

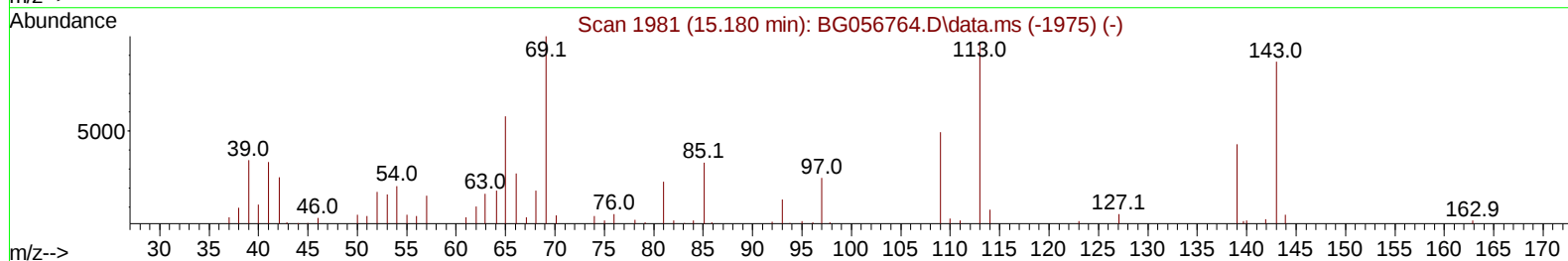
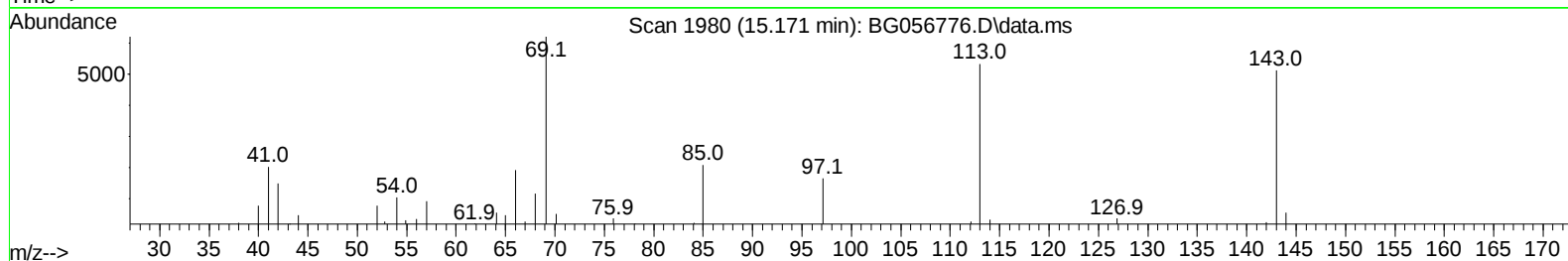
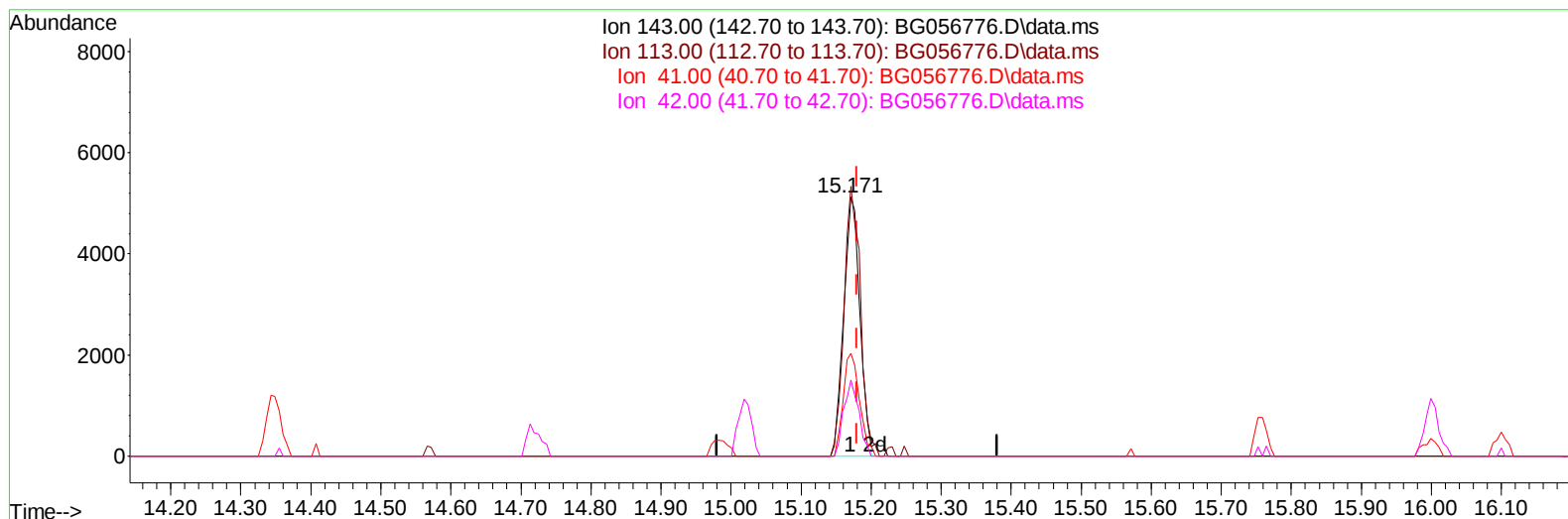
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030123\
 Data File : BG056776.D
 Acq On : 1 Mar 2023 23:19
 Operator : CG/JU
 Sample : 01649-14
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DBZW1

Manual IntegrationsAPPROVED

Quant Time: Mar 02 00:16:35 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Mar 02 00:12:28 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 03/02/2023
 Supervised By :Jagrut Upadhyay 03/02/2023



TIC: BG056776.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.171min (-0.009) 11.14 ng/u1

response 8327

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.30	104.12
41.00	46.10	39.64
42.00	27.10	29.42

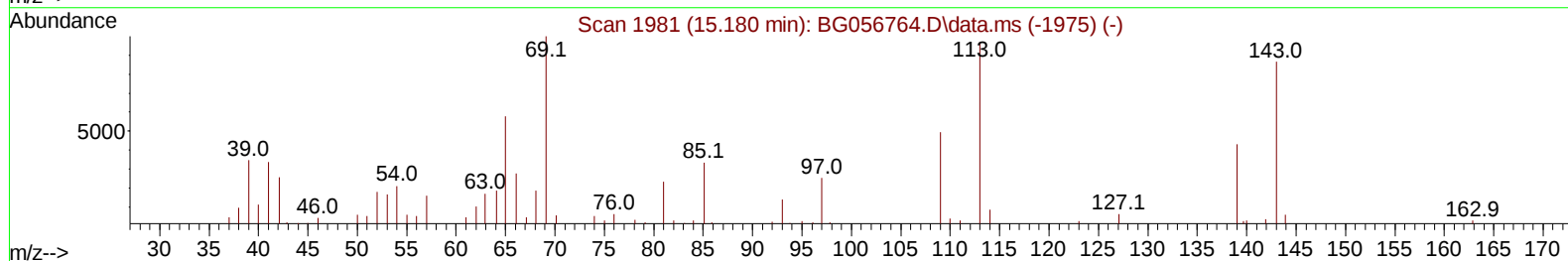
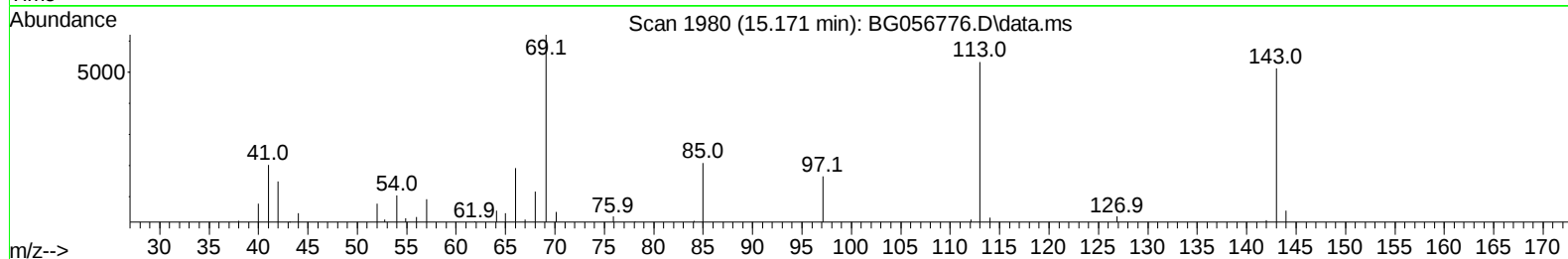
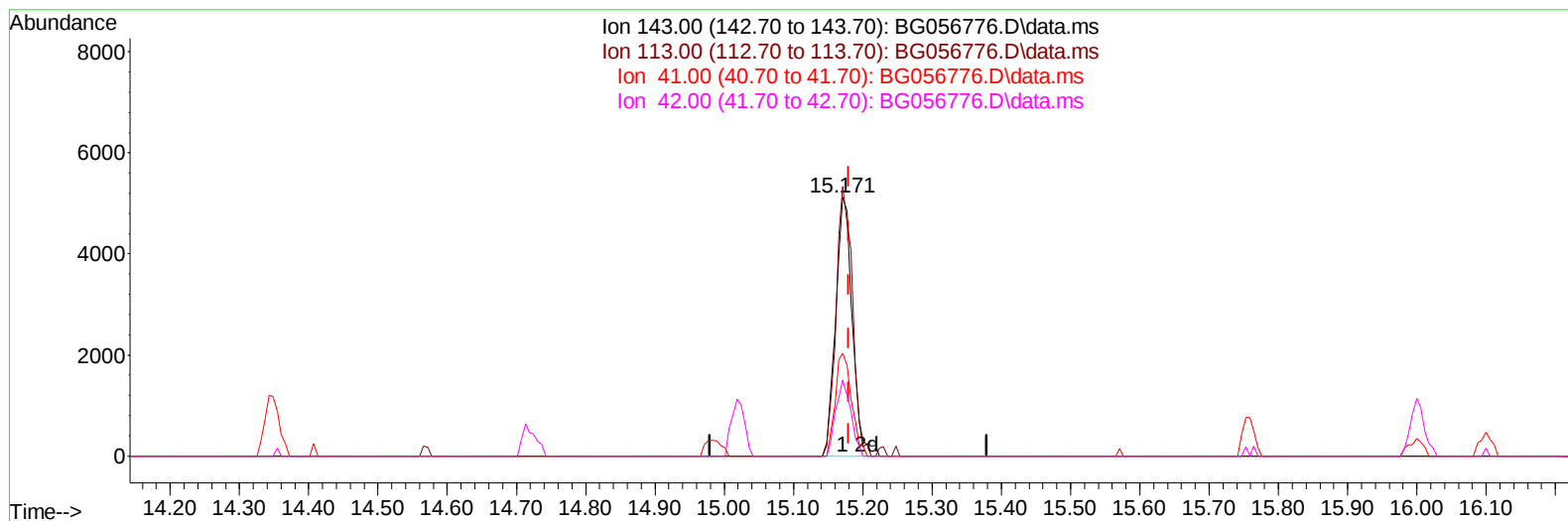
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TIC: BG056776.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.171min (-0.009) 11.45 ng/ul m

response 8556

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.30	104.12
41.00	46.10	39.64
42.00	27.10	29.42

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.414	152	15286	20.000	ng/ul	0.00
20) Naphthalene-d8	11.258	136	65221	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	15.018	164	53727	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.750	188	131971	20.000	ng/ul	0.00
79) Chrysene-d12	22.057	240	125871	20.000	ng/ul	0.00
88) Perylene-d12	25.582	264	134491	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.702	96	1029	3.734	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.539	99	19734	12.901	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.721	67	12297	13.179	ng/ul	0.00
11) 2-Chlorophenol-d4	7.938	132	13339	13.131	ng/ul	0.00
15) 4-Methylphenol-d8	9.102	113	12975	11.101	ng/ul	0.00
21) Nitrobenzene-d5	9.584	128	7442	13.616	ng/ul	0.00
24) 2-Nitrophenol-d4	10.318	143	8933	13.730	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.859	165	16565	13.515	ng/ul	0.00
31) 4-Chloroaniline-d4	11.376	131	14835	9.280	ng/ul	0.00
46) Dimethylphthalate-d6	14.407	166	60935	14.057	ng/ul	0.00
49) Acenaphthylene-d8	14.719	160	67838	13.769	ng/ul	0.00
54) 4-Nitrophenol-d4	15.171	143	8556m	11.448	ng/ul	0.00
60) Fluorene-d10	16.000	176	54309	13.544	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.099	200	10562	11.962	ng/ul	0.00
73) Anthracene-d10	17.850	188	90190	14.087	ng/ul	0.00
81) Pyrene-d10	20.107	212	114110	13.397	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.336	264	104997	14.464	ng/ul	-0.01

Target Compounds Qvalue

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

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