

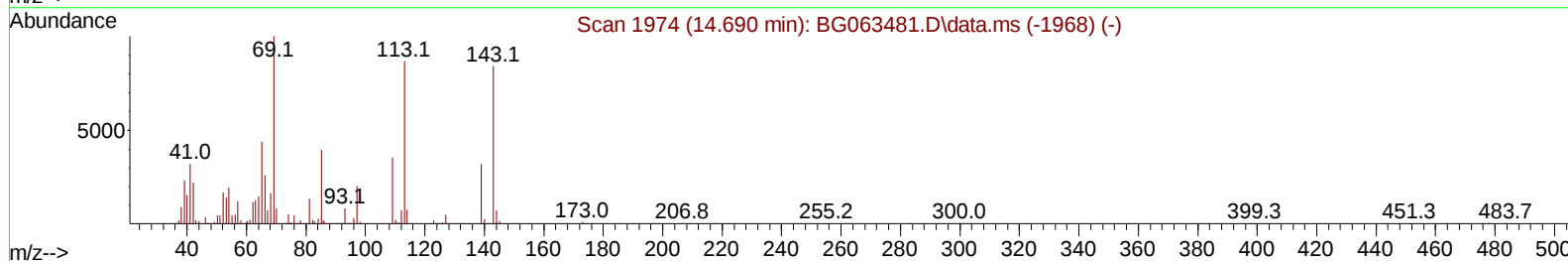
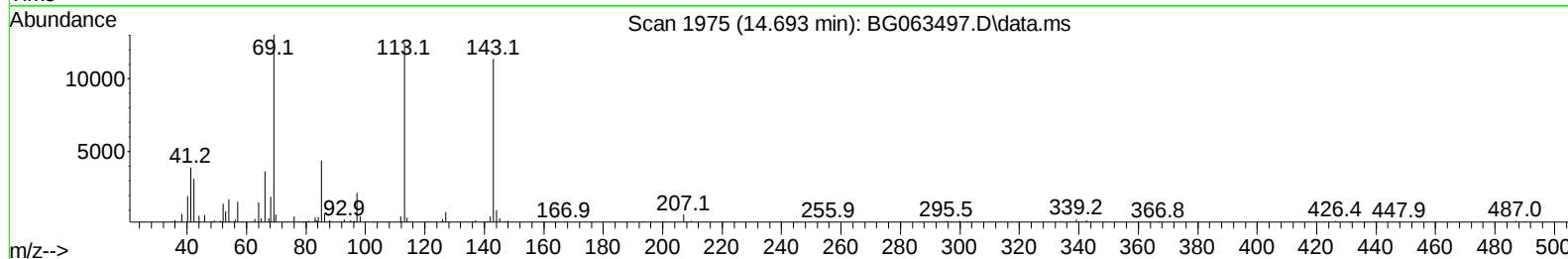
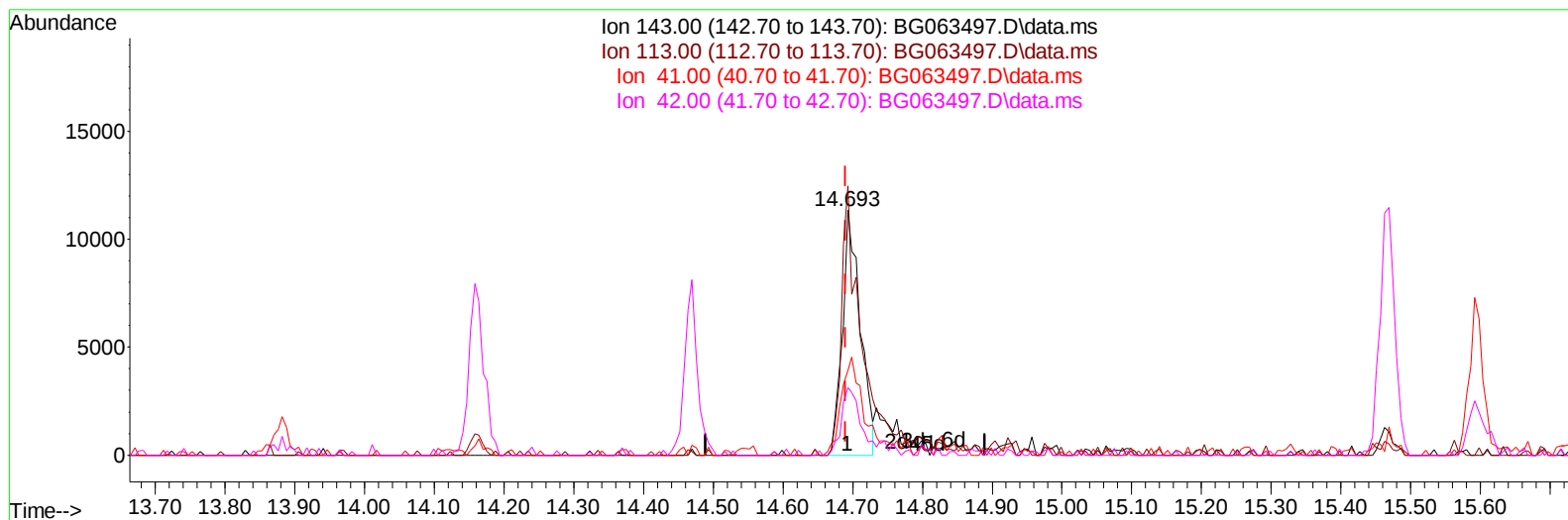
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112024\
Data File : BG063497.D
Acq On : 20 Nov 2024 23:32
Operator : RC/JU
Sample : P4904-11
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BG421

Manual IntegrationsAPPROVED

Quant Time: Nov 21 00:05:49 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112024.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 20 14:52:38 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/22/2024
Supervised By :mohammad ahmed 11/22/2024



TIC: BG063497.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.693min (+ 0.003) 6.60 ng/u1

response 20196

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	109.78
41.00	38.80	34.49
42.00	26.30	27.65

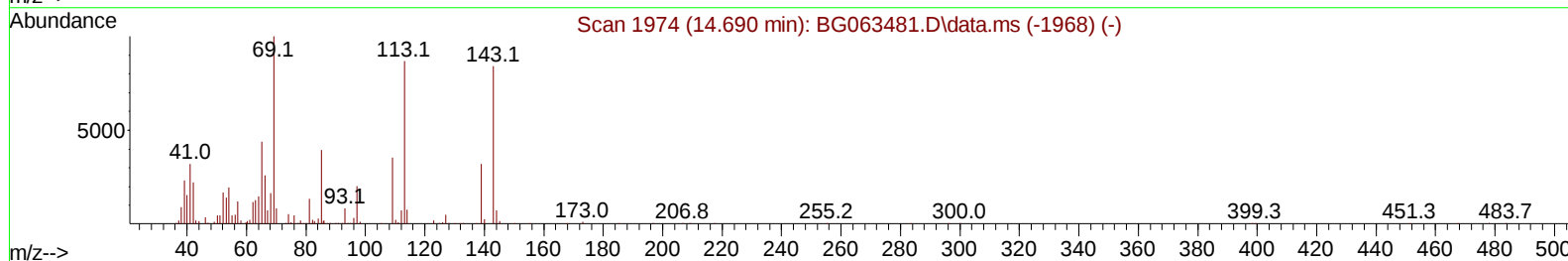
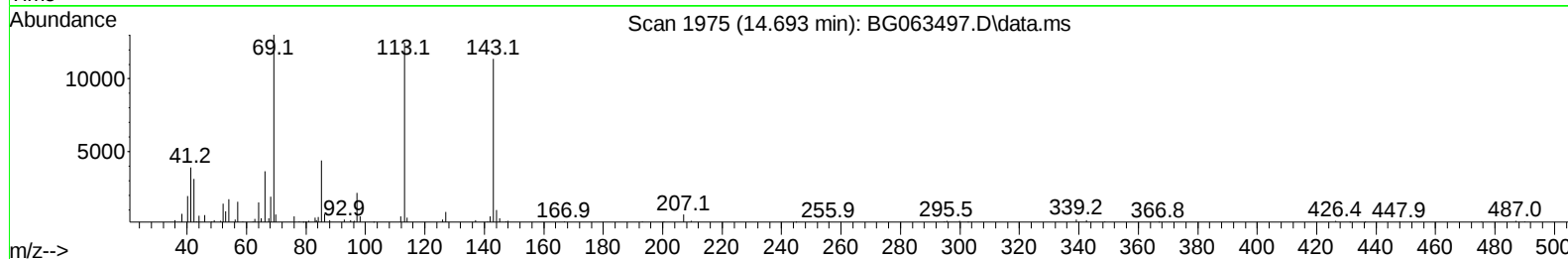
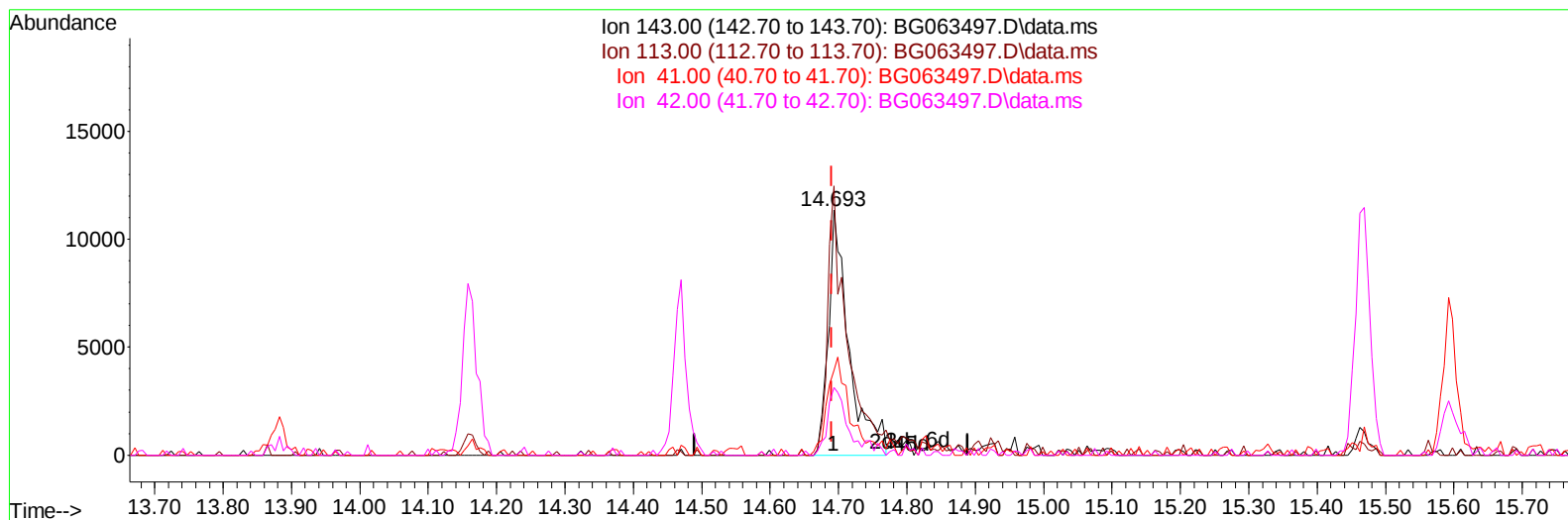
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TIC: BG063497.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.693min (+ 0.003) 7.76 ng/ul m

response 23729

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	109.78
41.00	38.80	34.49
42.00	26.30	27.65

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Quant Time: Nov 21 00:06:39 2024
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 QLast Update : Wed Nov 20 14:52:38 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.830	152	96507	20.000	ng/ul	0.00
20) Naphthalene-d8	10.621	136	377103	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.470	164	303337	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.225	188	730086	20.000	ng/ul	0.00
79) Chrysene-d12	21.479	240	787517	20.000	ng/ul	0.00
88) Perylene-d12	24.523	264	850953	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.300	96	13751	5.624	ng/uL	0.00
4) Pyridine-d5	3.712	84	46115	6.017	ng/ul	0.00
7) Phenol-d5	7.008	99	70844	8.926	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.161	67	181623	35.677	ng/ul	0.00
11) 2-Chlorophenol-d4	7.366	132	176660	31.406	ng/ul	0.00
15) 4-Methylphenol-d8	8.541	113	146209	22.044	ng/ul	0.00
21) Nitrobenzene-d5	8.982	128	99250	36.650	ng/ul	0.00
24) 2-Nitrophenol-d4	9.705	143	129643	38.752	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.251	165	224590	35.172	ng/ul	0.00
31) 4-Chloroaniline-d4	10.756	131	247109	30.596	ng/ul	0.00
46) Dimethylphthalate-d6	13.876	166	854714	41.507	ng/ul	0.00
49) Acenaphthylene-d8	14.164	160	903252	38.465	ng/ul	0.00
54) 4-Nitrophenol-d4	14.693	143	23729m	7.755	ng/ul	0.00
60) Fluorene-d10	15.468	176	767243	41.005	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	161591	39.369	ng/ul	0.00
73) Anthracene-d10	17.325	188	1440368	45.411	ng/ul	0.00
81) Pyrene-d10	19.622	212	1847330	46.797	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.317	264	1935254	46.888	ng/ul	0.00

Target Compounds Qvalue

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

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