

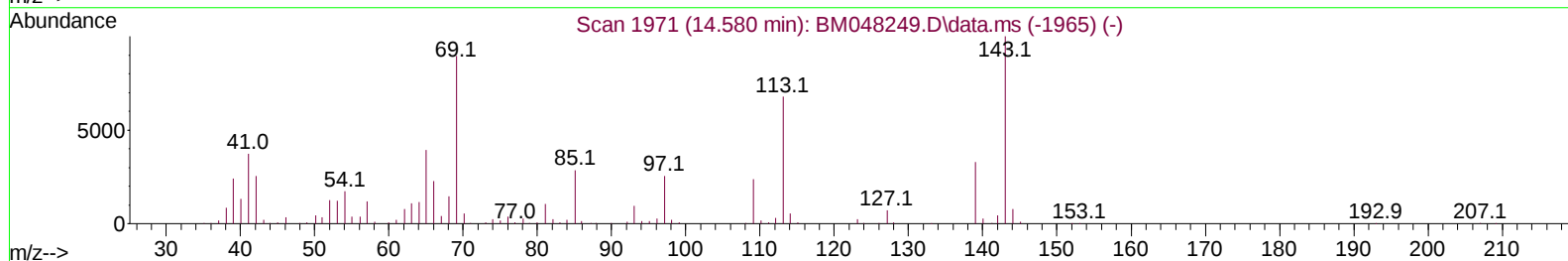
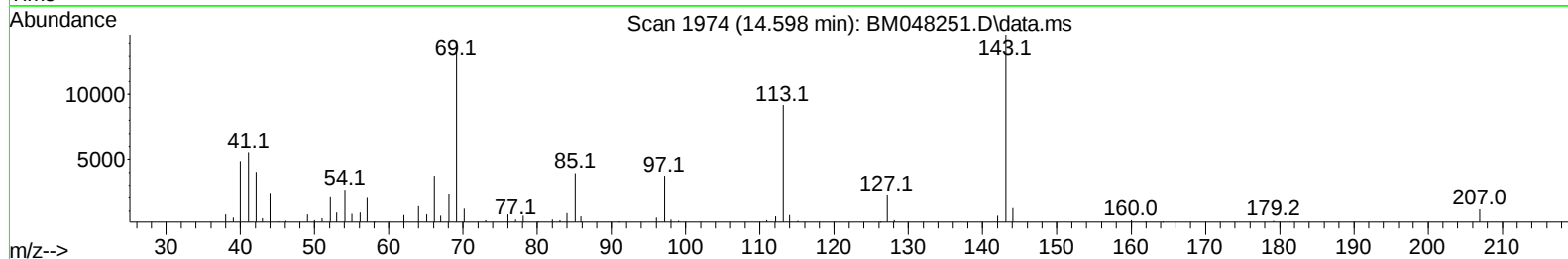
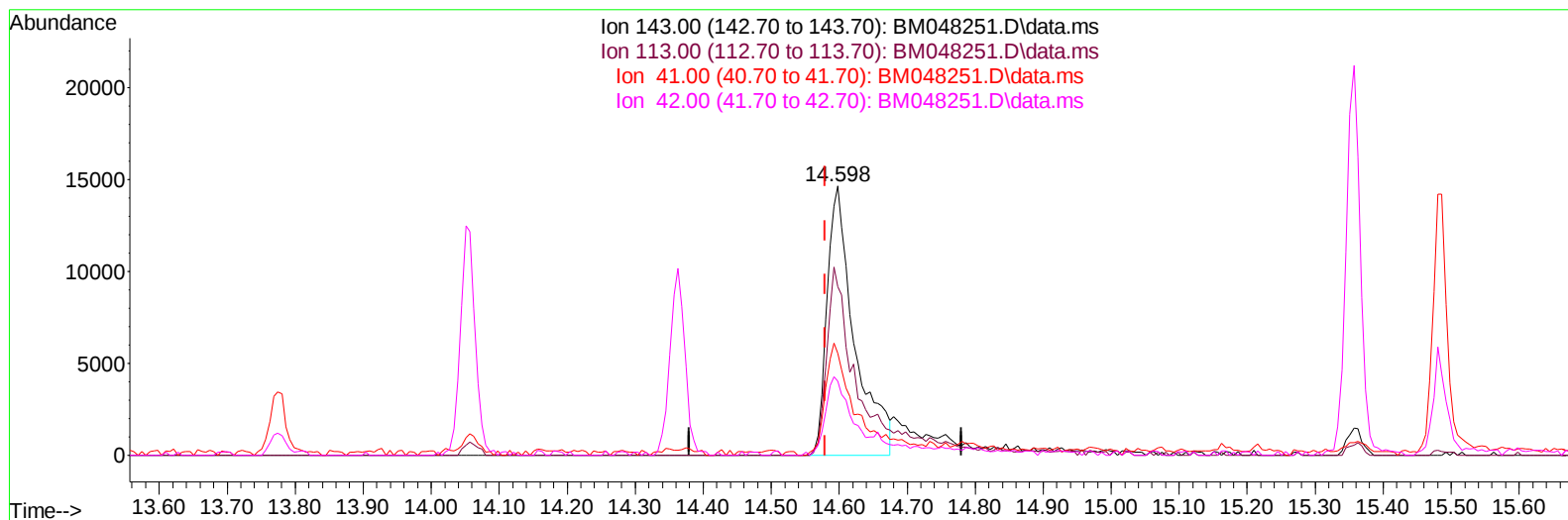
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102824\
Data File : BM048251.D
Acq On : 28 Oct 2024 12:02
Operator : RC/JU
Sample : P4559-13
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCNY8

Manual IntegrationsAPPROVED

Quant Time: Oct 28 12:40:24 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: BM048251.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.598min (+ 0.018) 5.35 ng/u1

response 40960

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	68.30	62.69
41.00	40.60	37.94
42.00	26.40	27.59

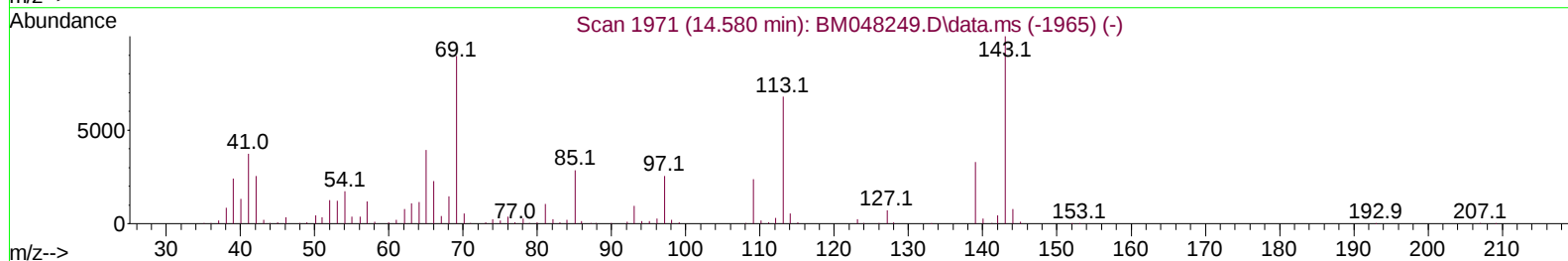
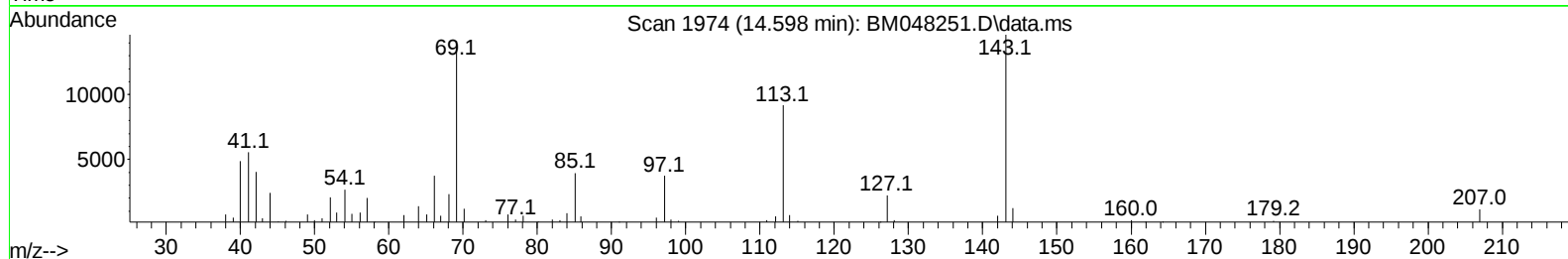
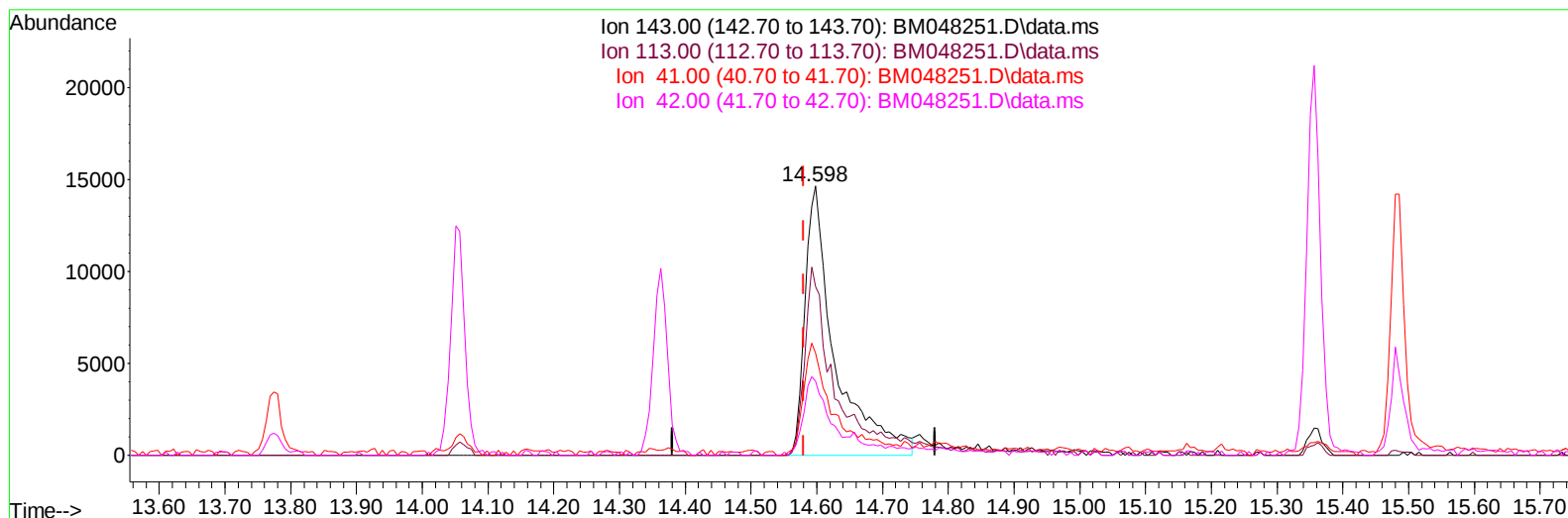
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TIC: BM048251.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.598min (+ 0.018) 6.07 ng/ul m

response 46507

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	68.30	62.69
41.00	40.60	37.94
42.00	26.40	27.59

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.722	152	196032	20.000	ng/ul	0.00
20) Naphthalene-d8	10.510	136	846644	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.363	164	560569	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.104	188	1167014	20.000	ng/ul	0.00
79) Chrysene-d12	21.327	240	994796	20.000	ng/ul	0.00
88) Perylene-d12	24.280	264	998394	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.169	96	30010	6.027	ng/uL	0.00
4) Pyridine-d5	3.593	84	107623	7.457	ng/ul	0.00
7) Phenol-d5	6.898	99	146825	8.718	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.051	67	380424	36.465	ng/ul	0.00
11) 2-Chlorophenol-d4	7.257	132	437275	31.928	ng/ul	0.00
15) 4-Methylphenol-d8	8.434	113	286672	21.202	ng/ul	0.00
21) Nitrobenzene-d5	8.875	128	246849	35.562	ng/ul	0.00
24) 2-Nitrophenol-d4	9.598	143	275522	36.640	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.139	165	454972	33.156	ng/ul	0.00
31) 4-Chloroaniline-d4	10.651	131	569817	29.905	ng/ul	0.00
46) Dimethylphthalate-d6	13.774	166	1585706	37.852	ng/ul	0.00
49) Acenaphthylene-d8	14.057	160	1724828	36.516	ng/ul	0.00
54) 4-Nitrophenol-d4	14.598	143	46507m	6.071	ng/ul	0.02
60) Fluorene-d10	15.357	176	1370189	38.708	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.486	200	254948	33.057	ng/ul	0.00
73) Anthracene-d10	17.204	188	2202566	39.913	ng/ul	0.00
81) Pyrene-d10	19.498	212	2500213	41.417	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.080	264	2160288	41.023	ng/ul	0.00

Target Compounds Qvalue

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

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