

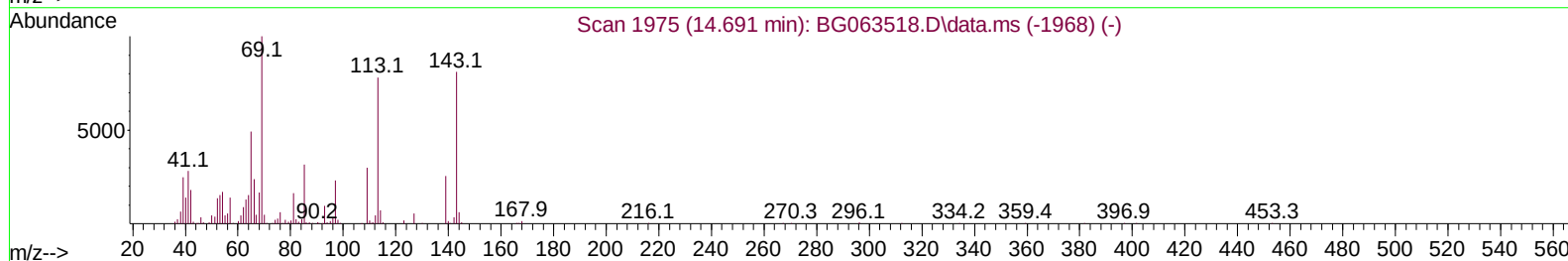
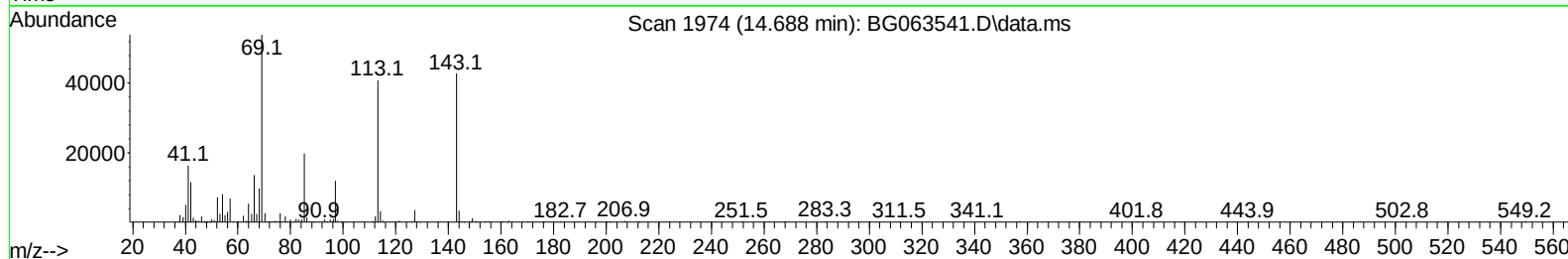
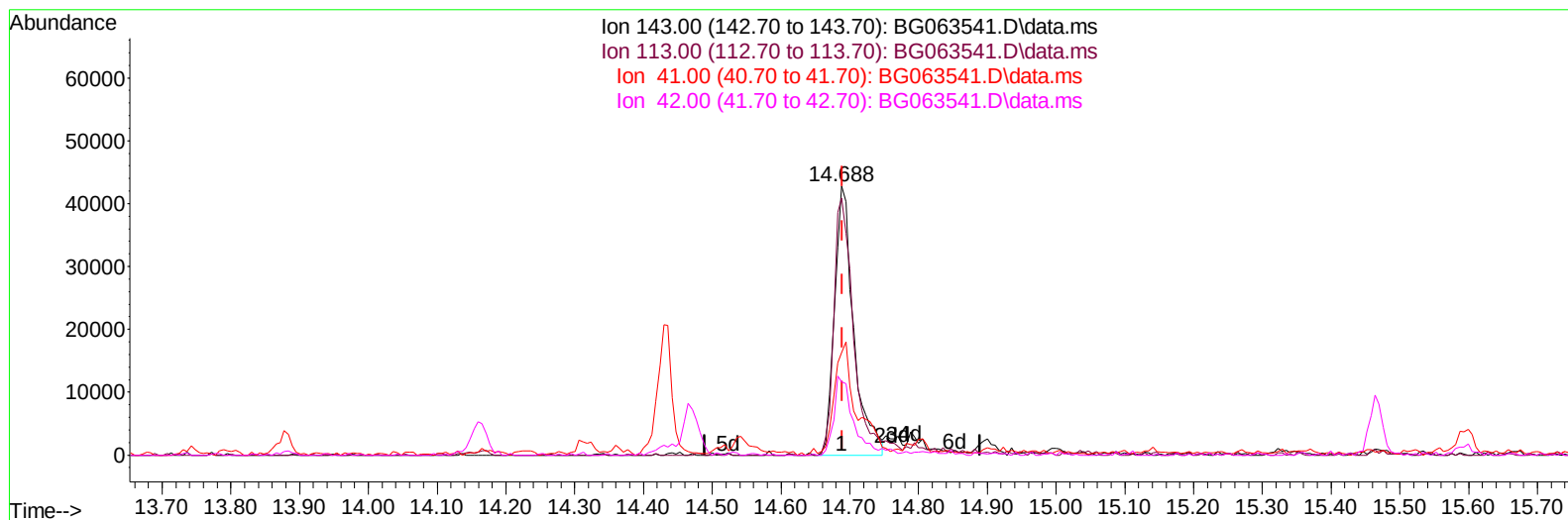
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112224\
Data File : BG063541.D
Acq On : 23 Nov 2024 5:08
Operator : RC/JU
Sample : P4881-07
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0TF7

Manual IntegrationsAPPROVED

Quant Time: Nov 23 05:42:11 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 :Jagrut Upadhyay 11/25/2024
Supervised By :mohammad ahmed 11/27/2024



TIC: BG063541.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.688min (-0.001) 20.80 ng/u1

response 83261

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	95.57
41.00	38.80	38.14
42.00	26.30	27.37

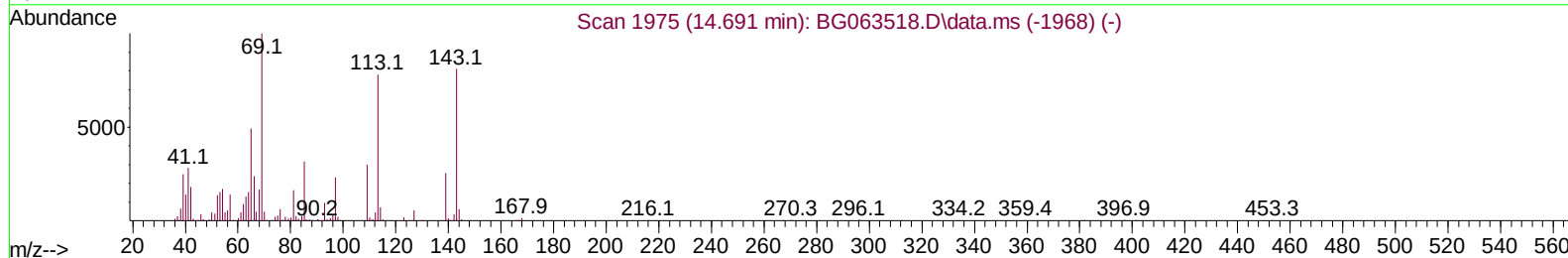
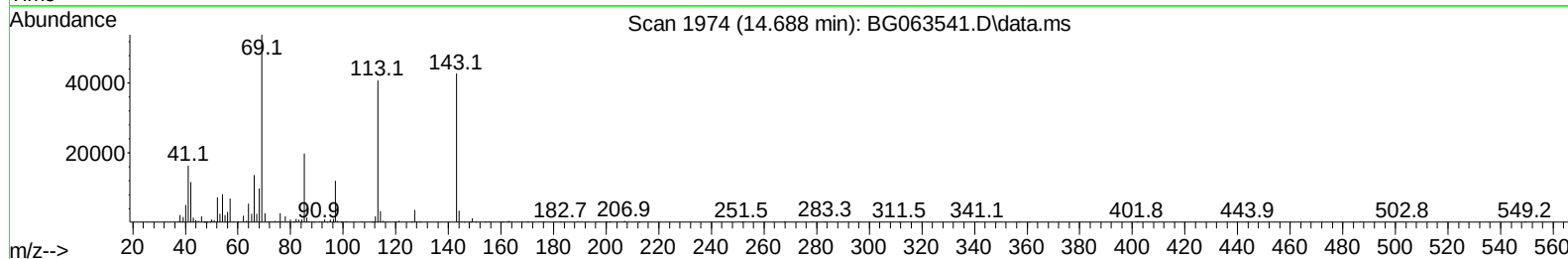
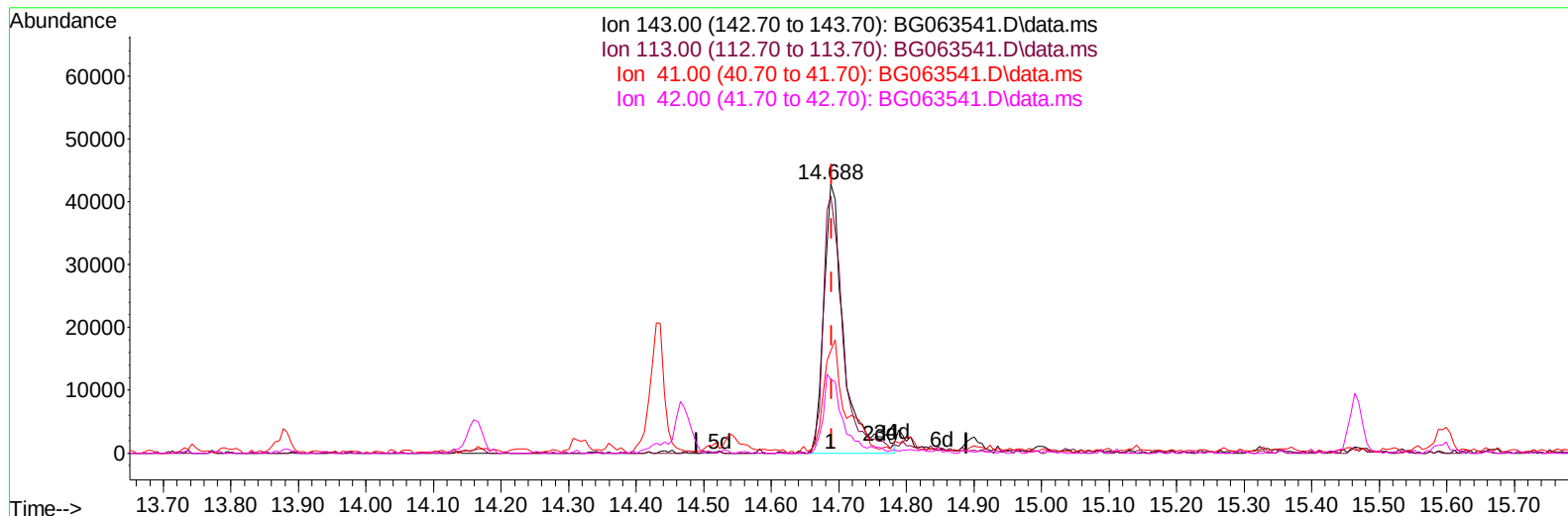
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(54) 4-Nitrophenol-d4 (S)

14.688min (-0.001) 22.07 ng/ul m

response 88361

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	95.57
41.00	38.80	38.14
42.00	26.30	27.37

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Quant Time: Nov 23 05:43:13 2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 20 14:52:38 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.832	152	121768	20.000	ng/ul	0.00
20) Naphthalene-d8	10.617	136	509118	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.471	164	396918	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.227	188	1007264	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	1080465	20.000	ng/ul	# 0.00
88) Perylene-d12	24.530	264	1164101	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.296	96	11361	3.682	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.003	99	208785	20.848	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.162	67	134136	20.883	ng/ul	0.00
11) 2-Chlorophenol-d4	7.362	132	150578	21.216	ng/ul	0.00
15) 4-Methylphenol-d8	8.537	113	179830	21.488	ng/ul	0.00
21) Nitrobenzene-d5	8.989	128	79145	21.647	ng/ul	0.00
24) 2-Nitrophenol-d4	9.706	143	97525	21.593	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.246	165	184097	21.355	ng/ul	0.00
31) 4-Chloroaniline-d4	10.758	131	162855	14.935	ng/ul	0.00
46) Dimethylphthalate-d6	13.877	166	620218	23.018	ng/ul	0.00
49) Acenaphthylene-d8	14.160	160	694202	22.593	ng/ul	0.00
54) 4-Nitrophenol-d4	14.688	143	88361m	22.070	ng/ul	0.00
60) Fluorene-d10	15.464	176	569685	23.268	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.593	200	108801	19.213	ng/ul	0.00
73) Anthracene-d10	17.321	188	956397	21.855	ng/ul	0.00
81) Pyrene-d10	19.624	212	1286118	23.747	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.318	264	1273334	22.552	ng/ul	0.00

Target Compounds Qvalue

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

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