

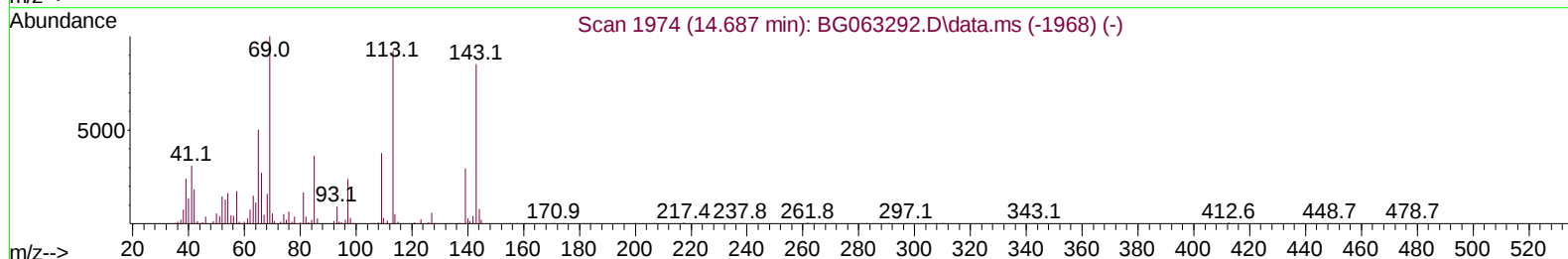
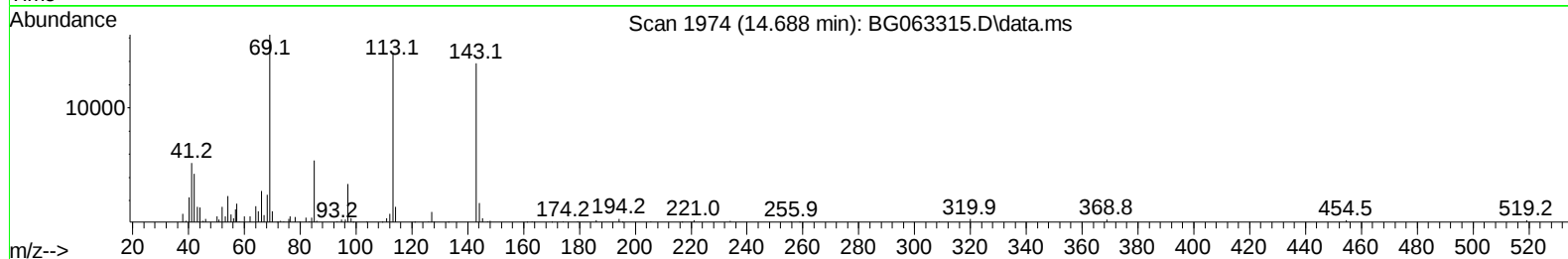
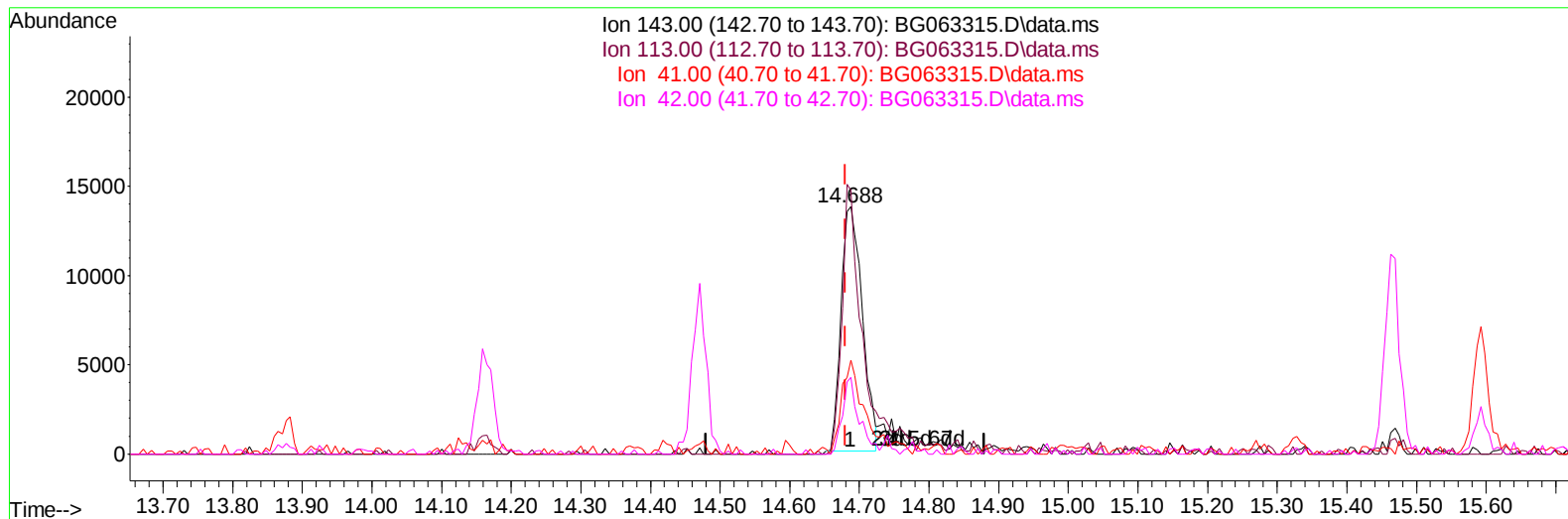
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111124\
Data File : BG063315.D
Acq On : 12 Nov 2024 8:19
Operator : RC/JU
Sample : P4714-15
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AH9

Manual IntegrationsAPPROVED

Quant Time: Nov 12 07:54:16 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: BG063315.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.688min (+ 0.008) 5.85 ng/u1

response 29115

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	123.00	106.40
41.00	42.60	38.01
42.00	30.90	30.97

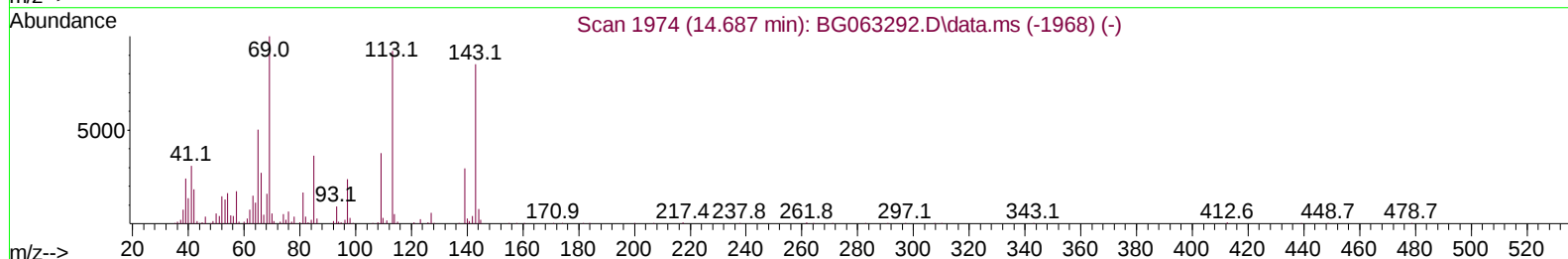
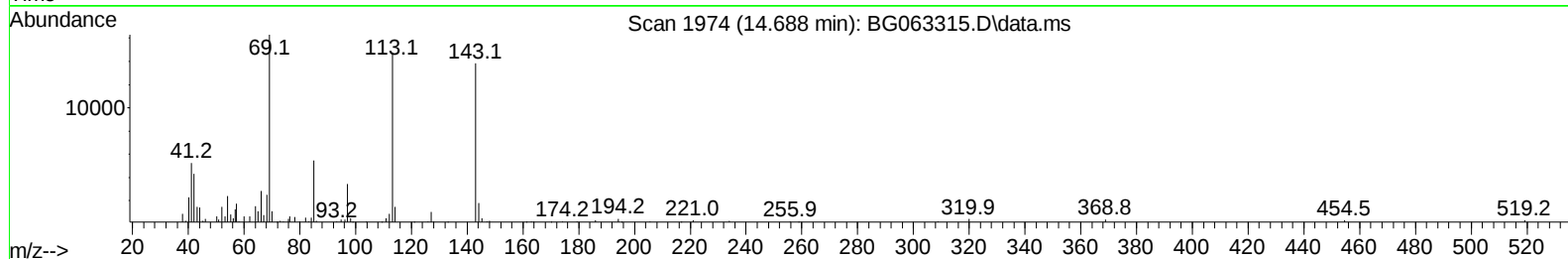
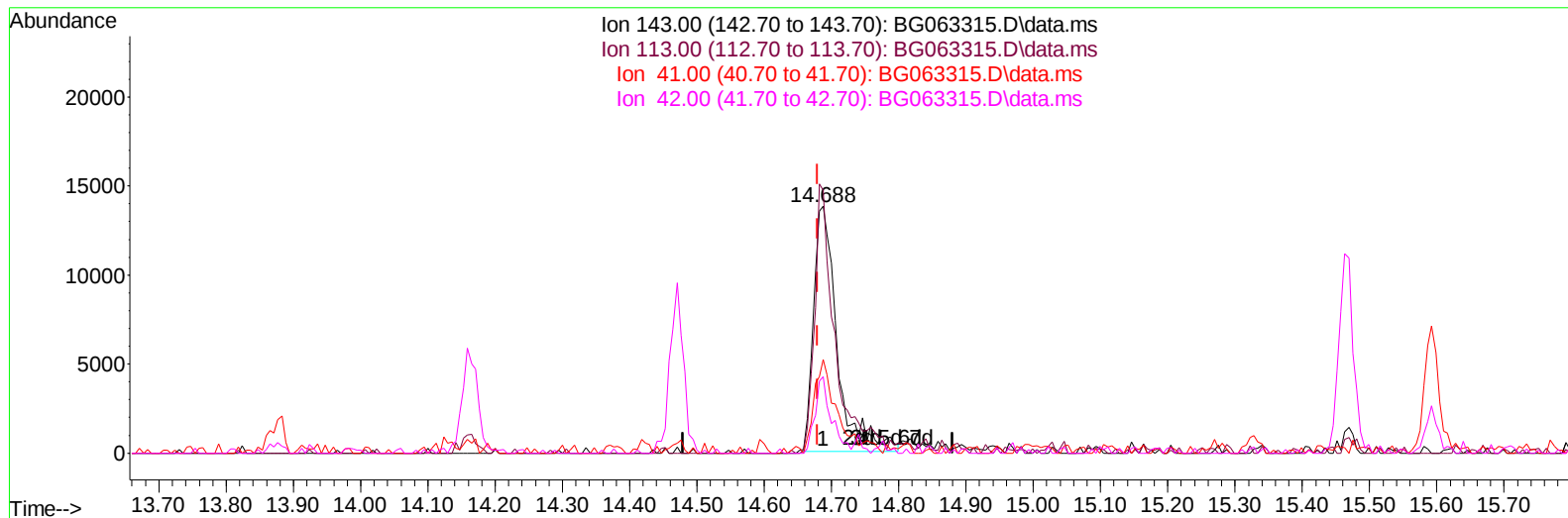
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TIC: BG063315.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.688min (+ 0.008) 6.77 ng/ul m

response 33666

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	123.00	106.40
41.00	42.60	38.01
42.00	30.90	30.97

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Reviewed By :Yogesh Patel 11/13/2024
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Quant Time: Nov 12 07:55:03 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	7.837	152	109971	20.000	ng/ul	0.00
20) Naphthalene-d8	10.622	136	530355	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.470	164	463829	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.220	188	1103421	20.000	ng/ul	0.00
79) Chrysene-d12	21.474	240	1129540	20.000	ng/ul	# 0.00
88) Perylene-d12	24.512	264	1217258	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.301	96	9339	3.840	ng/uL	0.00
4) Pyridine-d5	3.707	84	38495	4.789	ng/ul	0.00
7) Phenol-d5	7.009	99	69561	7.010	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.167	67	125594	21.578	ng/ul	0.00
11) 2-Chlorophenol-d4	7.367	132	140860	21.870	ng/ul	0.00
15) 4-Methylphenol-d8	8.542	113	136613	16.242	ng/ul	0.00
21) Nitrobenzene-d5	8.983	128	84441	22.647	ng/ul	0.00
24) 2-Nitrophenol-d4	9.711	143	102330	21.235	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.246	165	203477	21.516	ng/ul	0.00
31) 4-Chloroaniline-d4	10.757	131	228771	18.572	ng/ul	0.00
46) Dimethylphthalate-d6	13.877	166	776981	21.386	ng/ul	0.00
49) Acenaphthylene-d8	14.165	160	808962	22.749	ng/ul	0.00
54) 4-Nitrophenol-d4	14.688	143	33666m	6.768	ng/ul	0.00
60) Fluorene-d10	15.469	176	730392	23.820	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.593	200	164424	23.710	ng/ul	0.00
73) Anthracene-d10	17.320	188	1433206	30.221	ng/ul	0.00
81) Pyrene-d10	19.623	212	1785027	31.520	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.306	264	1806308	30.735	ng/ul	0.00

Target Compounds Qvalue

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

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