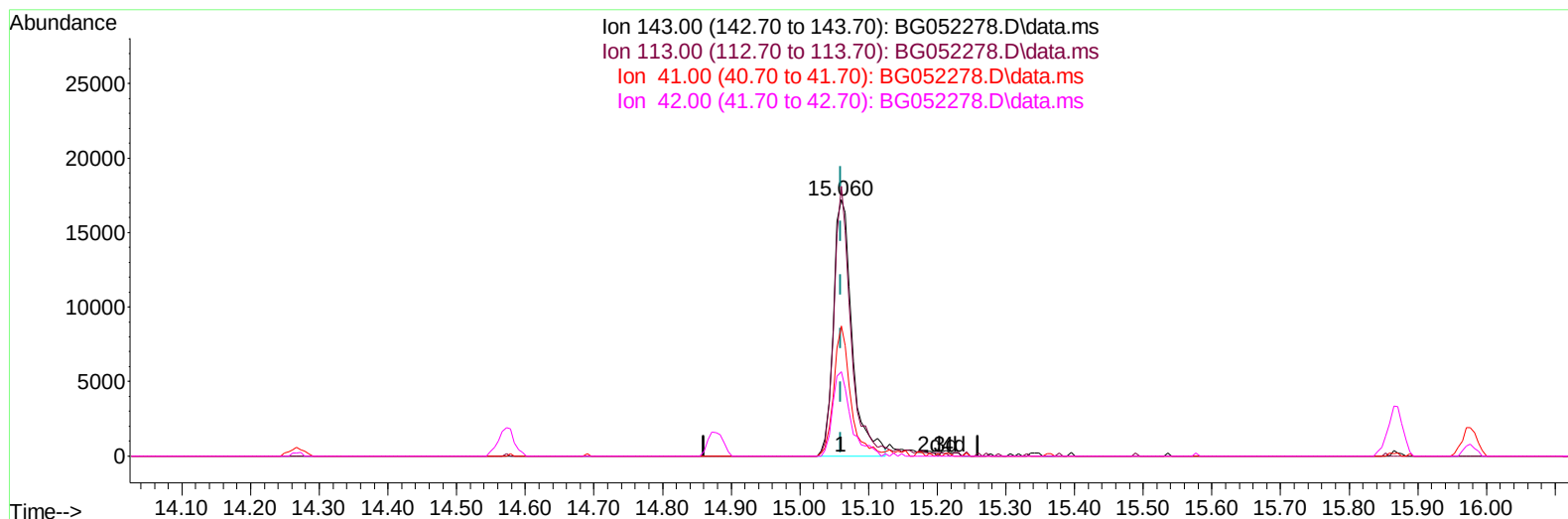
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TIC: BG052278.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.060min (0.000) 33.92 ng/ul m

response 32627

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	105.21#
41.00	44.40	50.82
42.00	29.70	33.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.240	152	29586	20.000	ng/ul	0.00
20) Naphthalene-d8	11.071	136	115559	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.878	164	80791	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.627	188	198367	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.939	240	212810	20.000	ng/ul	# 0.00
88) Perylene-d12	25.411	264	209386	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.564	96	4778	6.515	ng/uL	0.00
4) Pyridine-d5	3.993	84	69550	30.491	ng/ul	0.00
7) Phenol-d5	7.376	99	82611	30.697	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.541	67	52378	32.037	ng/ul	0.00
11) 2-Chlorophenol-d4	7.764	132	60496	32.260	ng/ul	0.00
15) 4-Methylphenol-d8	8.939	113	65852	32.246	ng/ul	0.00
21) Nitrobenzene-d5	9.403	128	30736	32.666	ng/ul	0.00
24) 2-Nitrophenol-d4	10.137	143	34908	33.951	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.684	165	64298	32.078	ng/ul	0.00
31) 4-Chloroaniline-d4	11.201	131	88074	34.641	ng/ul	0.00
46) Dimethylphthalate-d6	14.267	166	230610	36.314	ng/ul	0.00
49) Acenaphthylene-d8	14.573	160	283795	35.854	ng/ul	0.00
54) 4-Nitrophenol-d4	15.060	143	32627m	33.923	ng/ul	0.00
60) Fluorene-d10	15.871	176	203214	36.526	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.976	200	36242	29.594	ng/ul	0.00
73) Anthracene-d10	17.727	188	350782	37.509	ng/ul	0.00
81) Pyrene-d10	20.006	212	446652	35.020	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.176	264	422409	37.973	ng/ul	0.00

Target Compounds Qvalue

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

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