

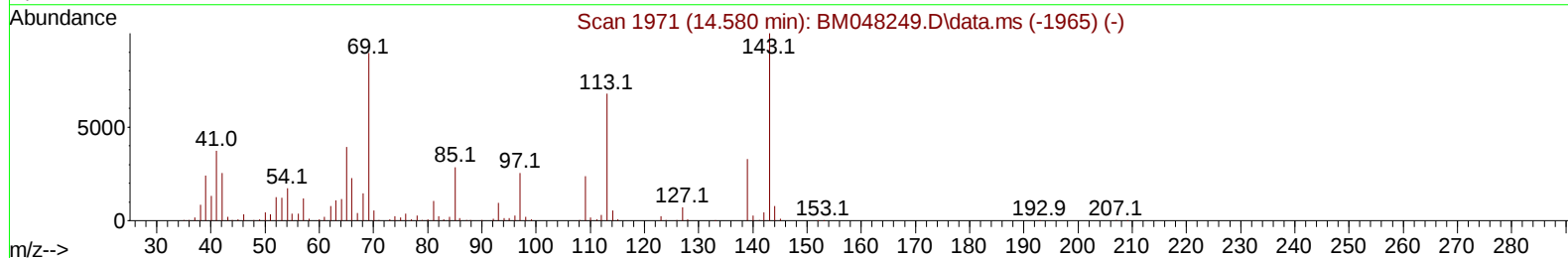
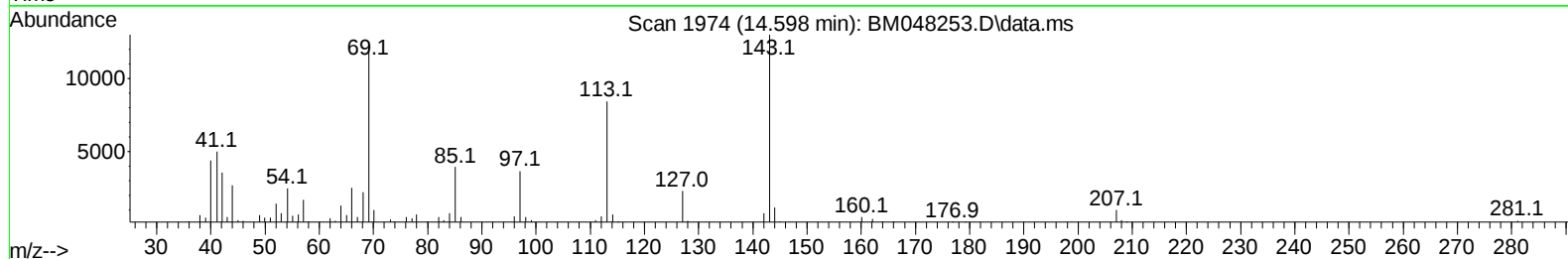
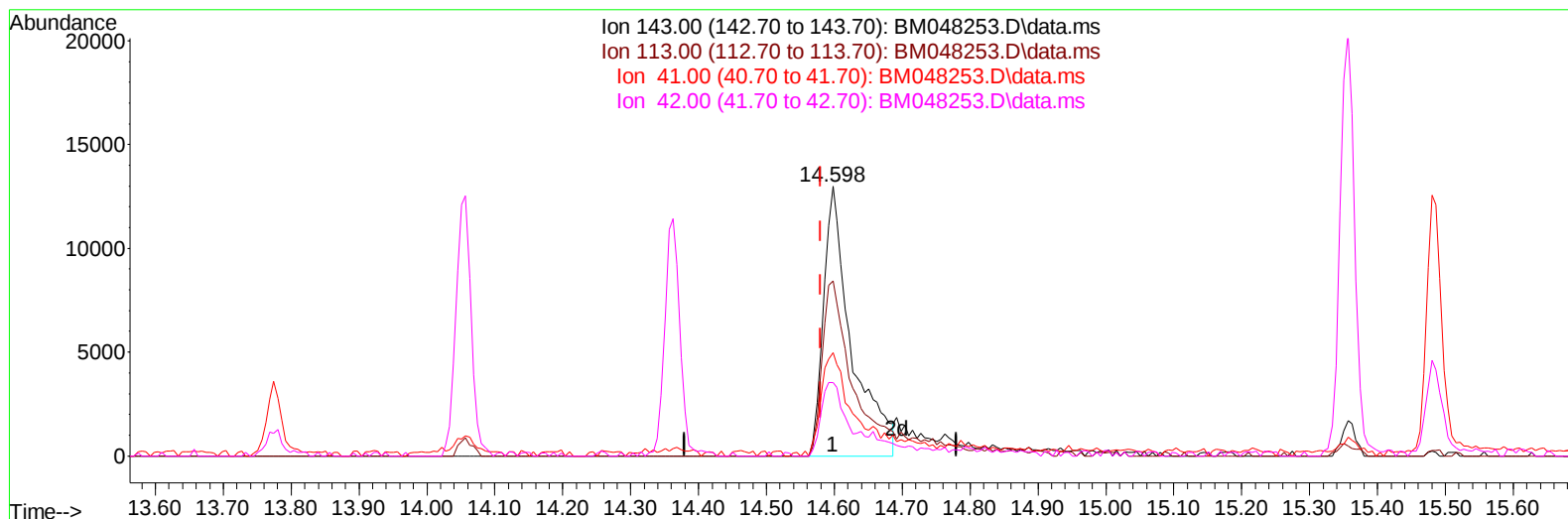
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102824\
Data File : BM048253.D
Acq On : 28 Oct 2024 13:21
Operator : RC/JU
Sample : P4559-15
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCNZ0

Manual IntegrationsAPPROVED

Quant Time: Oct 28 14:10:23 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM102624.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sun Oct 27 23:39:58 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/29/2024
Supervised By :mohammad ahmed 11/01/2024



TIC: BM048253.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.598min (+ 0.018) 4.24 ng/u1

response 36972

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	68.30	64.90
41.00	40.60	38.25
42.00	26.40	27.23

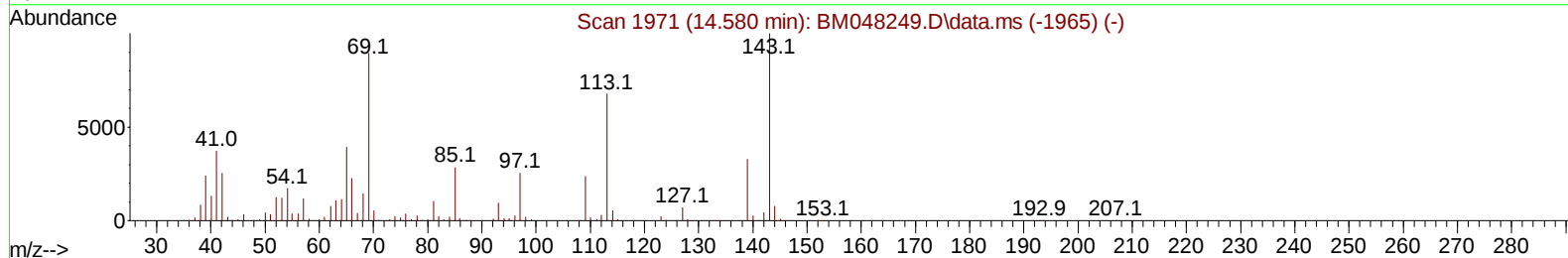
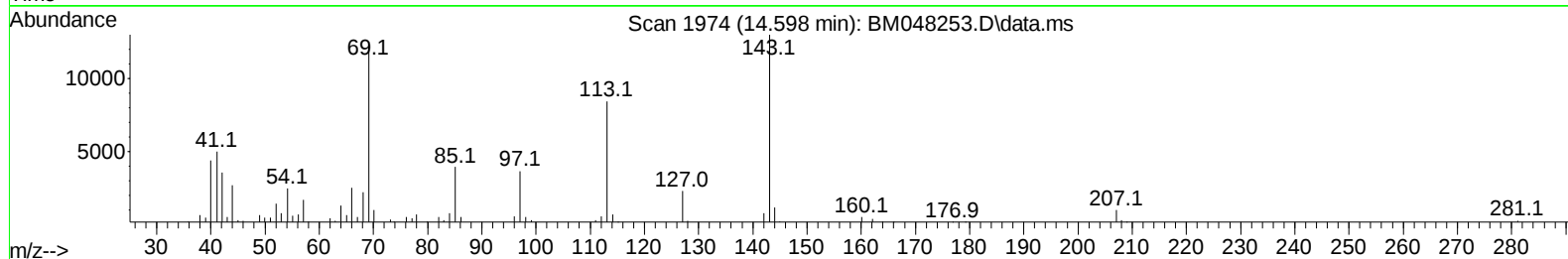
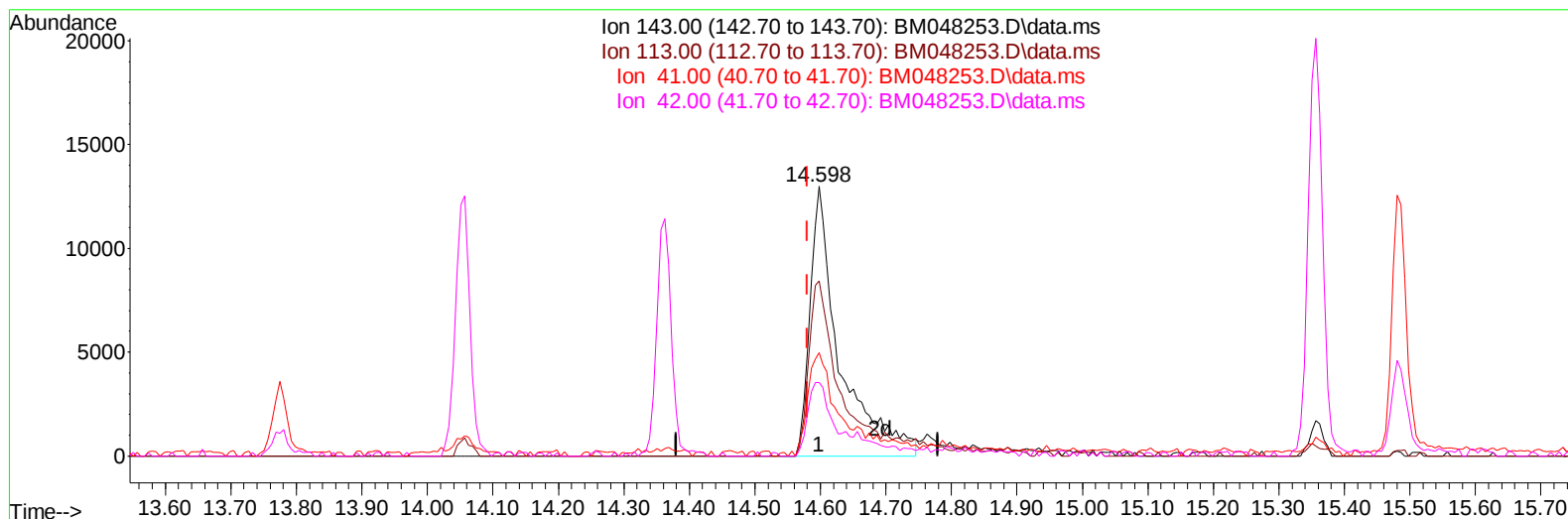
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TIC: BM048253.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.598min (+ 0.018) 4.70 ng/ul m

response 40911

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	68.30	64.90
41.00	40.60	38.25
42.00	26.40	27.23

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Instrument :
 BNA_M
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 GCNZ0

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/29/2024
 Supervised By :mohammad ahmed 11/01/2024

Quant Time: Oct 28 15:38:19 2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Oct 27 23:39:58 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.722	152	230411	20.000	ng/ul	0.00
20) Naphthalene-d8	10.510	136	983030	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.363	164	637522	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.104	188	1254715	20.000	ng/ul	0.00
79) Chrysene-d12	21.333	240	1016848	20.000	ng/ul	0.00
88) Perylene-d12	24.274	264	1058113	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.169	96	25699	4.391	ng/uL	0.00
4) Pyridine-d5	3.593	84	85101	5.017	ng/ul	0.00
7) Phenol-d5	6.898	99	143459	7.247	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.051	67	385087	31.405	ng/ul	0.00
11) 2-Chlorophenol-d4	7.257	132	443479	27.549	ng/ul	0.00
15) 4-Methylphenol-d8	8.434	113	289130	18.193	ng/ul	0.00
21) Nitrobenzene-d5	8.875	128	249670	30.978	ng/ul	0.00
24) 2-Nitrophenol-d4	9.598	143	277927	31.832	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.139	165	455083	28.563	ng/ul	0.00
31) 4-Chloroaniline-d4	10.651	131	525122	23.736	ng/ul	0.00
46) Dimethylphthalate-d6	13.774	166	1562770	32.802	ng/ul	0.00
49) Acenaphthylene-d8	14.057	160	1715295	31.931	ng/ul	0.00
54) 4-Nitrophenol-d4	14.598	143	40911m	4.696	ng/ul	0.02
60) Fluorene-d10	15.357	176	1346164	33.439	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.486	200	239666	28.903	ng/ul	0.00
73) Anthracene-d10	17.204	188	2105689	35.491	ng/ul	0.00
81) Pyrene-d10	19.498	212	2292685	37.155	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.080	264	2006331	35.949	ng/ul	0.00

Target Compounds Qvalue

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

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