

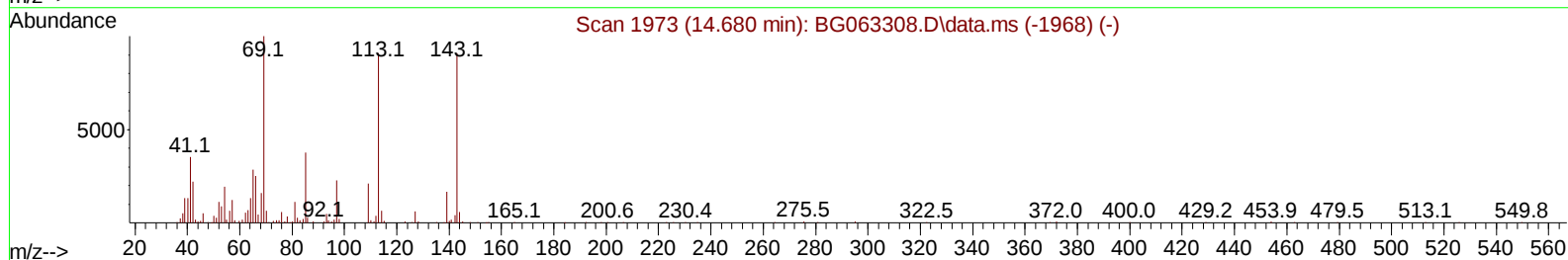
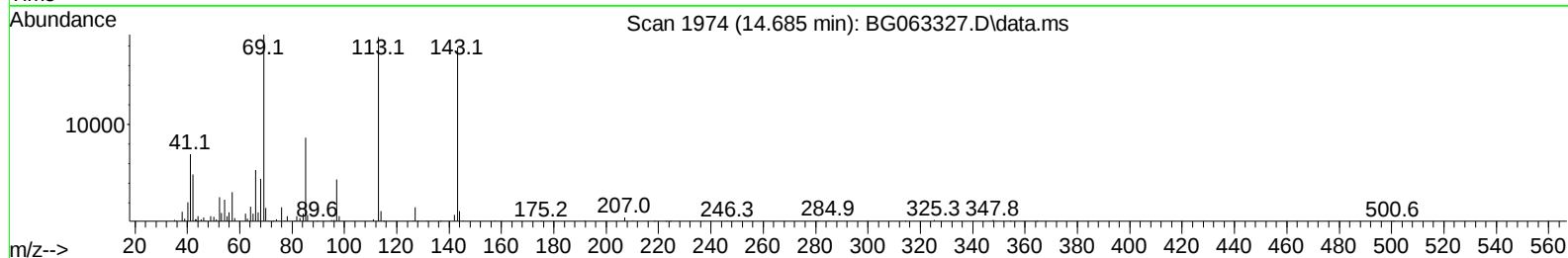
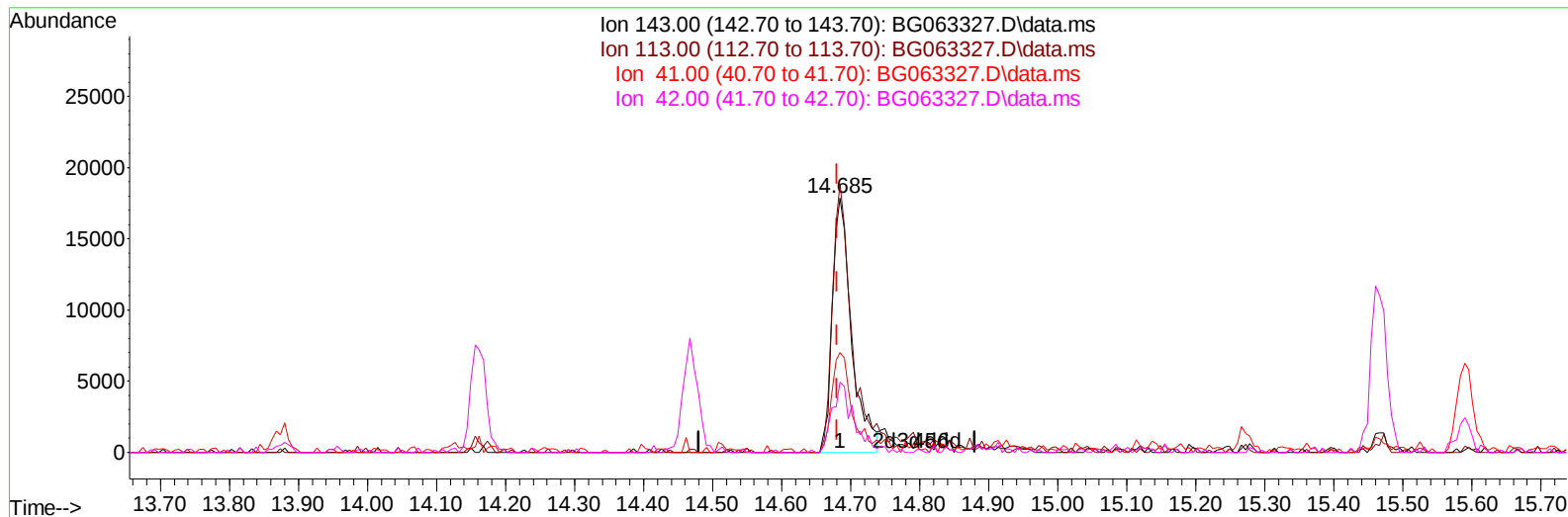
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111124\
Data File : BG063327.D
Acq On : 12 Nov 2024 15:55
Operator : RC/JU
Sample : P4801-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GCPA3

Manual IntegrationsAPPROVED

Quant Time: Nov 12 15:54:37 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: BG063327.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.685min (+ 0.005) 7.93 ng/u1

response 34337

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	123.00	105.55
41.00	42.60	39.17
42.00	30.90	27.43

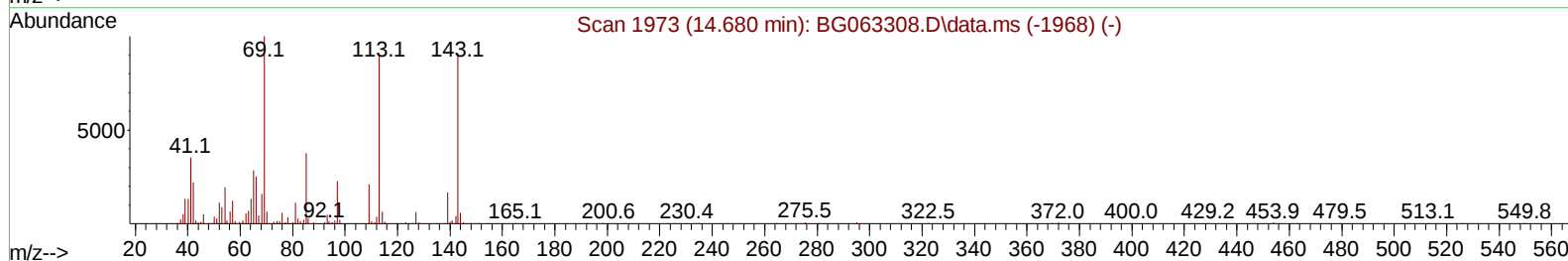
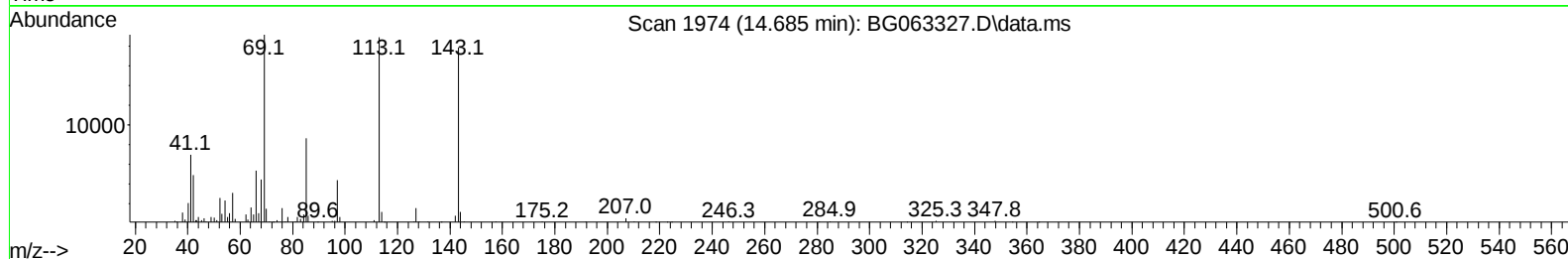
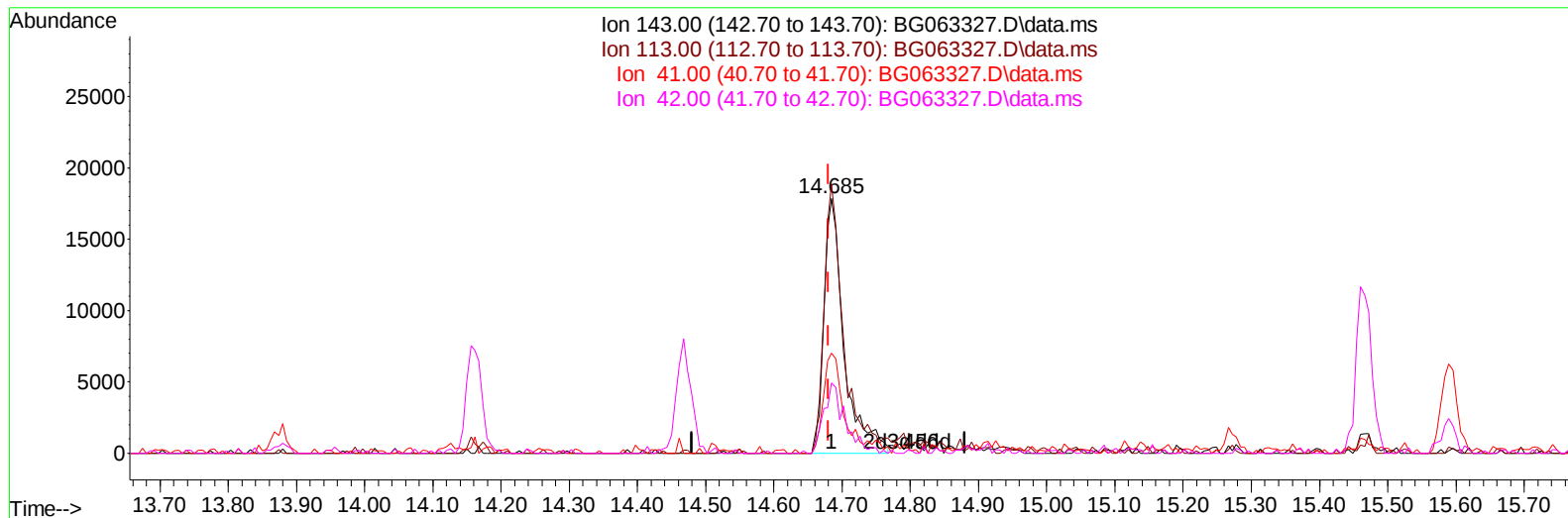
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(54) 4-Nitrophenol-d4 (S)

14.685min (+ 0.005) 8.38 ng/ul m

response 36267

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	123.00	105.55
41.00	42.60	39.17
42.00	30.90	27.43

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Instrument :
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 GCPA3

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/13/2024
 Supervised By :mohammad ahmed 11/14/2024

Quant Time: Nov 12 23:09:50 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

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.834	152	95994	20.000	ng/ul	0.00
20) Naphthalene-d8	10.625	136	456591	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.467	164	403646	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.223	188	964327	20.000	ng/ul	0.00
79) Chrysene-d12	21.471	240	977060	20.000	ng/ul	# 0.00
88) Perylene-d12	24.508	264	1035849	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.298	96	8570	4.037	ng/uL	0.00
4) Pyridine-d5	3.703	84	35867	5.112	ng/ul	0.00
7) Phenol-d5	7.005	99	75914	8.764	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.164	67	143134	28.172	ng/ul	0.00
11) 2-Chlorophenol-d4	7.370	132	157422	28.000	ng/ul	0.00
15) 4-Methylphenol-d8	8.539	113	152329	20.748	ng/ul	0.00
21) Nitrobenzene-d5	8.985	128	92486	28.812	ng/ul	0.00
24) 2-Nitrophenol-d4	9.708	143	119923	28.906	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.249	165	248154	30.479	ng/ul	0.00
31) 4-Chloroaniline-d4	10.760	131	238572	22.497	ng/ul	0.00
46) Dimethylphthalate-d6	13.880	166	900817	28.491	ng/ul	0.00
49) Acenaphthylene-d8	14.162	160	966627	31.235	ng/ul	0.00
54) 4-Nitrophenol-d4	14.685	143	36267m	8.378	ng/ul	0.00
60) Fluorene-d10	15.466	176	837158	31.373	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.589	200	178124	29.391	ng/ul	0.00
73) Anthracene-d10	17.323	188	1524661	36.787	ng/ul	0.00
81) Pyrene-d10	19.620	212	1863169	38.034	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.303	264	1904431	38.079	ng/ul	0.00

Target Compounds Qvalue

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

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