

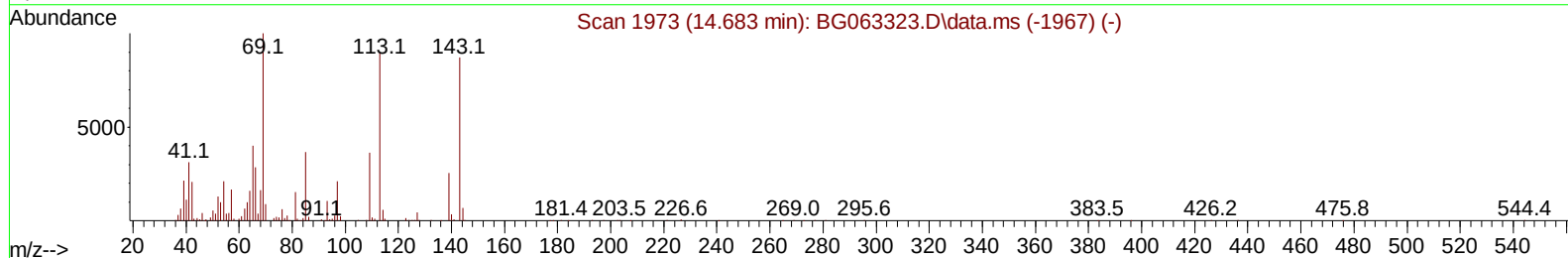
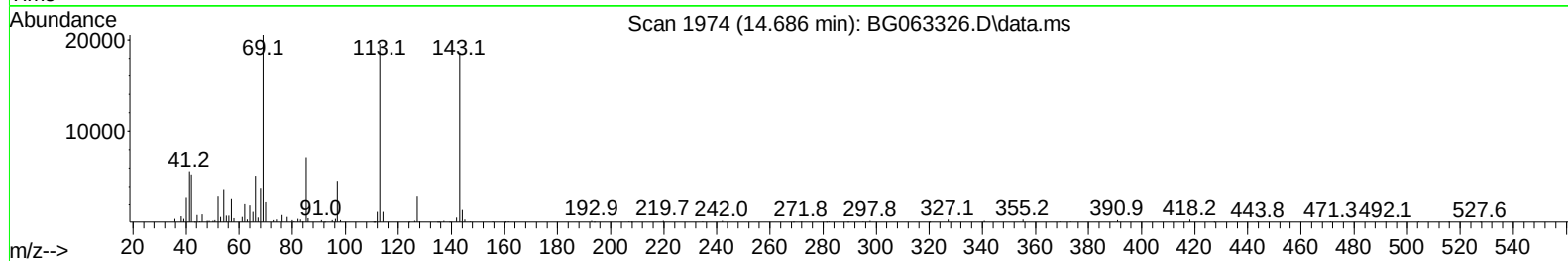
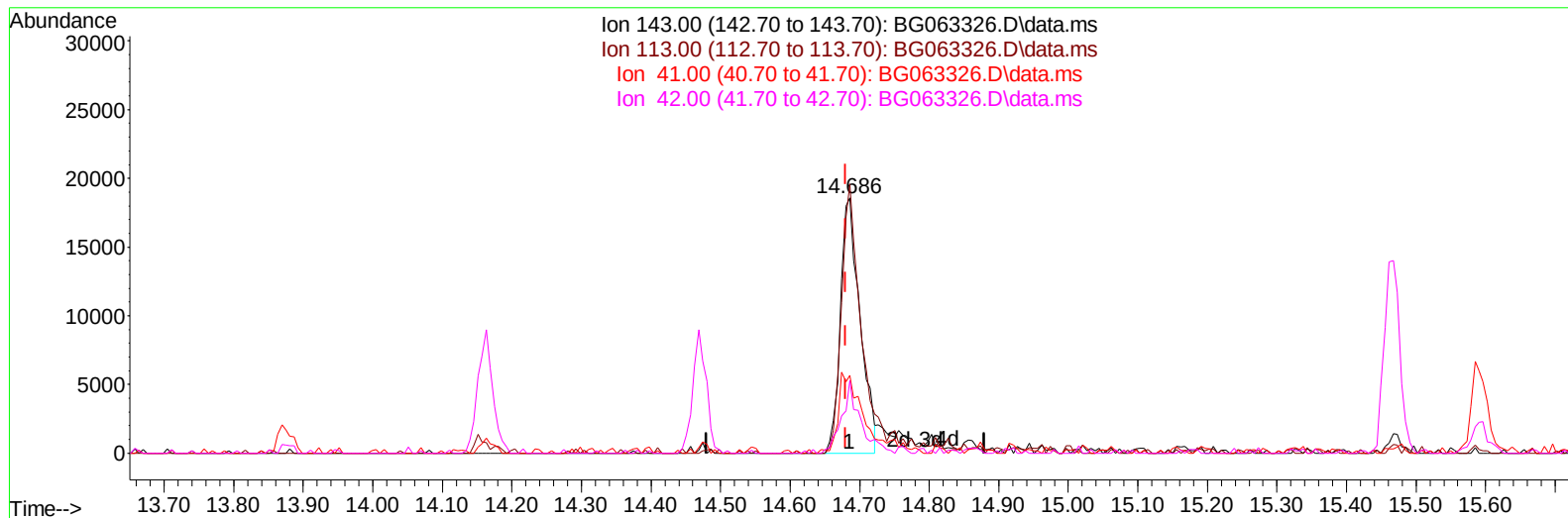
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111124\
Data File : BG063326.D
Acq On : 12 Nov 2024 15:17
Operator : RC/JU
Sample : P4801-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GCPA2

Manual IntegrationsAPPROVED

Quant Time: Nov 12 15:49:23 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 11/13/2024
Supervised By :mohammad ahmed 11/14/2024



TIC: BG063326.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.686min (+ 0.006) 8.03 ng/u1

response 36558

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	123.00	105.38
41.00	42.60	30.51#
42.00	30.90	28.56

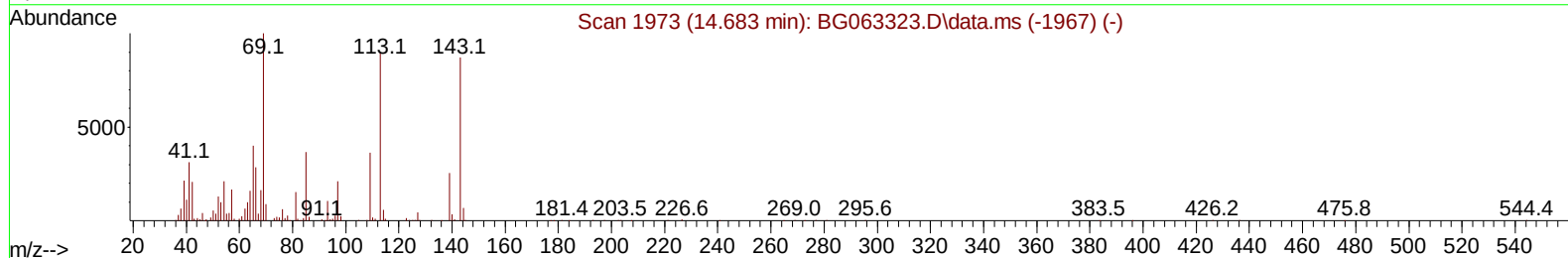
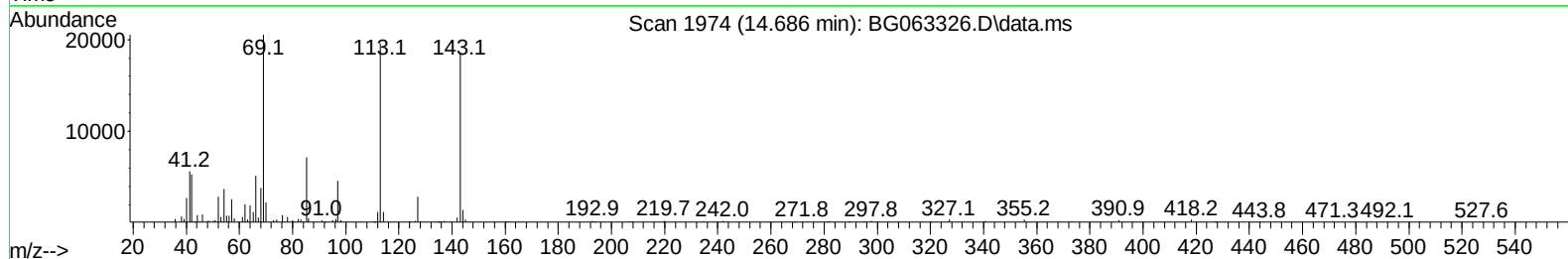
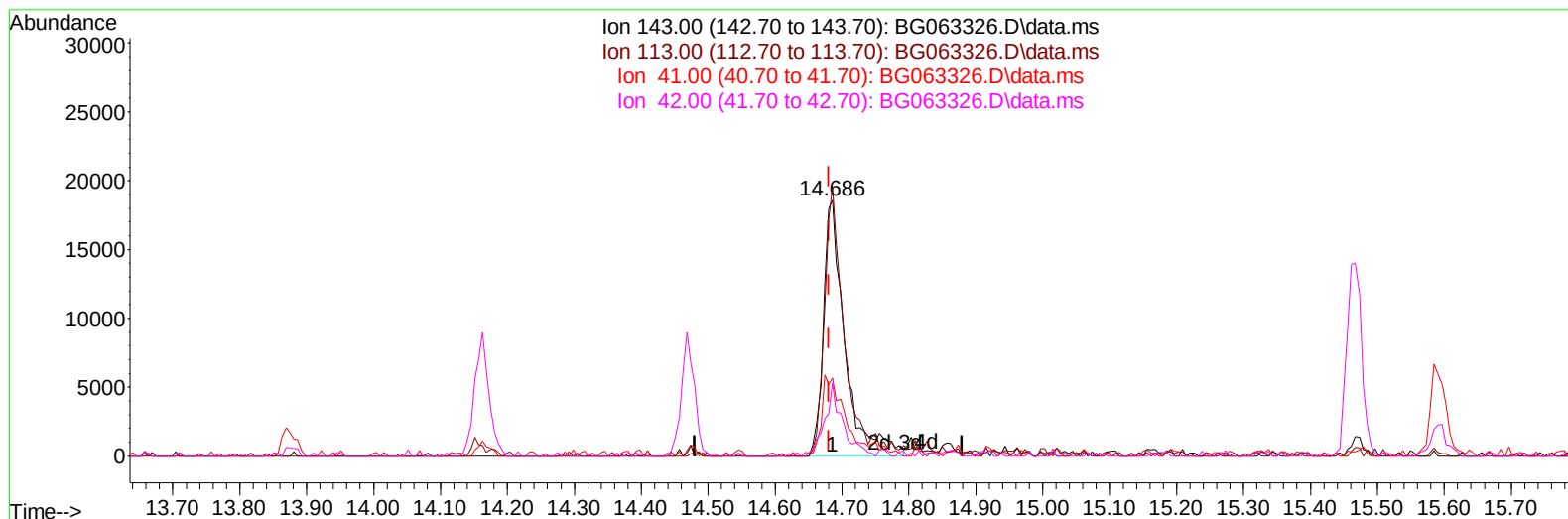
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(54) 4-Nitrophenol-d4 (S)

14.686min (+ 0.006) 9.17 ng/ul m

response 41775

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	123.00	105.38
41.00	42.60	30.51#
42.00	30.90	28.56

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.835	152	100162	20.000	ng/ul	0.00
20) Naphthalene-d8	10.626	136	486542	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.468	164	424612	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.224	188	995829	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.478	240	1030527	20.000	ng/ul	# 0.00
88) Perylene-d12	24.510	264	1083726	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	11804	5.329	ng/uL	0.00
4) Pyridine-d5	3.705	84	59112	8.074	ng/ul	0.00
7) Phenol-d5	7.007	99	96574	10.685	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.165	67	166024	31.317	ng/ul	0.00
11) 2-Chlorophenol-d4	7.371	132	174409	29.730	ng/ul	0.00
15) 4-Methylphenol-d8	8.540	113	181299	23.666	ng/ul	0.00
21) Nitrobenzene-d5	8.987	128	106843	31.236	ng/ul	0.00
24) 2-Nitrophenol-d4	9.703	143	134690	30.467	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.250	165	267856	30.874	ng/ul	0.00
31) 4-Chloroaniline-d4	10.761	131	289546	25.623	ng/ul	0.00
46) Dimethylphthalate-d6	13.875	166	1017330	30.587	ng/ul	0.00
49) Acenaphthylene-d8	14.163	160	1094303	33.615	ng/ul	0.00
54) 4-Nitrophenol-d4	14.686	143	41775m	9.174	ng/ul	0.00
60) Fluorene-d10	15.467	176	950106	33.847	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.591	200	191052	30.527	ng/ul	0.00
73) Anthracene-d10	17.318	188	1638462	38.282	ng/ul	0.00
81) Pyrene-d10	19.621	212	1985350	38.425	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.304	264	2042941	39.044	ng/ul	0.00

Target Compounds Qvalue

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

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