

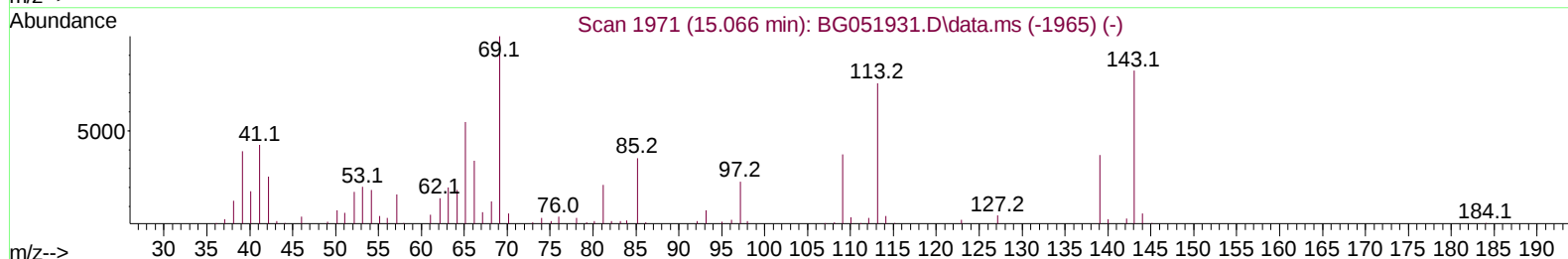
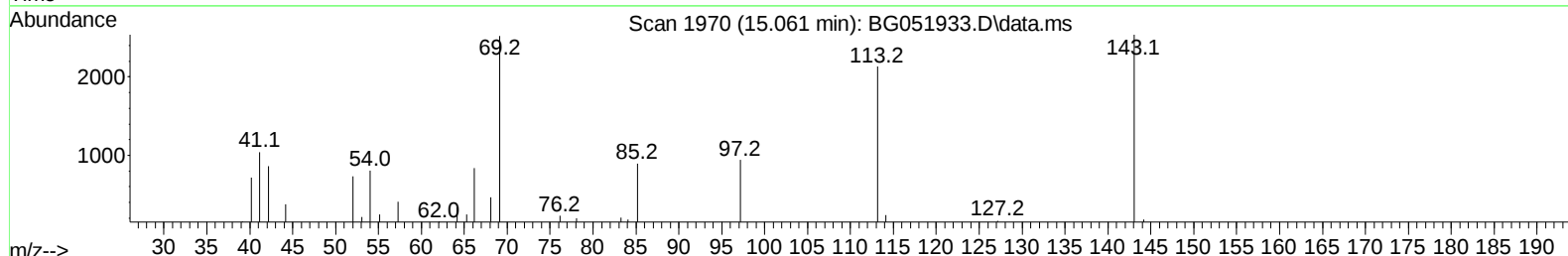
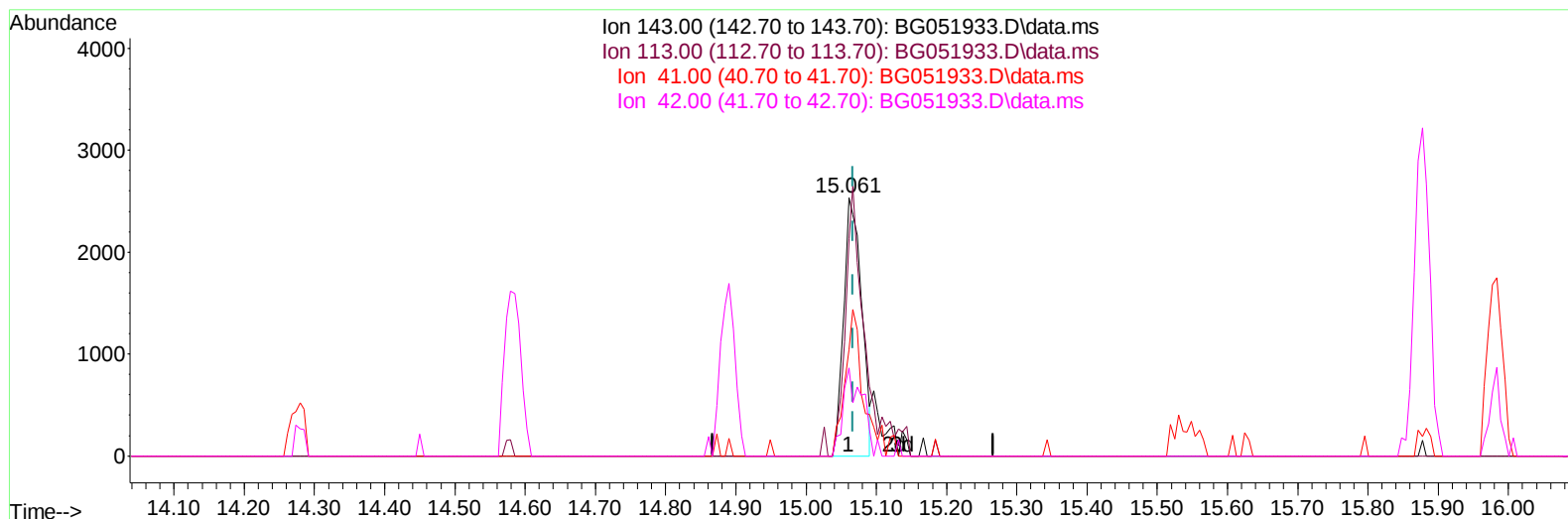
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG011022\
Data File : BG051933.D
Acq On : 10 Jan 2022 16:54
Operator : CG/JU
Sample : N1027-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0J37

Manual IntegrationsAPPROVED

Quant Time: Jan 11 00:05:26 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jan 06 14:43:02 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/11/2022
Supervised By :mohammad ahmed 01/12/2022



TIC: BG051933.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.061min (-0.006) 5.57 ng/u1

response 4494

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	84.03
41.00	44.40	41.17
42.00	29.70	33.99

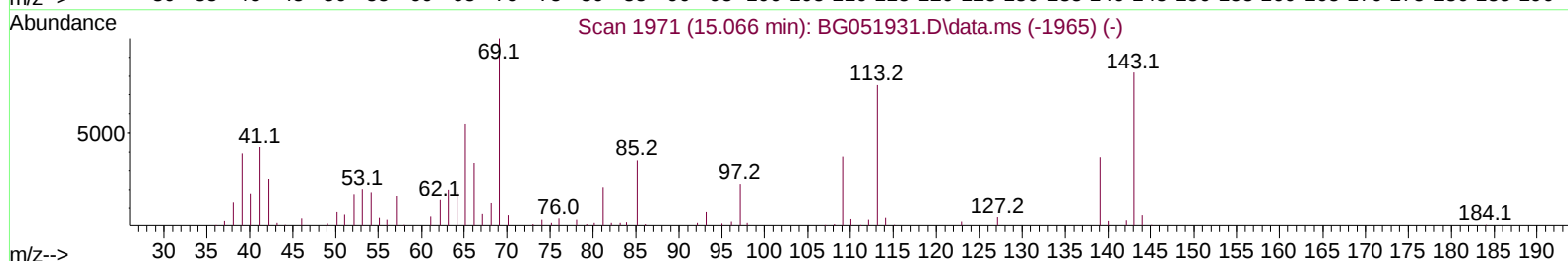
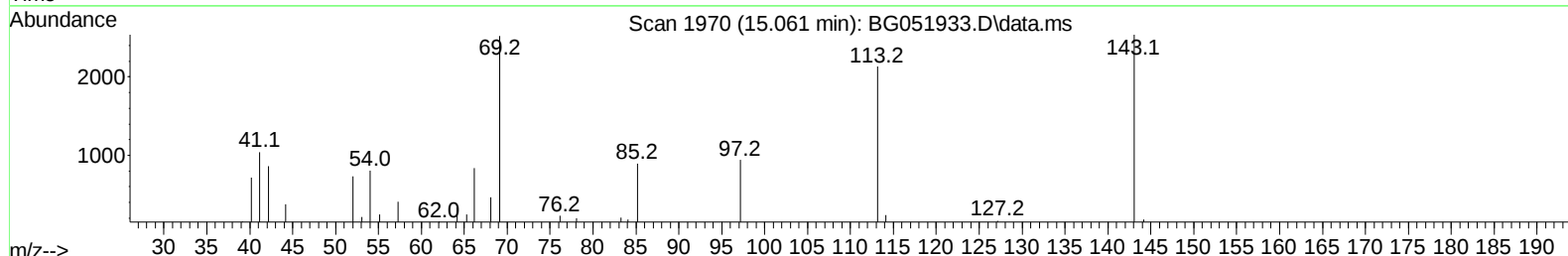
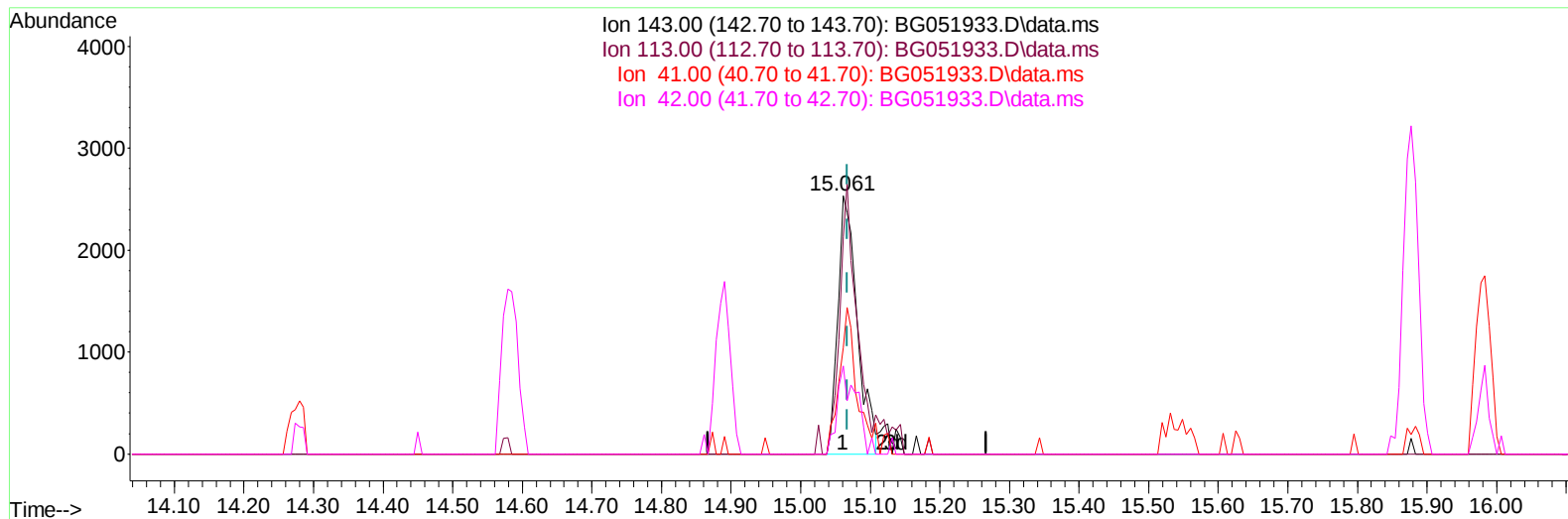
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TIC: BG051933.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.061min (-0.006) 6.12 ng/ul m

response 4936

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	84.03
41.00	44.40	41.17
42.00	29.70	33.99

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.246	152	19324	20.000	ng/ul	0.00
20) Naphthalene-d8	11.084	136	77820	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.890	164	58392	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.640	188	173370	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.957	240	192368	20.000	ng/ul	# 0.00
88) Perylene-d12	25.441	264	190197	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.570	96	2735	5.057	ng/uL	0.00
4) Pyridine-d5	4.011	84	5775	3.435	ng/ul	0.00
7) Phenol-d5	7.383	99	10949	5.789	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.547	67	38720	32.250	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.771	132	30915	24.689	ng/ul	0.00
15) 4-Methylphenol-d8	8.946	113	20629	14.031	ng/ul	0.00
21) Nitrobenzene-d5	9.415	128	21361	34.738	ng/ul	0.00
24) 2-Nitrophenol-d4	10.144	143	21961	32.292	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.696	165	39856	29.983	ng/ul	0.00
31) 4-Chloroaniline-d4	11.213	131	27075	14.625	ng/ul	0.00
46) Dimethylphthalate-d6	14.279	166	186966	38.737	ng/ul	0.00
49) Acenaphthylene-d8	14.585	160	206083	35.347	ng/ul	0.00
54) 4-Nitrophenol-d4	15.061	143	4936m	6.117	ng/ul	0.00
60) Fluorene-d10	15.877	176	161742	38.435	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.983	200	29301	26.791	ng/ul	0.00
73) Anthracene-d10	17.734	188	342136	41.601	ng/ul	0.00
81) Pyrene-d10	20.013	212	452109	40.837	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.200	264	410620	41.009	ng/ul	0.00

Target Compounds Qvalue

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

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