

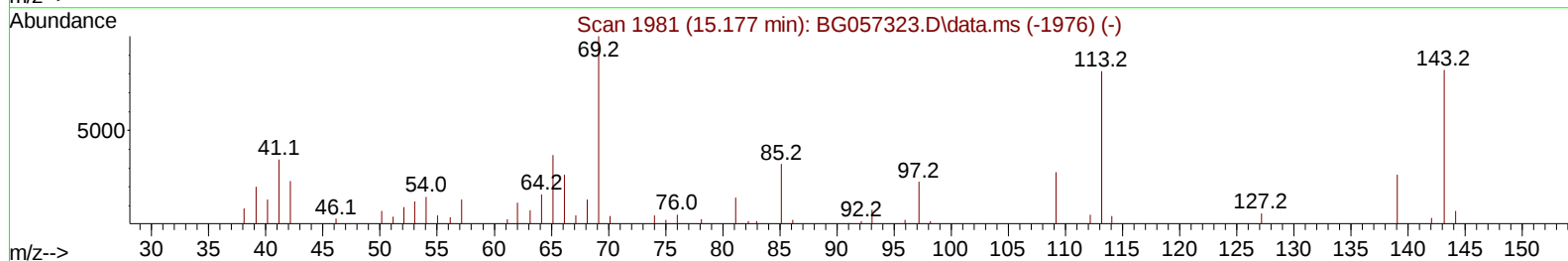
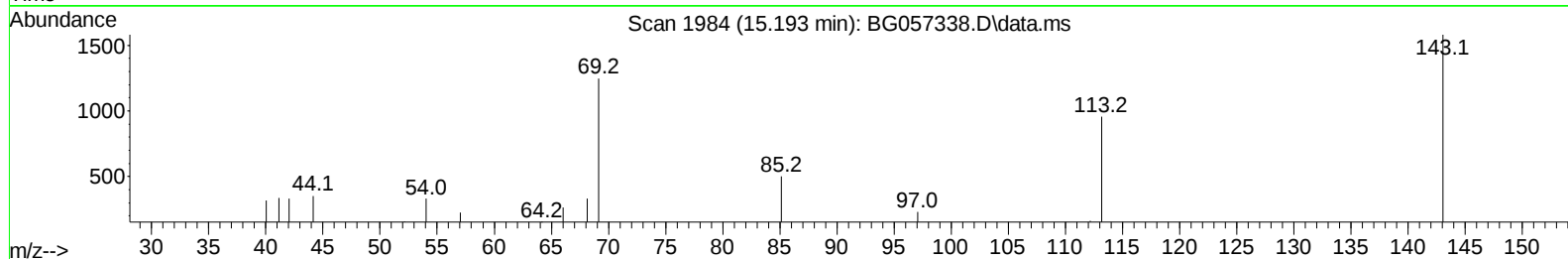
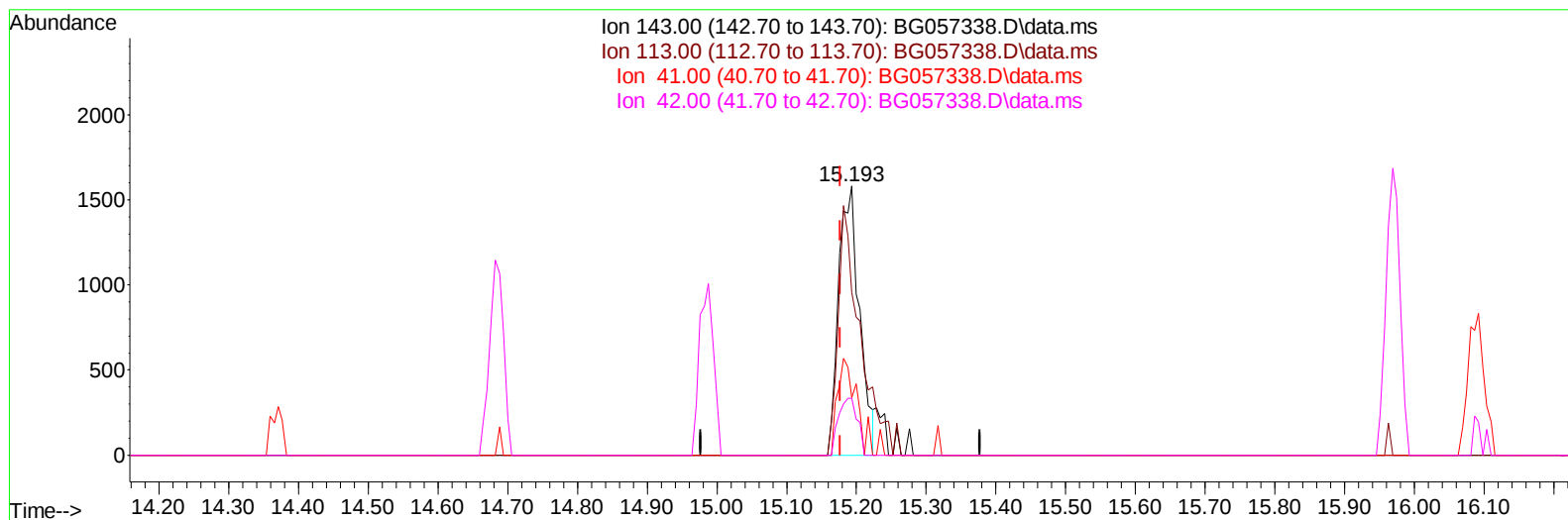
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG050823\
Data File : BG057338.D
Acq On : 8 May 2023 20:12
Operator : CG/JU
Sample : 02609-01
Misc :
ALS Vial : 16 Sample Multiplier: 3

Instrument :
BNA_G
ClientSampleId :
JREC2

Manual IntegrationsAPPROVED

Quant Time: May 09 00:11:59 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG042823.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon May 08 11:47:25 2023
Response via : Initial Calibration

Reviewed By :Christian Giraldo 05/09/2023
Supervised By :Jagrut Upadhyay 05/09/2023



TIC: BG057338.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.193min (+ 0.016) 4.91 ng/u1

response 3257

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	95.60	60.44#
41.00	44.40	21.46#
42.00	28.40	21.01#

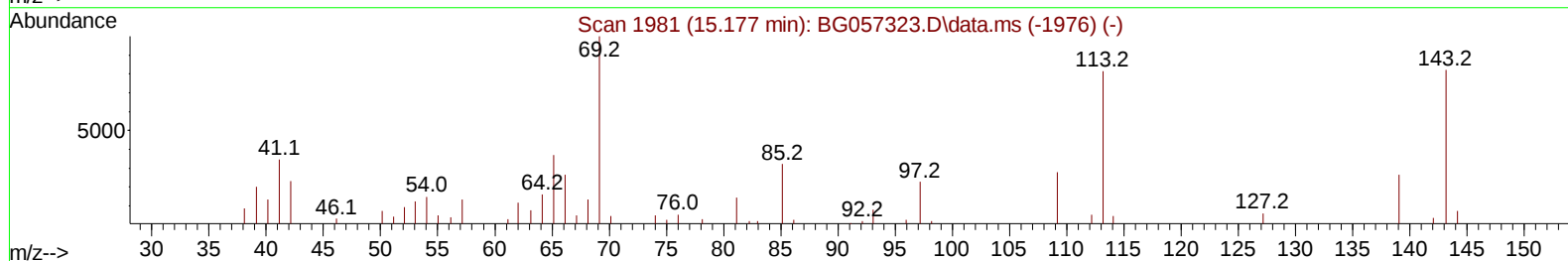
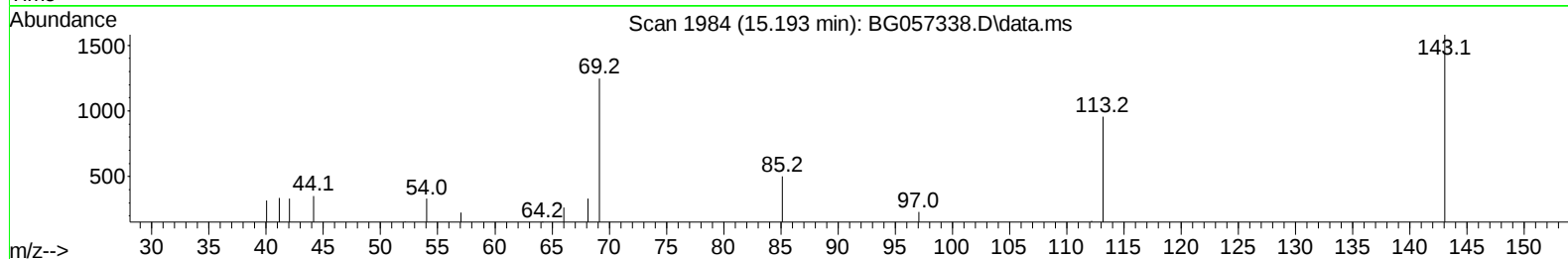
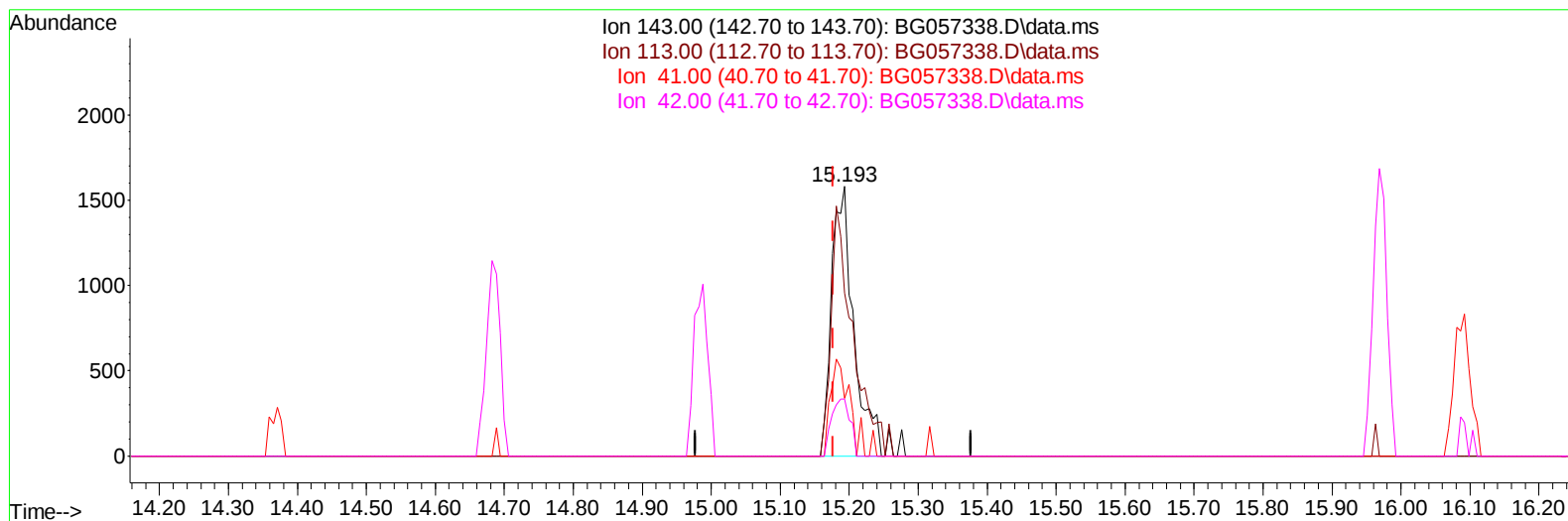
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Data File : BG057338.D
Acq On : 8 May 2023 20:12
Operator : CG/JU
Sample : 02609-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
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TIC: BG057338.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.193min (+ 0.016) 5.30 ng/ul m

response 3518

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	95.60	60.44#
41.00	44.40	21.46#
42.00	28.40	21.01#

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 Sample : 02609-01
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 ALS Vial : 16 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 05/09/2023
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Quant Time: May 09 00:58:13 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG042823.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon May 08 11:47:25 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.373	152	15444	20.000	ng/ul	0.00
20) Naphthalene-d8	11.211	136	69139	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.988	164	57520	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.725	188	144941	20.000	ng/ul	0.00
79) Chrysene-d12	22.037	240	131591	20.000	ng/ul	# 0.00
88) Perylene-d12	25.556	264	133893	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.650	96	1401	3.992	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.510	99	8337	5.708	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.674	67	20104	23.403	ng/ul	0.00
11) 2-Chlorophenol-d4	7.903	132	20778	21.588	ng/ul	0.00
15) 4-Methylphenol-d8	9.072	113	14042	12.650	ng/ul	0.00
21) Nitrobenzene-d5	9.548	128	14700	26.698	ng/ul	0.00
24) 2-Nitrophenol-d4	10.282	143	17178	26.733	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.829	165	30050	24.240	ng/ul	0.00
31) 4-Chloroaniline-d4	11.346	131	9564	5.821	ng/ul	0.00
46) Dimethylphthalate-d6	14.371	166	125753	27.591	ng/ul	0.00
49) Acenaphthylene-d8	14.682	160	136969	26.739	ng/ul	0.00
54) 4-Nitrophenol-d4	15.193	143	3518m	5.303	ng/ul	0.02
60) Fluorene-d10	15.969	176	115395	27.178	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.092	200	21174	24.657	ng/ul	0.00
73) Anthracene-d10	17.825	188	205202	31.017	ng/ul	0.00
81) Pyrene-d10	20.093	212	251467	32.195	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.315	264	226134	33.150	ng/ul	0.00

Target Compounds Qvalue

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

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