

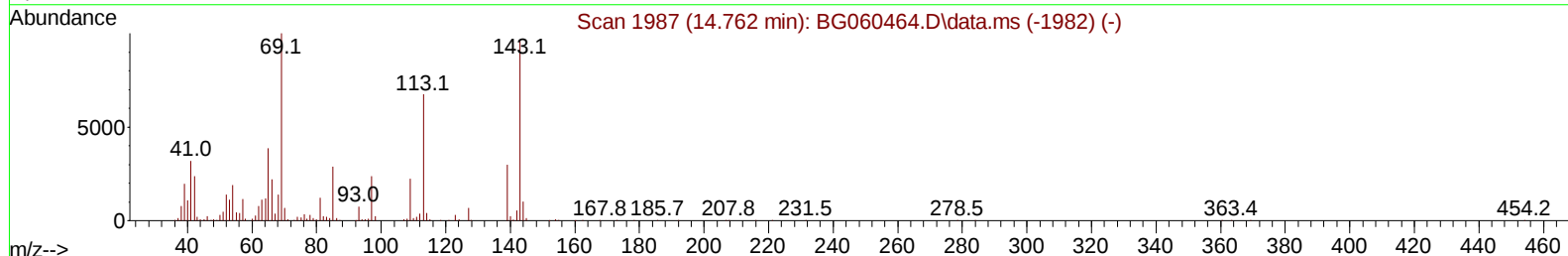
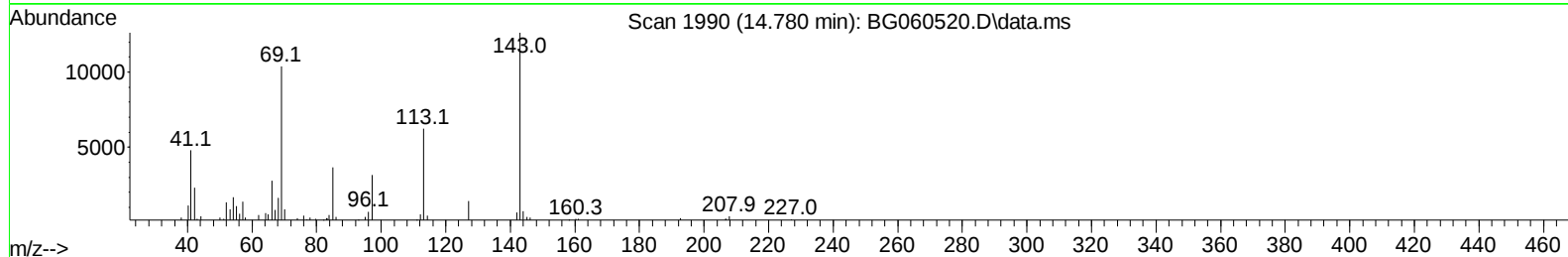
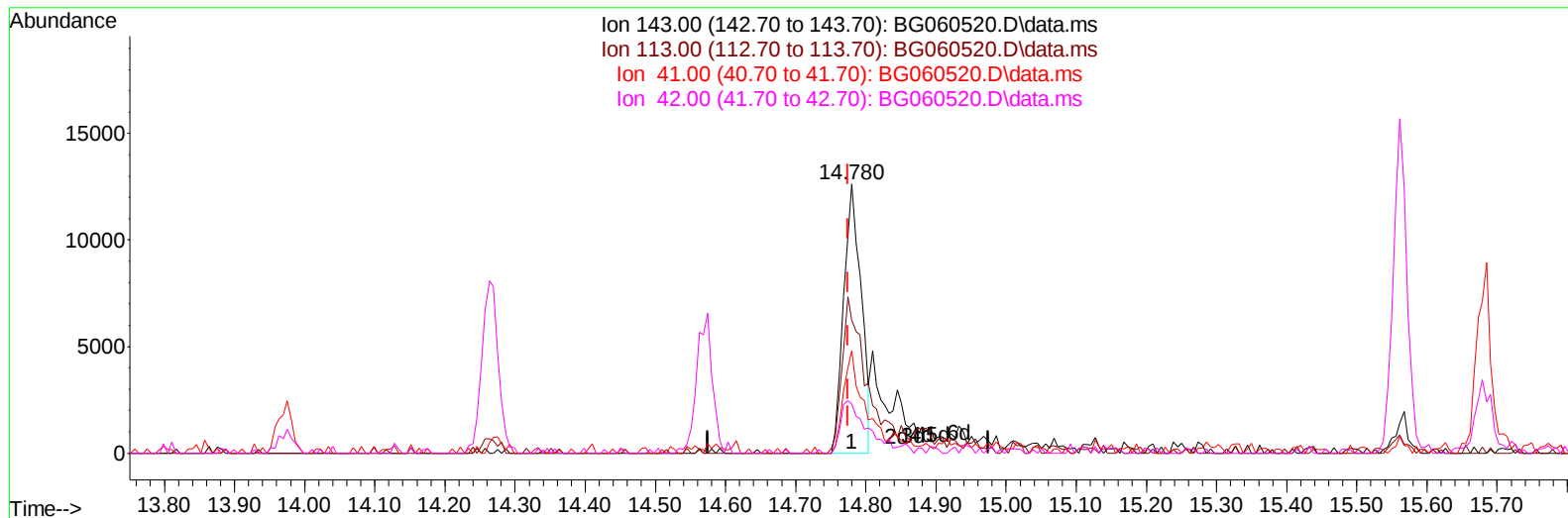
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG020124\
Data File : BG060520.D
Acq On : 31 Jan 2024 22:37
Operator : MA/JU
Sample : P1289-06
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AG3

Manual IntegrationsAPPROVED

Quant Time: Feb 01 01:18:34 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG012624.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jan 31 22:04:13 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 02/01/2024
Supervised By :Sohil Jodhani 02/02/2024



TIC: BG060520.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.780min (+ 0.005) 3.57 ng/u1

response 21986

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.60	49.33
41.00	29.80	38.07#
42.00	22.70	18.26

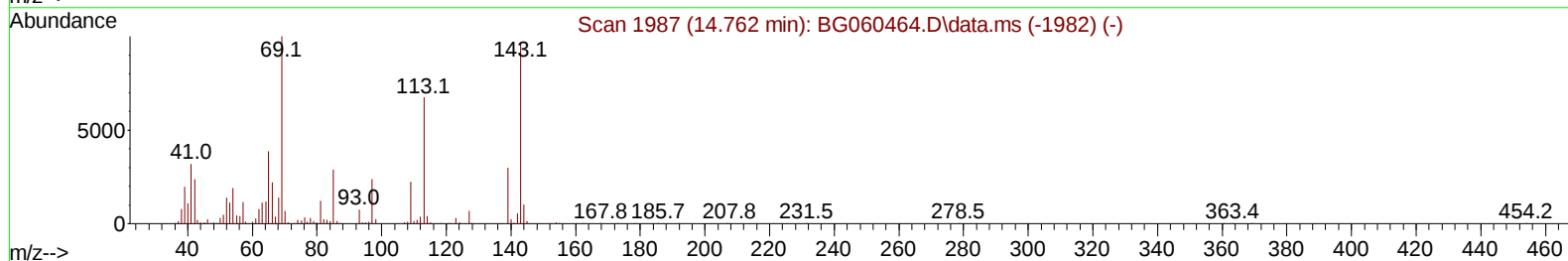
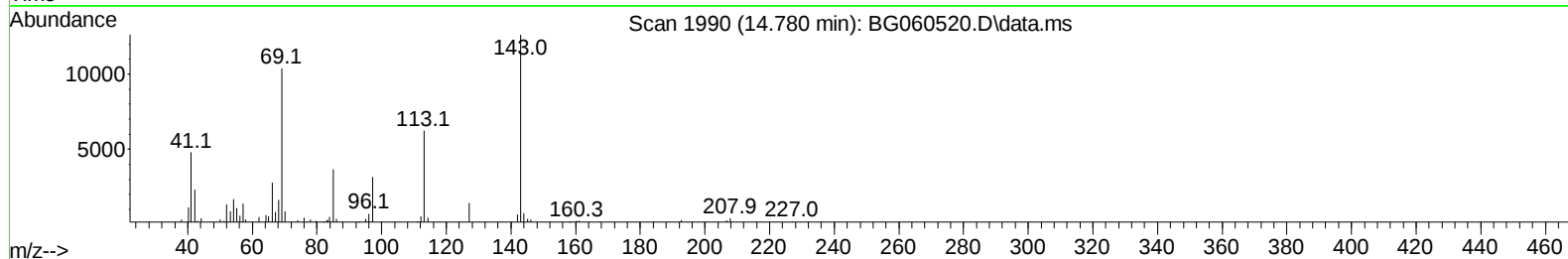
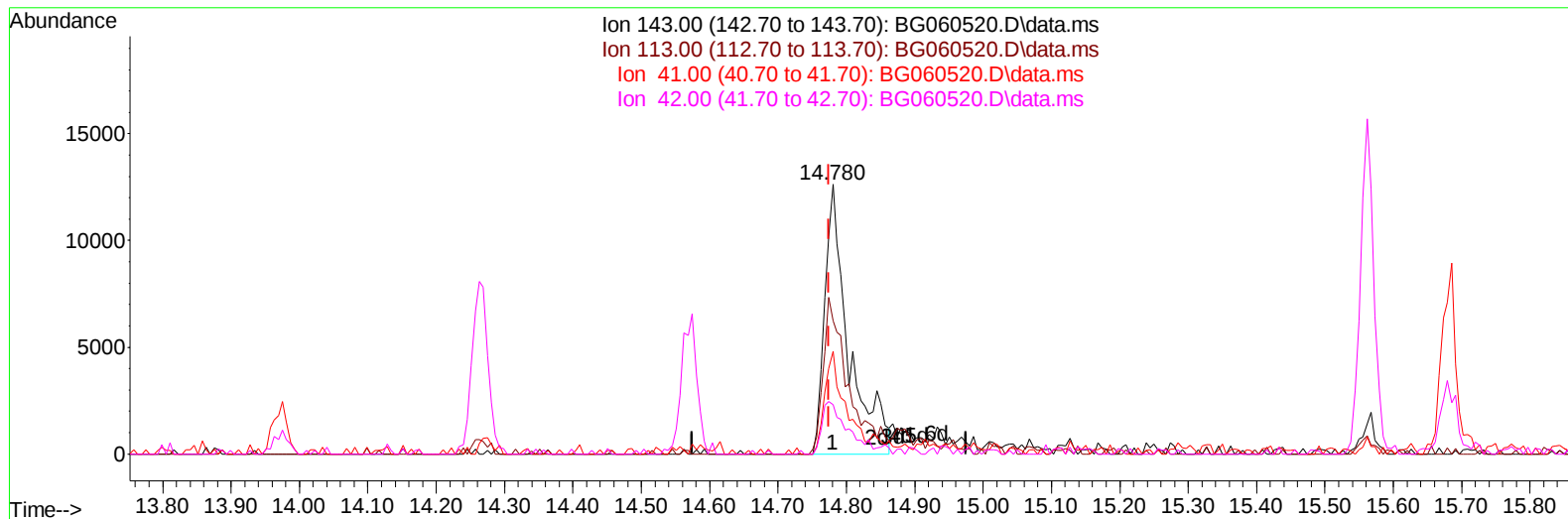
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TIC: BG060520.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.780min (+ 0.005) 4.96 ng/ul m

response 30591

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.60	49.33
41.00	29.80	38.07#
42.00	22.70	18.26

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 Sample : P1289-06
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Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 02/01/2024
 Supervised By :Sohil Jodhani 02/02/2024

Quant Time: Feb 01 01:22:05 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG012624.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jan 31 22:04:13 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.935	152	131858	20.000	ng/ul	0.00
20) Naphthalene-d8	10.738	136	609124	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.569	164	525987	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.318	188	1455218	20.000	ng/ul	0.00
79) Chrysene-d12	21.578	240	1424275	20.000	ng/ul	0.00
88) Perylene-d12	24.751	264	1552165	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	12862	4.888	ng/uL	0.00
4) Pyridine-d5	3.787	84	49844	5.988	ng/ul	0.00
7) Phenol-d5	7.101	99	78385	6.259	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.260	67	249786	31.741	ng/ul	0.00
11) 2-Chlorophenol-d4	7.465	132	208932	24.984	ng/ul	0.00
15) 4-Methylphenol-d8	8.640	113	164298	15.928	ng/ul	0.00
21) Nitrobenzene-d5	9.093	128	159221	35.434	ng/ul	0.00
24) 2-Nitrophenol-d4	9.815	143	190293	36.248	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.362	165	362635	33.215	ng/ul	0.00
31) 4-Chloroaniline-d4	10.873	131	371119	25.912	ng/ul	0.00
46) Dimethylphthalate-d6	13.975	166	1542475	35.838	ng/ul	0.00
49) Acenaphthylene-d8	14.263	160	1692755	35.062	ng/ul	0.00
54) 4-Nitrophenol-d4	14.780	143	30591m	4.962	ng/ul	0.00
60) Fluorene-d10	15.562	176	1418754	37.620	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.679	200	259043	28.748	ng/ul	0.00
73) Anthracene-d10	17.418	188	2569272	37.245	ng/ul	0.00
81) Pyrene-d10	19.710	212	3122987	36.585	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.539	264	3048883	37.445	ng/ul	0.00

Target Compounds Qvalue

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

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