

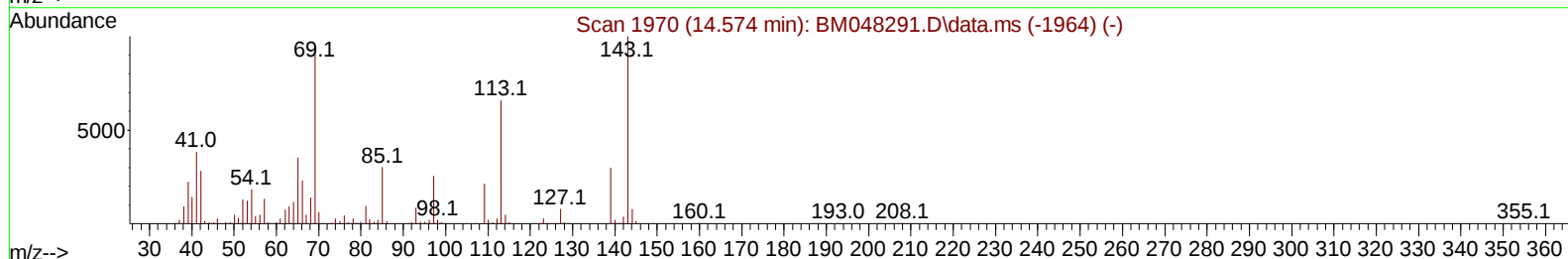
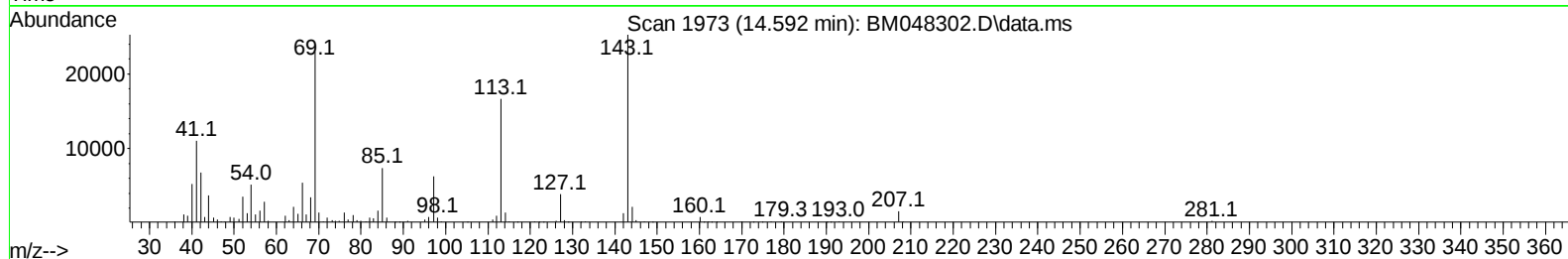
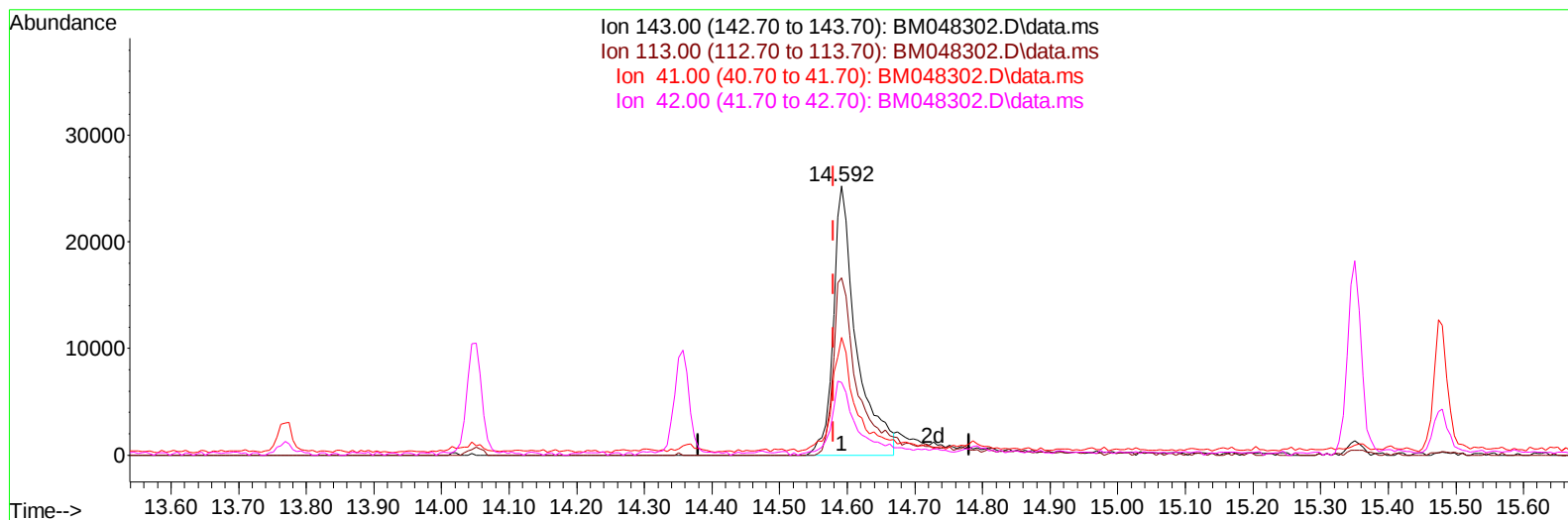
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102824\
Data File : BM048302.D
Acq On : 30 Oct 2024 00:36
Operator : RC/JU
Sample : P4557-01
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
E1L51

Manual IntegrationsAPPROVED

Quant Time: Oct 30 01:11:04 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 :Jagrut Upadhyay 10/30/2024
Supervised By :mohammad ahmed 11/01/2024



TIC: BM048302.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.592min (+ 0.012) 8.14 ng/u1

response 59587

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	68.30	65.82
41.00	40.60	43.67
42.00	26.40	27.04

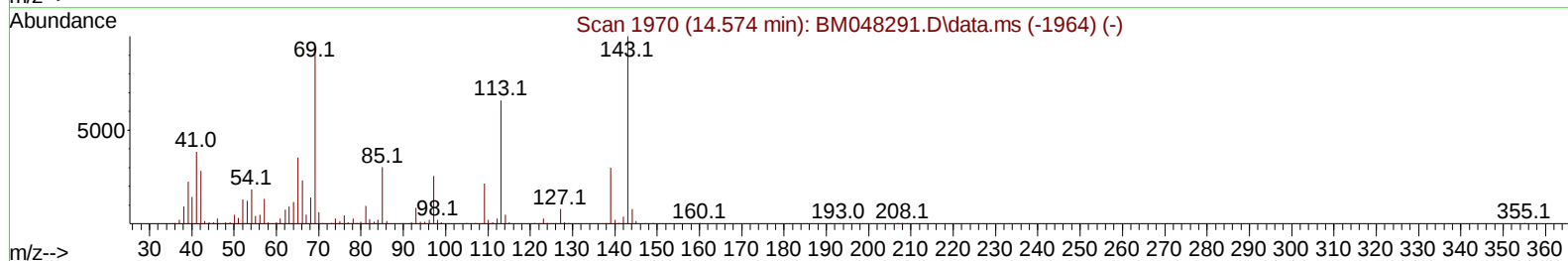
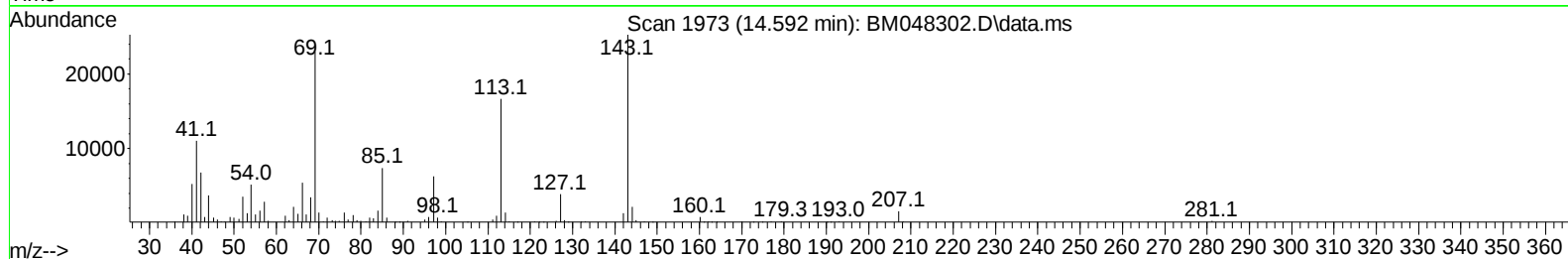
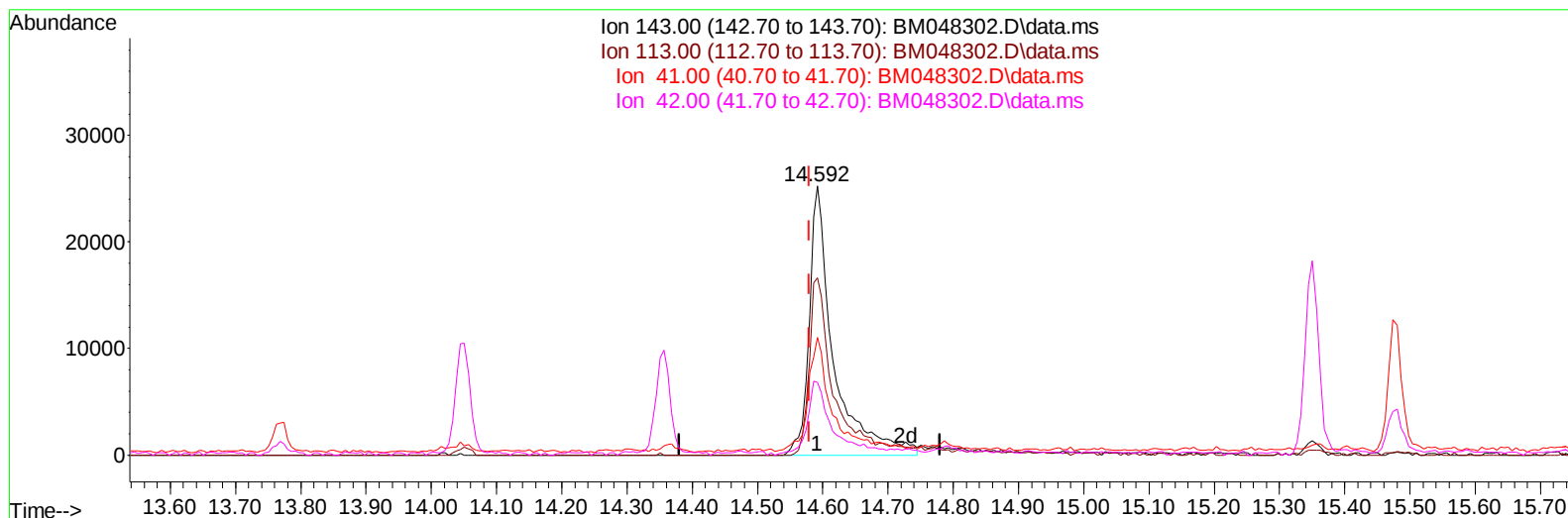
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TIC: BM048302.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.592min (+ 0.012) 8.96 ng/ul m

response 65598

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	68.30	65.82
41.00	40.60	43.67
42.00	26.40	27.04

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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 10/30/2024
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Quant Time: Oct 30 01:11:46 2024
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 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	7.716	152	219574	20.000	ng/ul	0.00
20) Naphthalene-d8	10.504	136	887212	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.357	164	535427	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.098	188	1031954	20.000	ng/ul	-0.01
79) Chrysene-d12	21.321	240	902299	20.000	ng/ul	-0.01
88) Perylene-d12	24.268	264	1005788	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.169	96	23431	4.201	ng/uL	0.00
4) Pyridine-d5	3.587	84	205429	12.708	ng/ul	0.00
7) Phenol-d5	6.892	99	249239	13.213	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.045	67	343876	29.428	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.251	132	455145	29.669	ng/ul	0.00
15) 4-Methylphenol-d8	8.428	113	352891	23.301	ng/ul	-0.01
21) Nitrobenzene-d5	8.875	128	217464	29.897	ng/ul	0.00
24) 2-Nitrophenol-d4	9.592	143	246796	31.320	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.133	165	432100	30.050	ng/ul	0.00
31) 4-Chloroaniline-d4	10.645	131	345650	17.311	ng/ul	0.00
46) Dimethylphthalate-d6	13.768	166	1316147	32.893	ng/ul	-0.01
49) Acenaphthylene-d8	14.051	160	1455460	32.261	ng/ul	0.00
54) 4-Nitrophenol-d4	14.592	143	65598m	8.965	ng/ul	0.01
60) Fluorene-d10	15.351	176	1122558	33.202	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.480	200	197178	28.913	ng/ul	0.00
73) Anthracene-d10	17.198	188	1789948	36.681	ng/ul	-0.01
81) Pyrene-d10	19.492	212	2023542	36.957	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.062	264	1984671	37.411	ng/ul	-0.02

Target Compounds Qvalue

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

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