

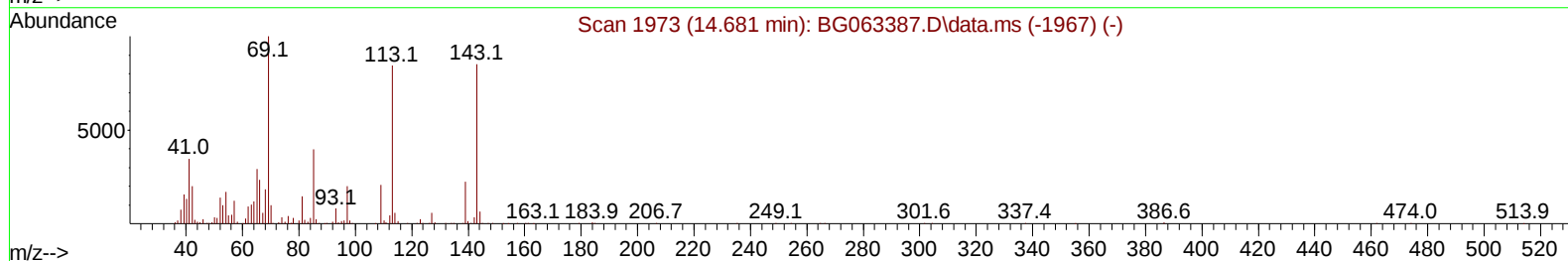
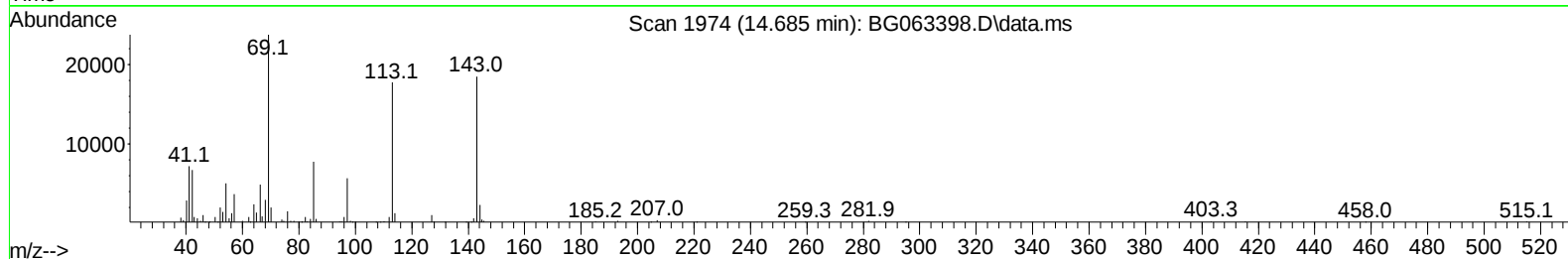
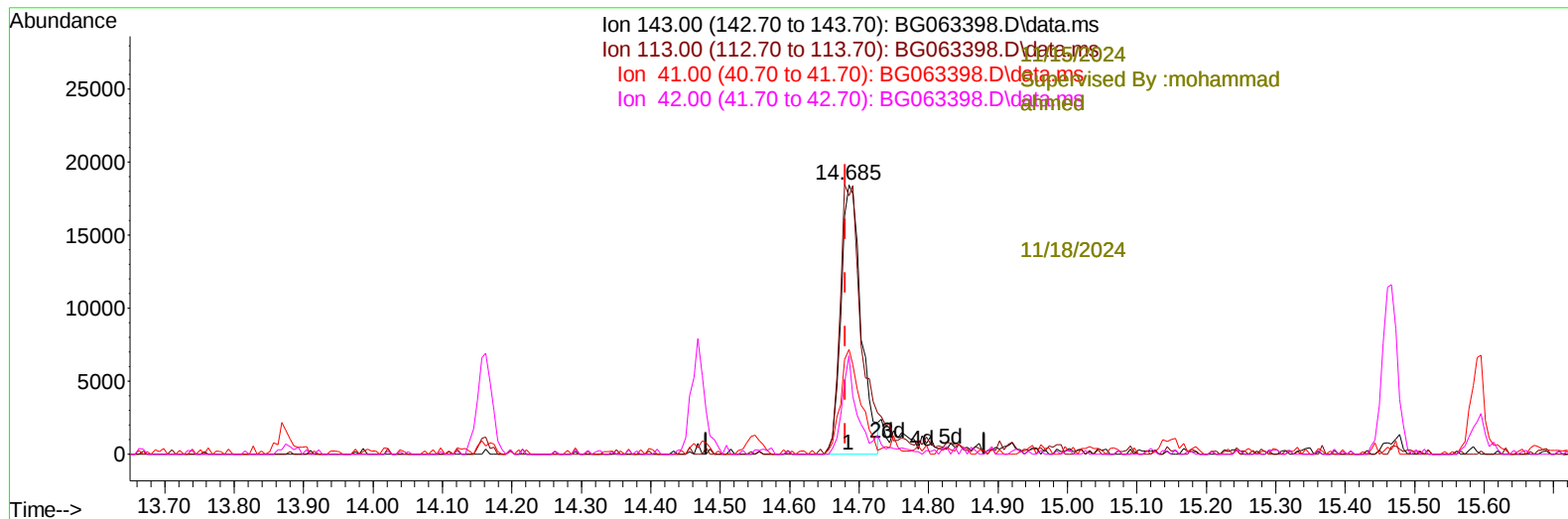
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111324\
Data File : BG063398.D
Acq On : 14 Nov 2024 19:19
Operator : RC/JU
Sample : P4715-13
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COAK6

Manual IntegrationsAPPROVED

Quant Time: Nov 14 22:51:23 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 11/15/2024
Supervised By :mohammad ahmed 11/18/2024



TIC: BG063398.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.685min (+ 0.005) 9.24 ng/u1

response 37093

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	123.00	95.93#
41.00	42.60	38.82
42.00	30.90	36.47

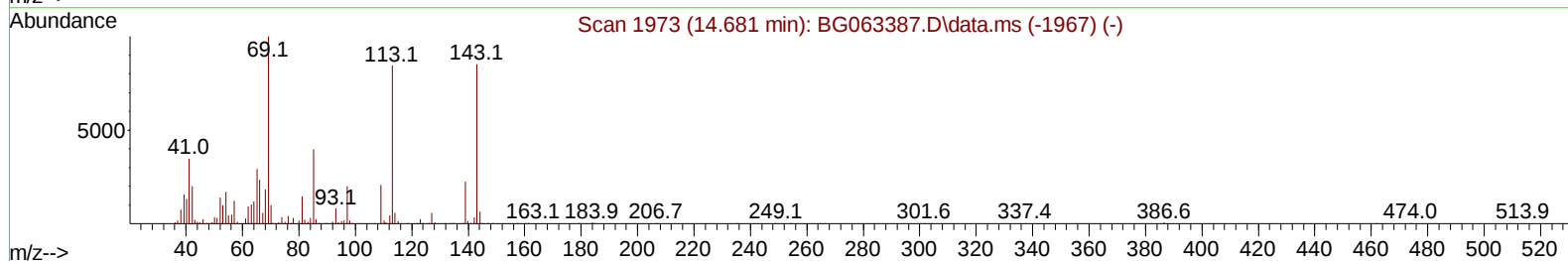
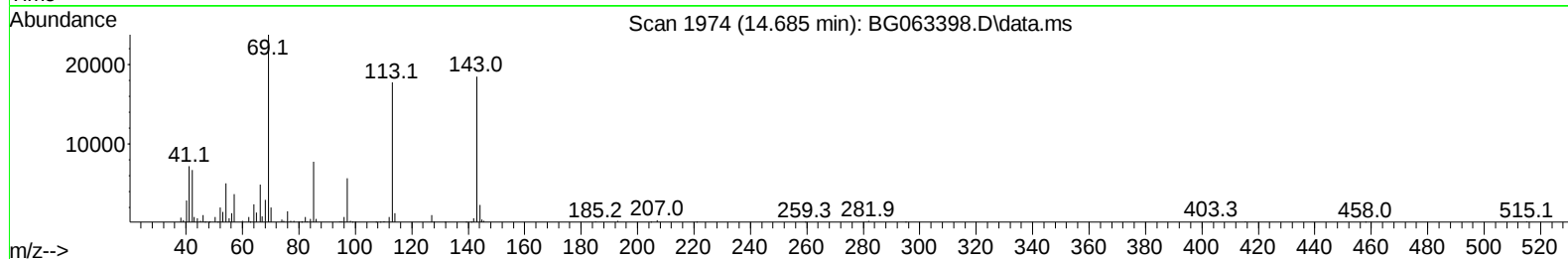
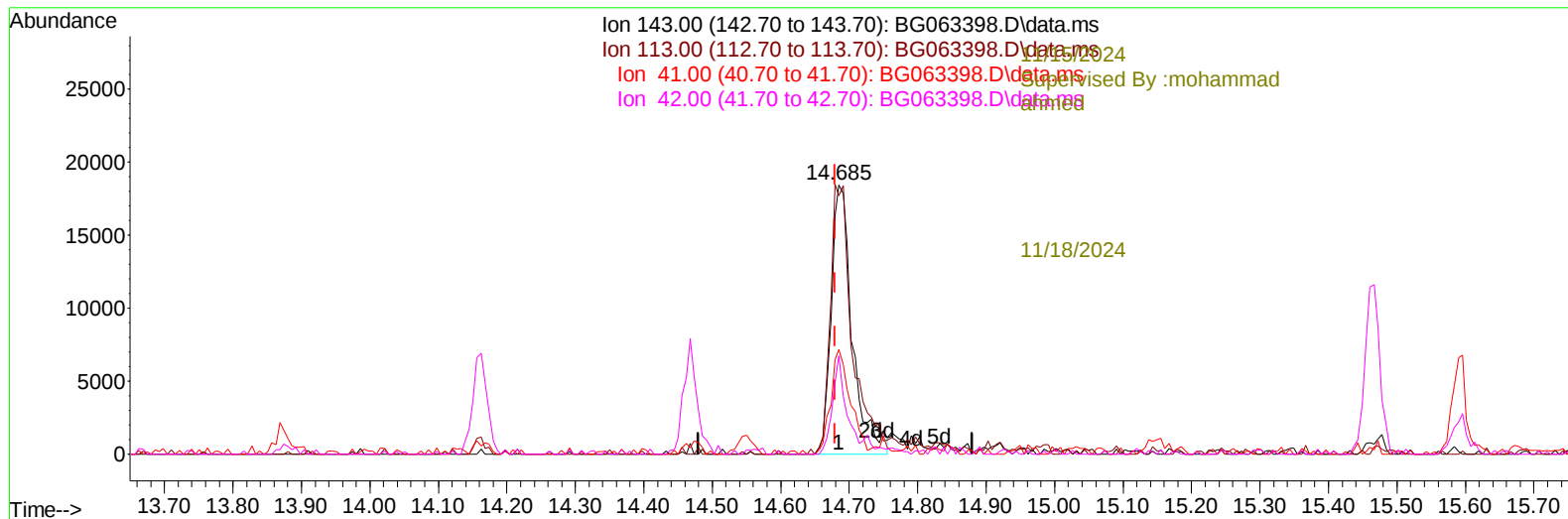
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TIC: BG063398.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.685min (+ 0.005) 9.86 ng/ul m

response 39570

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	123.00	95.93#
41.00	42.60	38.82
42.00	30.90	36.47

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 C0AK6

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/15/2024
 Supervised By :mohammad ahmed 11/18/2024

Quant Time: Nov 14 22:52:08 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)
-----11/15/2024						
Supervised By :mohammad						
ahmed						
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.834	152	108061	20.000	ng/ul	0.00
20) Naphthalene-d8	10.619	136	449436	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.468	164	374255	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.223	188	929600	20.000	ng/ul	0.00
79) Chrysene-d12	21.471	240	973018	20.000	ng/ul	# 0.00
88) Perylene-d12	24.509	264	1031721	20.000	ng/ul	0.0011/18/2024
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.304	96	10542	4.411	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.006	99	116391	11.937	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.165	67	189479	33.129	ng/ul	0.00
11) 2-Chlorophenol-d4	7.370	132	209324	33.074	ng/ul	0.00
15) 4-Methylphenol-d8	8.539	113	196450	23.770	ng/ul	0.00
21) Nitrobenzene-d5	8.986	128	111337	35.237	ng/ul	0.00
24) 2-Nitrophenol-d4	9.709	143	138536	33.924	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.249	165	253225	31.597	ng/ul	0.00
31) 4-Chloroaniline-d4	10.760	131	160567	15.382	ng/ul	0.00
46) Dimethylphthalate-d6	13.874	166	903947	30.835	ng/ul	0.00
49) Acenaphthylene-d8	14.162	160	963614	33.583	ng/ul	0.00
54) 4-Nitrophenol-d4	14.685	143	39570m	9.859	ng/ul	0.00
60) Fluorene-d10	15.467	176	819870	33.138	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.590	200	172975	29.607	ng/ul	0.00
73) Anthracene-d10	17.317	188	1466004	36.693	ng/ul	0.00
81) Pyrene-d10	19.620	212	1879381	38.524	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.303	264	1860886	37.358	ng/ul	0.00

Target Compounds Qvalue

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

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