

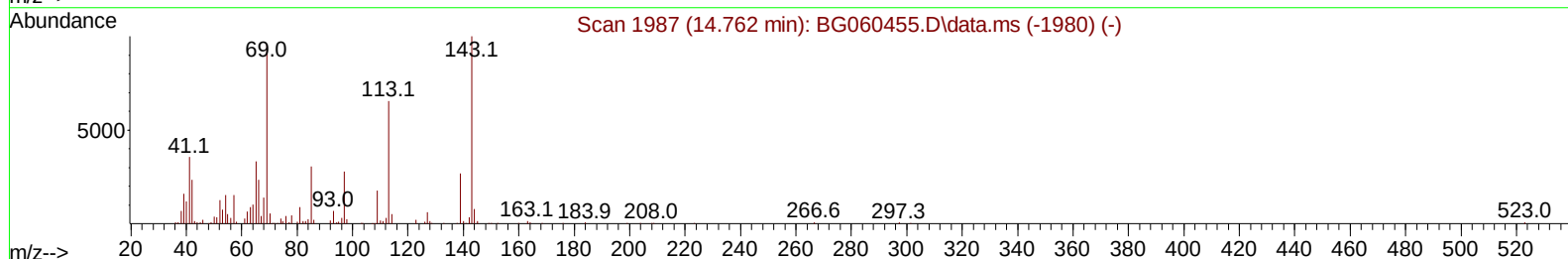
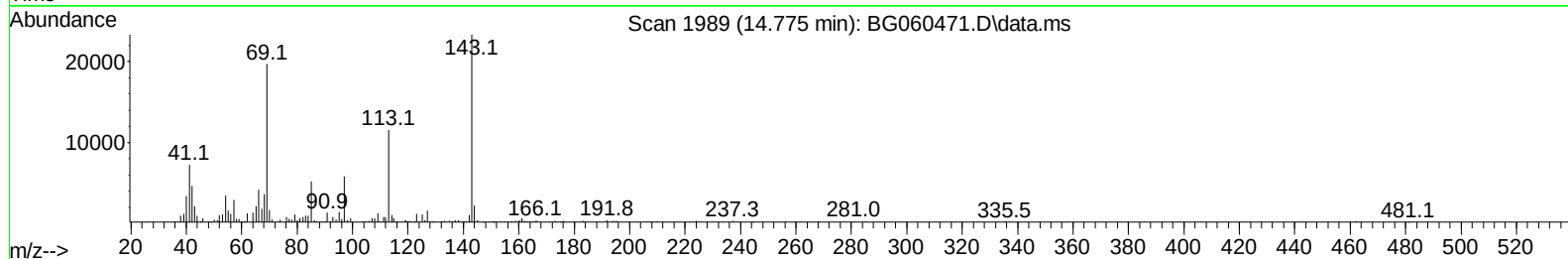
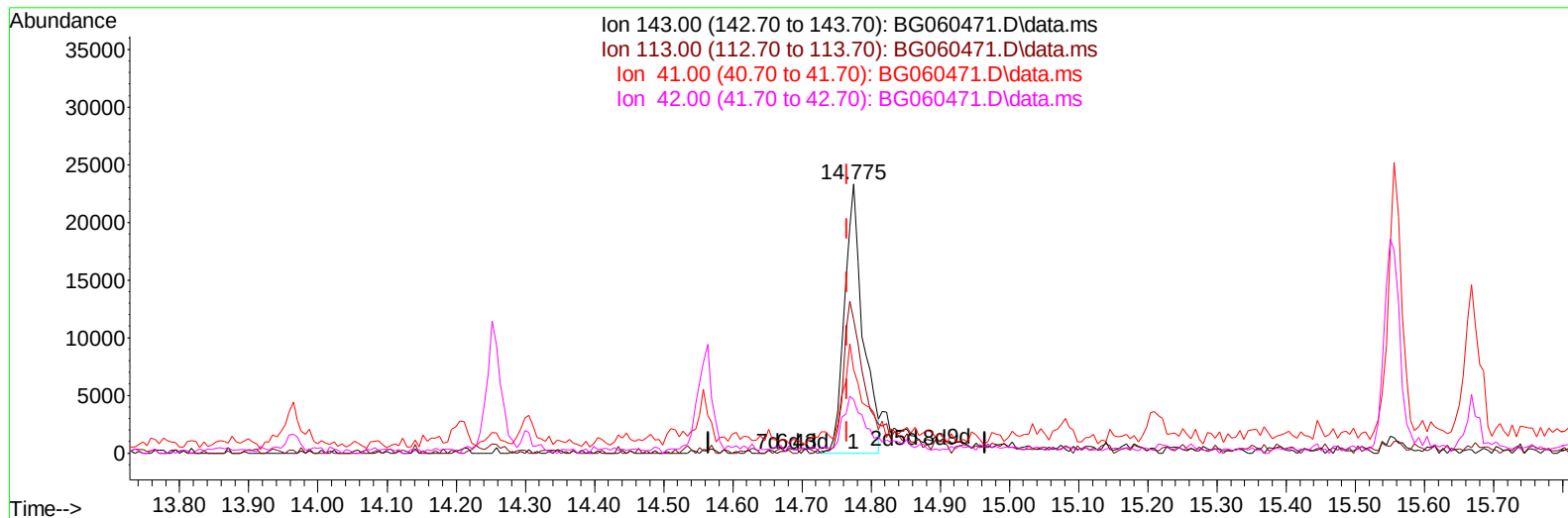
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012924\
Data File : BG060471.D
Acq On : 29 Jan 2024 21:01
Operator : MA/JU
Sample : P1239-16
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0FC6

Manual IntegrationsAPPROVED

Quant Time: Jan 30 01:49:14 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG012624.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jan 26 06:06:34 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/30/2024
Supervised By :mohammad ahmed 01/31/2024



TIC: BG060471.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.775min (+ 0.010) 5.30 ng/ul

response 43159

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.60	49.60
41.00	29.80	30.96
42.00	22.70	19.89

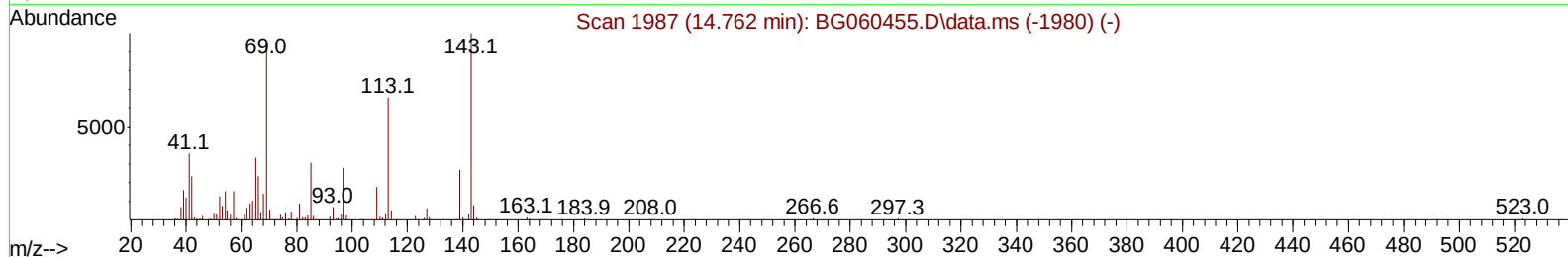
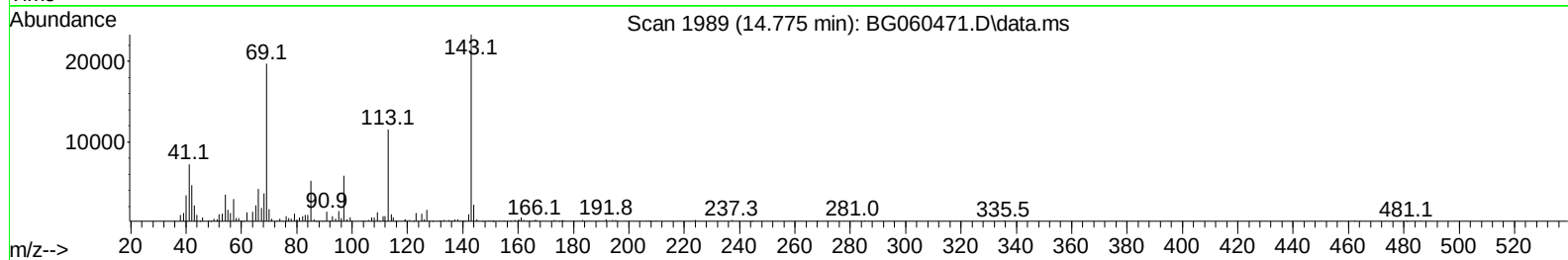
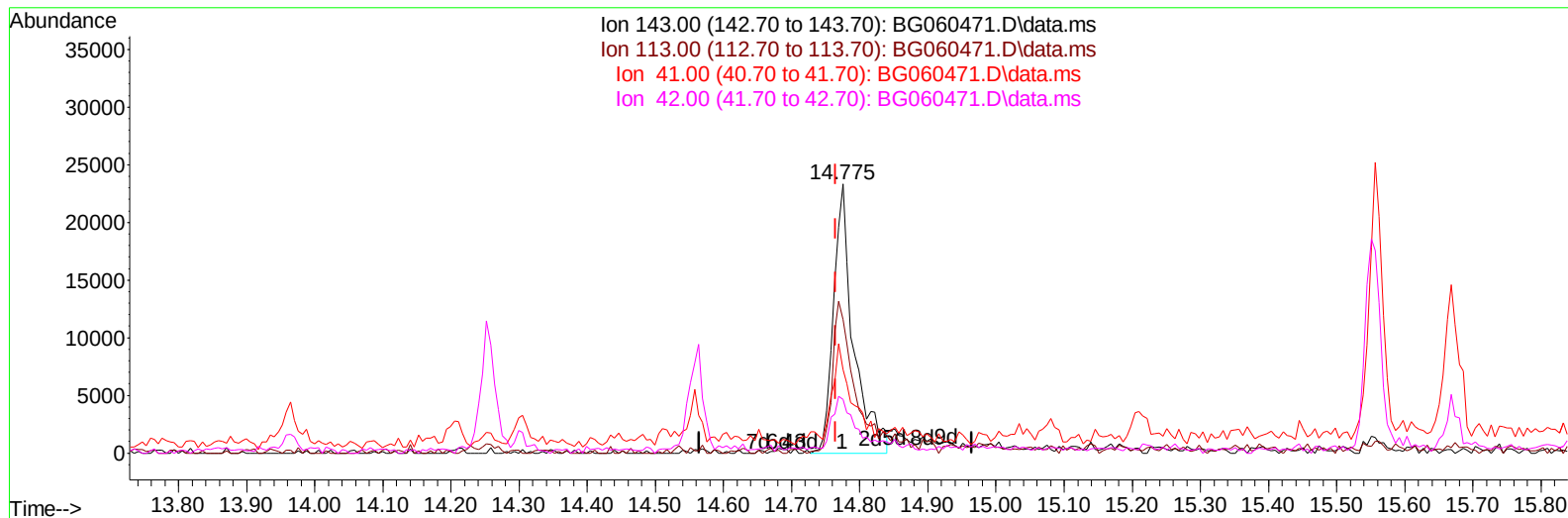
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TIC: BG060471.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.775min (+ 0.010) 5.86 ng/ul m

response 47734

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.60	49.60
41.00	29.80	30.96
42.00	22.70	19.89

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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 01/30/2024
 Supervised By :mohammad ahmed 01/31/2024

Quant Time: Jan 30 02:04:19 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG012624.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jan 26 06:06:34 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.924	152	182600	20.000	ng/ul	0.00
20) Naphthalene-d8	10.727	136	818373	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.563	164	694586	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.307	188	1609361	20.000	ng/ul	0.00
79) Chrysene-d12	21.573	240	1479120	20.000	ng/ul	0.00
88) Perylene-d12	24.734	264	1646421	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.365	96	19015	5.218	ng/uL	0.00
4) Pyridine-d5	3.770	84	122818	10.655	ng/ul	0.00
7) Phenol-d5	7.090	99	105383	6.077	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.249	67	300529	27.577	ng/ul	0.00
11) 2-Chlorophenol-d4	7.454	132	269333	23.257	ng/ul	0.00
15) 4-Methylphenol-d8	8.629	113	217535	15.228	ng/ul	0.00
21) Nitrobenzene-d5	9.088	128	186854	30.951	ng/ul	0.00
24) 2-Nitrophenol-d4	9.810	143	222329	31.522	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.345	165	459601	31.333	ng/ul	0.00
31) 4-Chloroaniline-d4	10.862	131	380799	19.789	ng/ul	0.00
46) Dimethylphthalate-d6	13.964	166	1807017	31.793	ng/ul	0.00
49) Acenaphthylene-d8	14.258	160	2003093	31.419	ng/ul	0.00
54) 4-Nitrophenol-d4	14.775	143	47734m	5.864	ng/ul	0.01
60) Fluorene-d10	15.556	176	1778543	35.713	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.674	200	335699	33.686	ng/ul	0.00
73) Anthracene-d10	17.407	188	2689943	35.260	ng/ul	0.00
81) Pyrene-d10	19.705	212	3153468	35.572	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.522	264	3150328	36.476	ng/ul	0.01
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.394	88	43201	10.715	ng/uL	98

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

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