

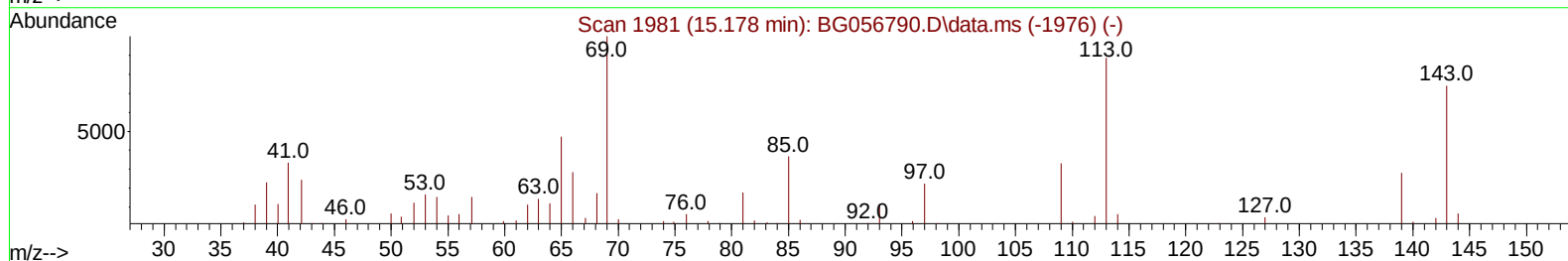
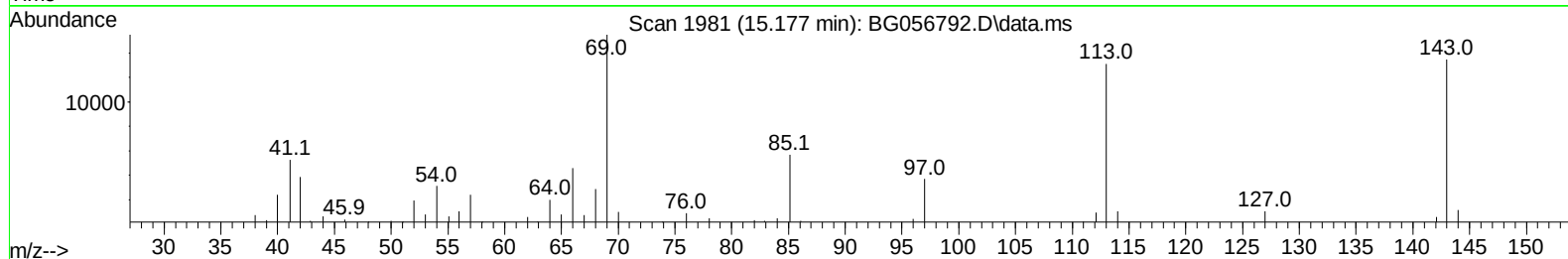
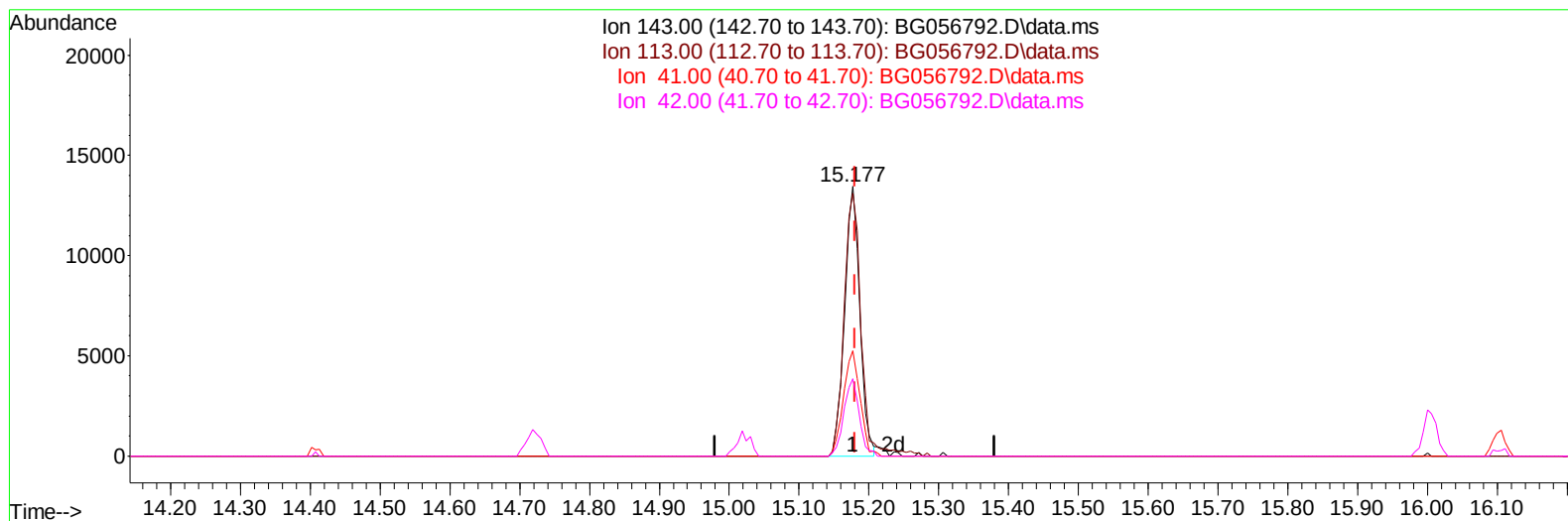
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030223\
Data File : BG056792.D
Acq On : 2 Mar 2023 16:03
Operator : CG/JU
Sample : PB151151BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK151

Manual IntegrationsAPPROVED

Quant Time: Mar 03 00:59:28 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 03/03/2023
Supervised By :Jagrut Upadhyay 03/03/2023



TIC: BG056792.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.177min (-0.003) 31.85 ng/ul

response 20187

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.30	97.47
41.00	46.10	39.13
42.00	27.10	28.66

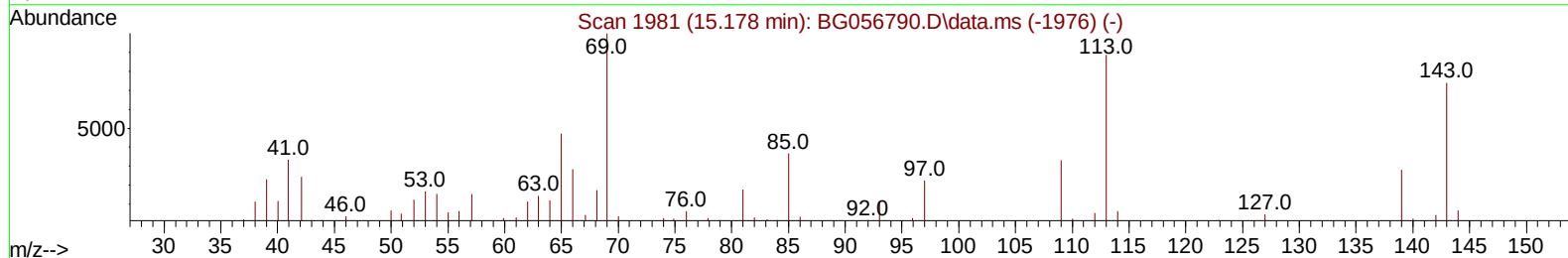
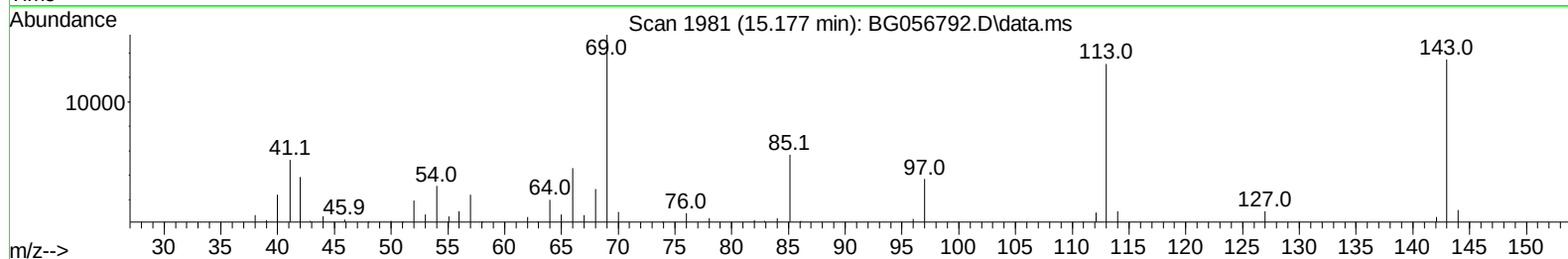
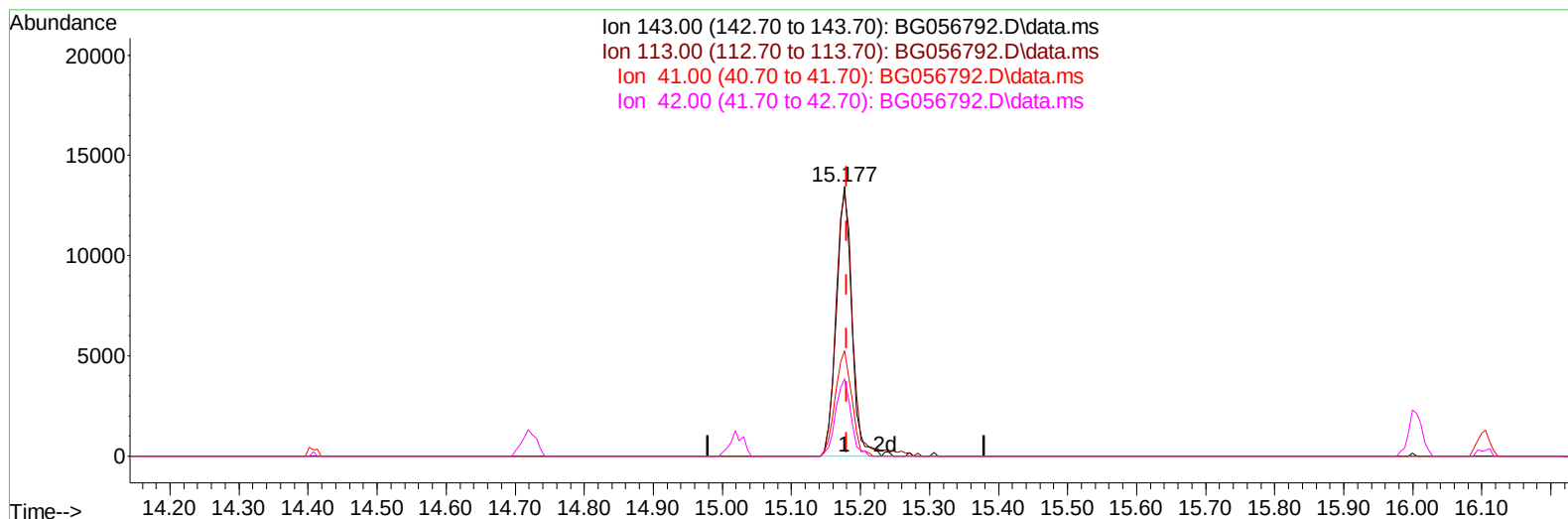
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TIC: BG056792.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.177min (-0.003) 32.50 ng/ul m

response 20594

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.30	97.47
41.00	46.10	39.13
42.00	27.10	28.66

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.420	152	12335	20.000	ng/ul	0.00
20) Naphthalene-d8	11.258	136	52581	20.000	ng/ul	0.00
38) Acenaphthene-d10	15.024	164	45557	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.756	188	120882	20.000	ng/ul	0.00
79) Chrysene-d12	22.063	240	119656	20.000	ng/ul	0.00
88) Perylene-d12	25.588	264	126738	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.702	96	2458	11.054	ng/uL	0.00
4) Pyridine-d5	4.137	84	34587	38.535	ng/ul	0.00
7) Phenol-d5	7.539	99	40457	32.776	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.721	67	26785	35.574	ng/ul	0.00
11) 2-Chlorophenol-d4	7.939	132	27260	33.256	ng/ul	0.00
15) 4-Methylphenol-d8	9.108	113	30585	32.428	ng/ul	0.00
21) Nitrobenzene-d5	9.595	128	15674	35.570	ng/ul	0.00
24) 2-Nitrophenol-d4	10.324	143	17885	34.097	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.859	165	34017	34.425	ng/ul	0.00
31) 4-Chloroaniline-d4	11.376	131	44883	34.824	ng/ul	0.00
46) Dimethylphthalate-d6	14.407	166	127345	34.646	ng/ul	0.00
49) Acenaphthylene-d8	14.719	160	143106	34.256	ng/ul	0.00
54) 4-Nitrophenol-d4	15.177	143	20594m	32.497	ng/ul	0.00
60) Fluorene-d10	16.006	176	117548	34.572	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.105	200	23909	29.563	ng/ul	0.00
73) Anthracene-d10	17.856	188	202402	34.514	ng/ul	0.00
81) Pyrene-d10	20.113	212	246823	30.484	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.348	264	236646	34.593	ng/ul	0.00

Target Compounds Qvalue

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

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