

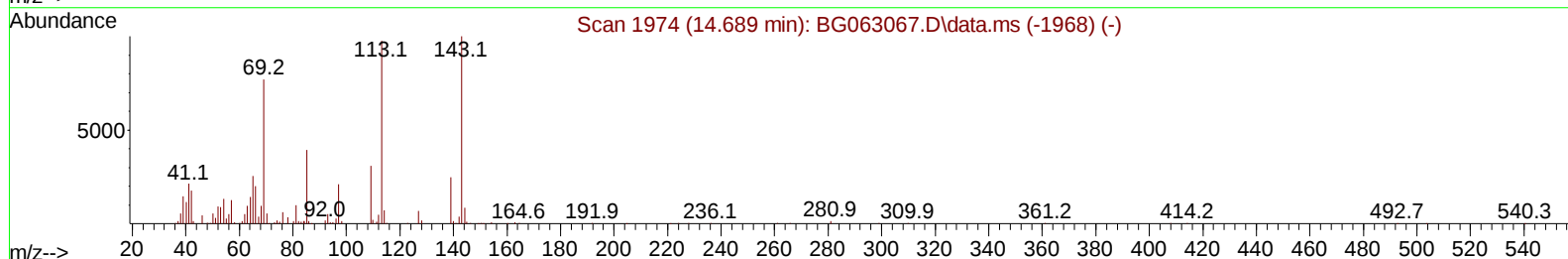
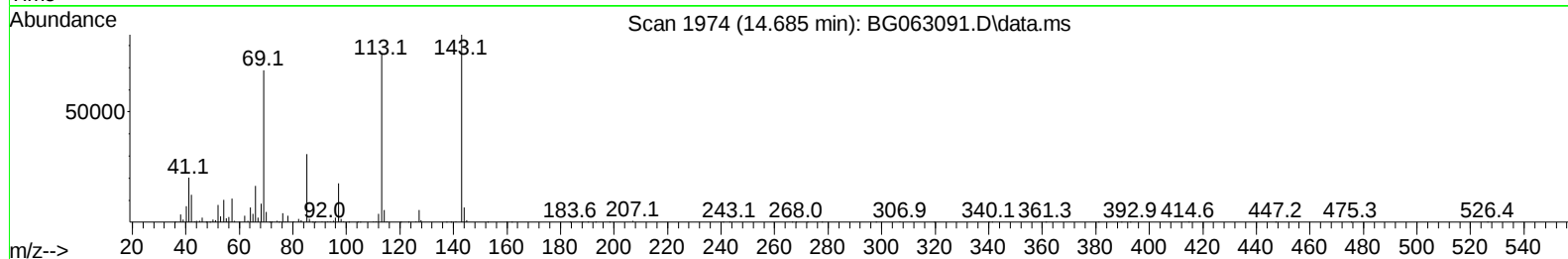
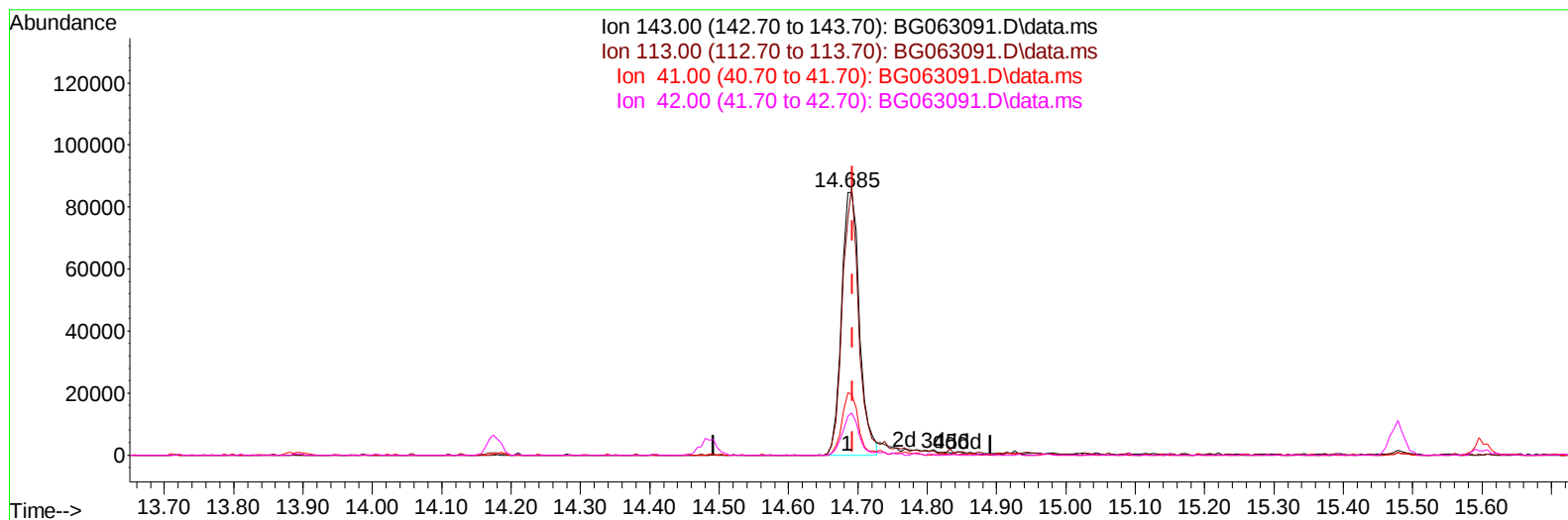
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092724\
Data File : BG063091.D
Acq On : 28 Sep 2024 3:03
Operator : RC/JU
Sample : PB163431BL
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK431

Manual IntegrationsAPPROVED

Quant Time: Sep 28 03:49:35 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092524.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Sep 25 13:16:45 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 09/30/2024
Supervised By :mohammad ahmed 10/01/2024



TIC: BG063091.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.685min (-0.007) 27.80 ng/u1

response 151544

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	120.10	89.98#
41.00	30.30	23.75#
42.00	17.10	14.91

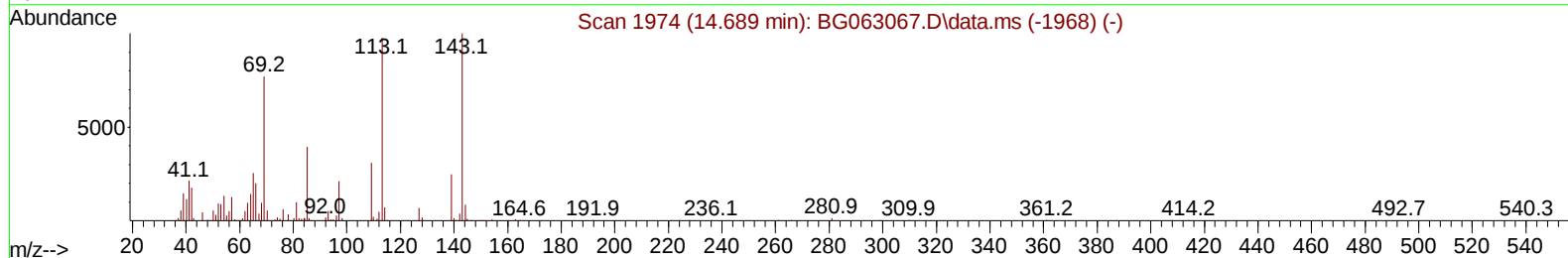
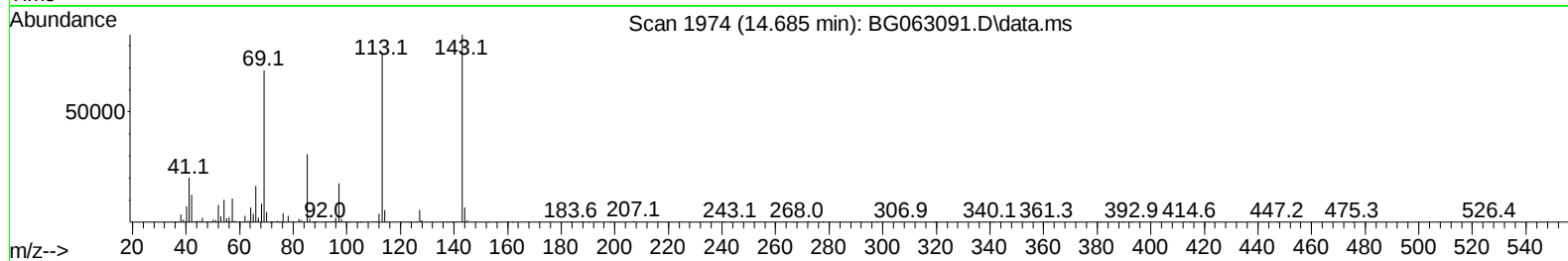
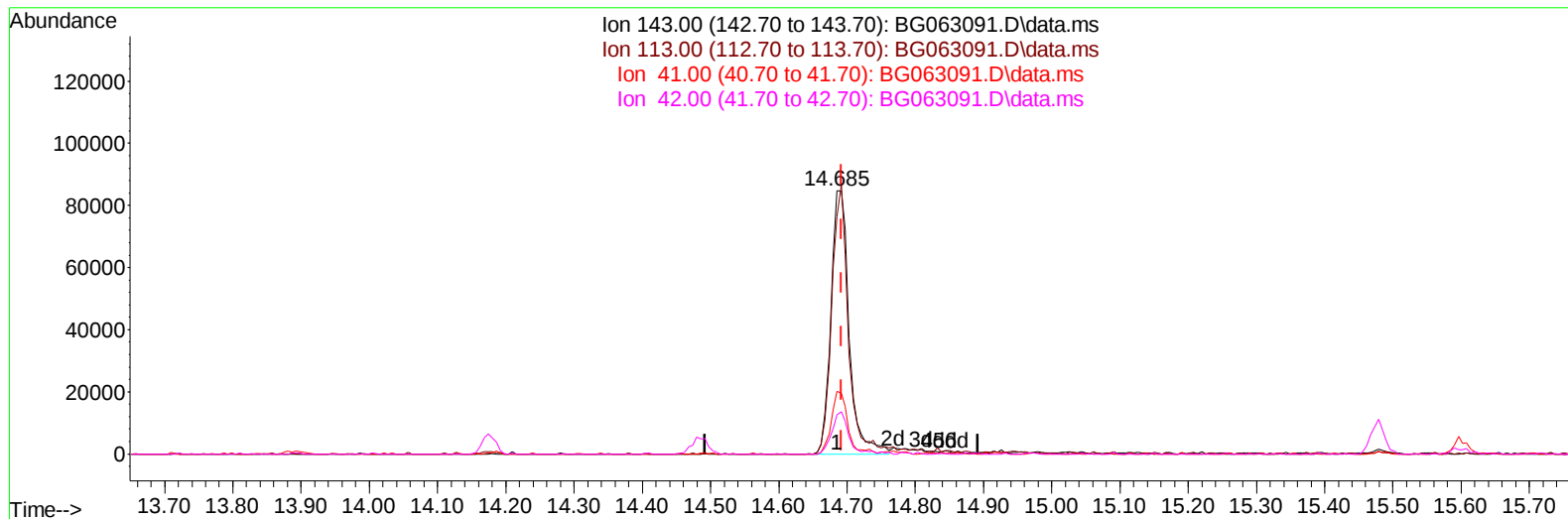
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TIC: BG063091.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.685min (-0.007) 28.84 ng/ul m

response 157225

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	120.10	89.98#
41.00	30.30	23.75#
42.00	17.10	14.91

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Quant Time: Sep 28 03:50:06 2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 25 13:16:45 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.846	152	109508	20.000	ng/ul	0.00
20) Naphthalene-d8	10.637	136	488351	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.486	164	454629	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.235	188	1216548	20.000	ng/ul	0.00
79) Chrysene-d12	21.483	240	1302661	20.000	ng/ul	# 0.00
88) Perylene-d12	24.533	264	1480376	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.311	96	11202	5.897	ng/uL	0.00
4) Pyridine-d5	3.722	84	159542	34.348	ng/ul	0.00
7) Phenol-d5	7.018	99	229267	33.896	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.177	67	95753	36.237	ng/ul	0.00
11) 2-Chlorophenol-d4	7.382	132	215399	32.411	ng/ul	0.00
15) 4-Methylphenol-d8	8.551	113	217244	35.579	ng/ul	0.00
21) Nitrobenzene-d5	9.004	128	116268	32.378	ng/ul	0.00
24) 2-Nitrophenol-d4	9.721	143	149069	32.504	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.261	165	295594	30.881	ng/ul	0.00
31) 4-Chloroaniline-d4	10.772	131	349632	31.129	ng/ul	0.00
46) Dimethylphthalate-d6	13.892	166	1093279	30.245	ng/ul	0.00
49) Acenaphthylene-d8	14.174	160	1190189	31.071	ng/ul	0.00
54) 4-Nitrophenol-d4	14.685	143	157225m	28.840	ng/ul	0.00
60) Fluorene-d10	15.479	176	1027885	31.035	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.602	200	212581	25.004	ng/ul	0.00
73) Anthracene-d10	17.335	188	1848219	31.666	ng/ul	0.00
81) Pyrene-d10	19.633	212	2424656	33.056	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.321	264	2499230	31.521	ng/ul	0.00

Target Compounds Qvalue

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

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