

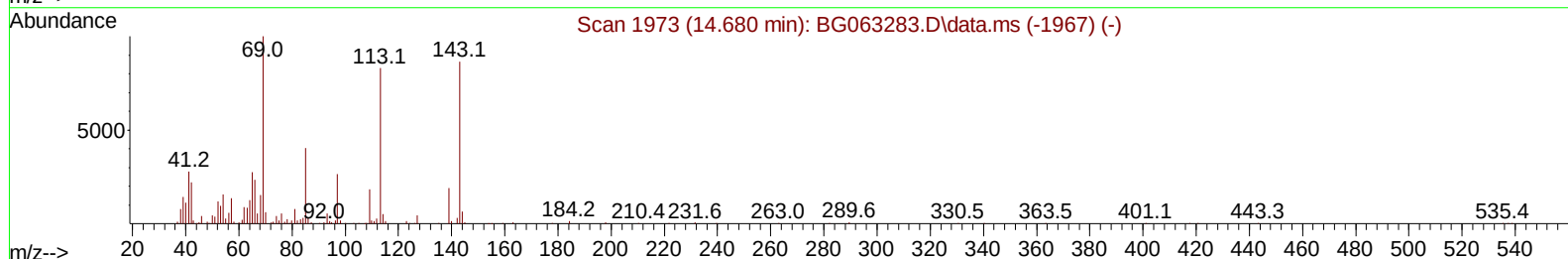
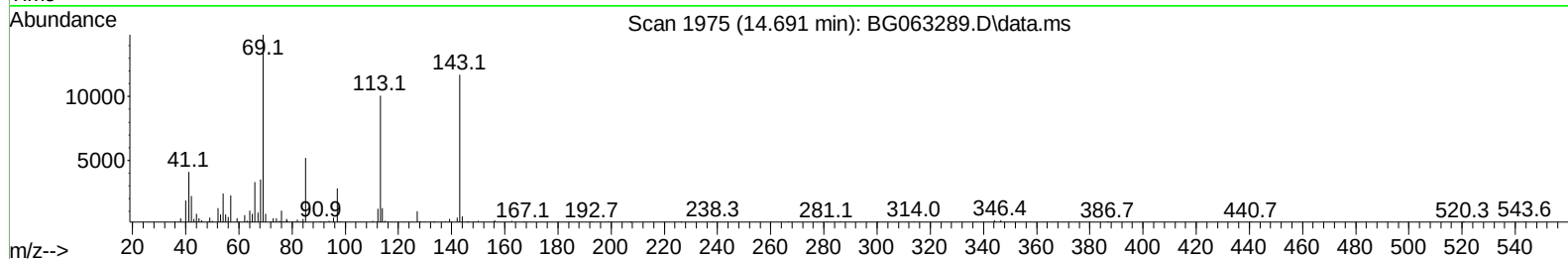
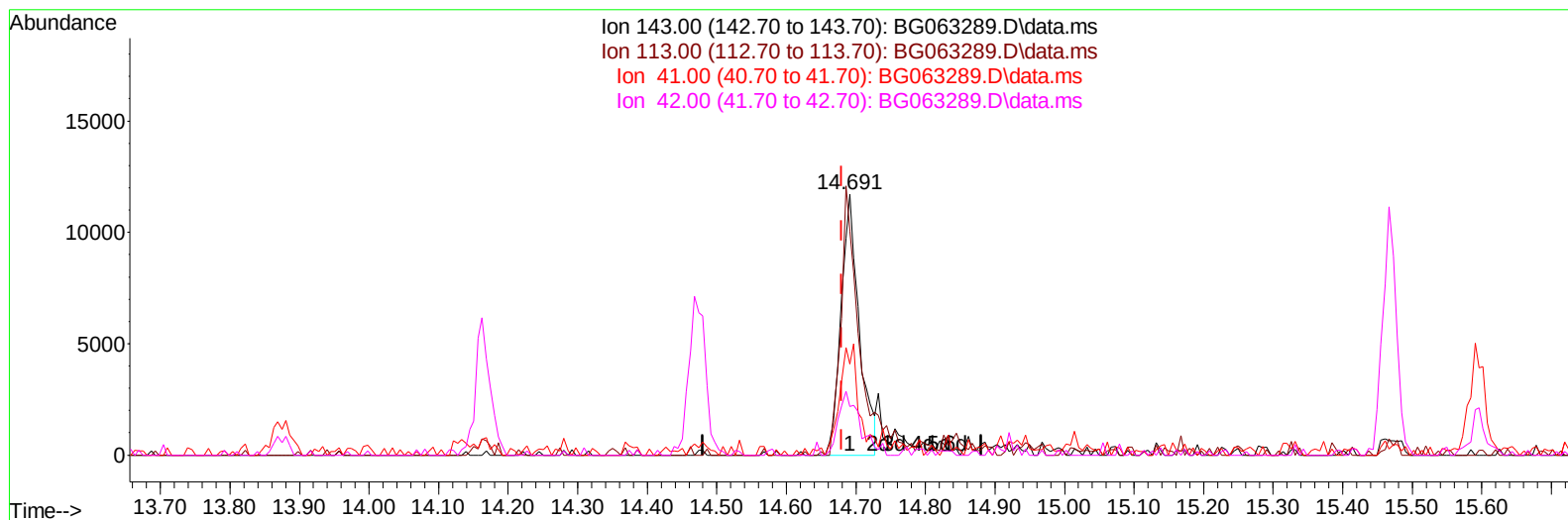
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111124\
Data File : BG063289.D
Acq On : 11 Nov 2024 14:56
Operator : RC/JU
Sample : P4625-16
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AC6

Manual IntegrationsAPPROVED

Quant Time: Nov 11 15:17:21 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110624.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 06 15:45:52 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/12/2024
Supervised By :mohammad ahmed 11/14/2024



TIC: BG063289.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.691min (+ 0.011) 5.08 ng/ul

response 21448

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	123.00	85.97#
41.00	42.60	34.92
42.00	30.90	18.58#

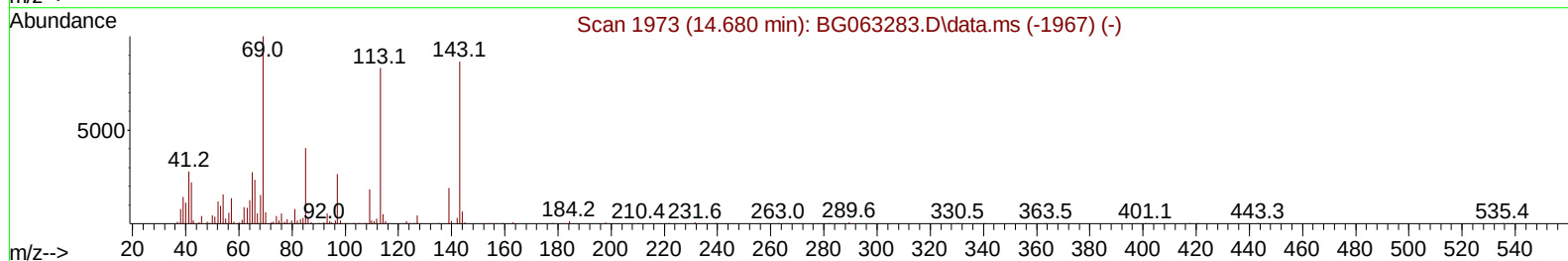
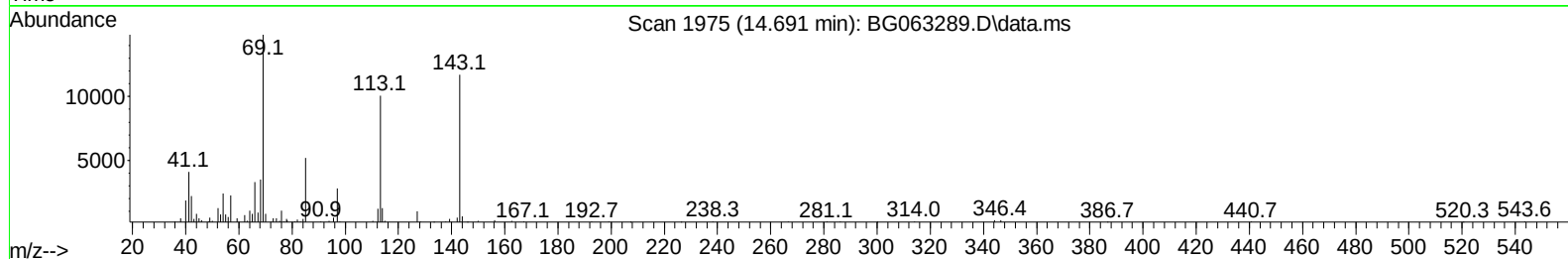
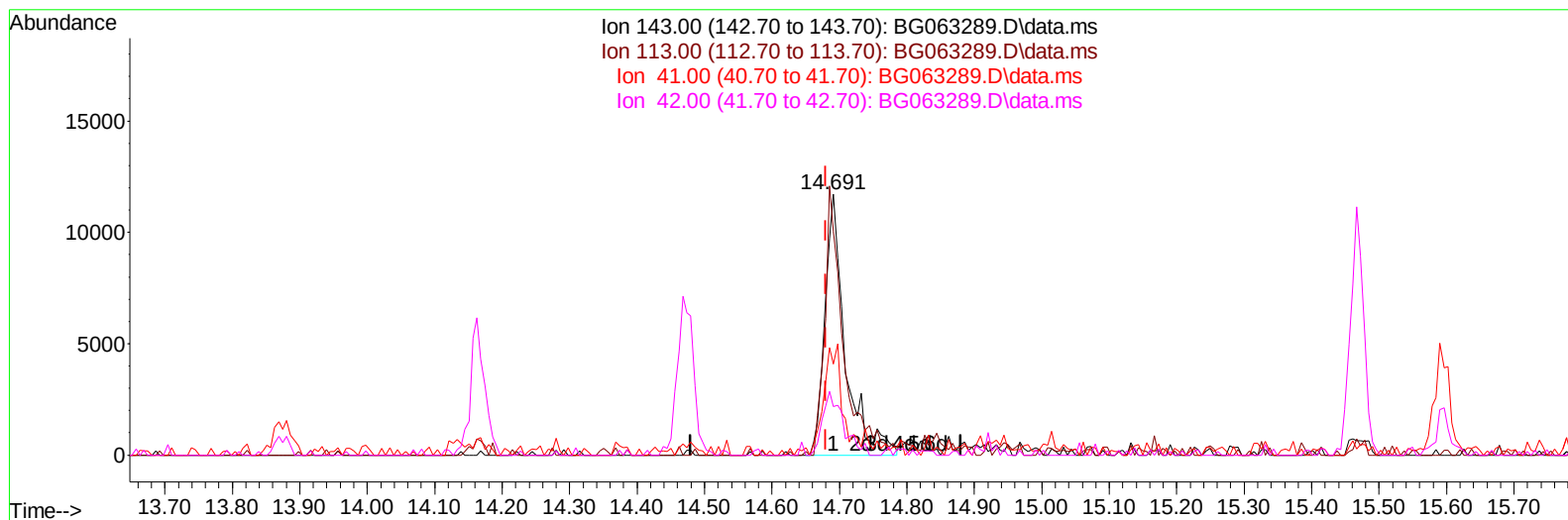
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TIC: BG063289.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.691min (+ 0.011) 5.83 ng/ul m

response 24579

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	123.00	85.97#
41.00	42.60	34.92
42.00	30.90	18.58#

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 QLast Update : Wed Nov 06 15:45:52 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.835	152	92876	20.000	ng/ul	0.00
20) Naphthalene-d8	10.625	136	452055	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.474	164	393396	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.224	188	922452	20.000	ng/ul	0.00
79) Chrysene-d12	21.477	240	935515	20.000	ng/ul	# 0.00
88) Perylene-d12	24.509	264	1005203	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	10866	5.290	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.012	99	66801	7.971	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.165	67	129392	26.322	ng/ul	0.00
11) 2-Chlorophenol-d4	7.370	132	139816	25.703	ng/ul	0.00
15) 4-Methylphenol-d8	8.540	113	135686	19.102	ng/ul	0.00
21) Nitrobenzene-d5	8.986	128	86332	27.165	ng/ul	0.00
24) 2-Nitrophenol-d4	9.709	143	106701	25.977	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.249	165	212463	26.357	ng/ul	0.00
31) 4-Chloroaniline-d4	10.761	131	223028	21.242	ng/ul	0.00
46) Dimethylphthalate-d6	13.880	166	756491	24.550	ng/ul	0.00
49) Acenaphthylene-d8	14.162	160	816226	27.062	ng/ul	0.00
54) 4-Nitrophenol-d4	14.691	143	24579m	5.826	ng/ul	0.01
60) Fluorene-d10	15.467	176	704155	27.076	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.590	200	126109	21.753	ng/ul	0.00
73) Anthracene-d10	17.323	188	1162451	29.321	ng/ul	0.00
81) Pyrene-d10	19.621	212	1466688	31.270	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.304	264	1447221	29.820	ng/ul	0.00

Target Compounds Qvalue

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

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