

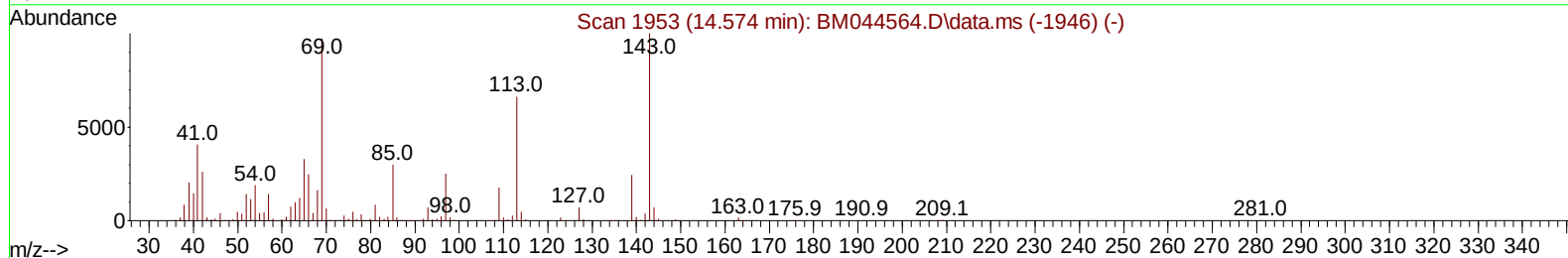
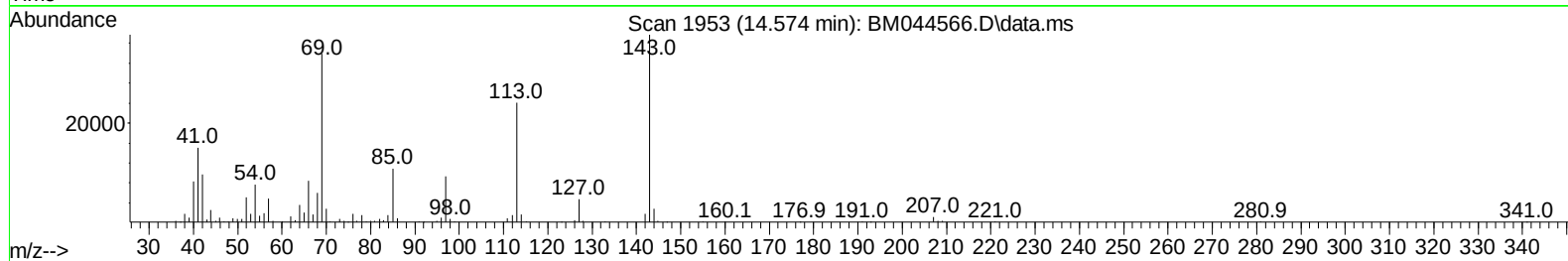
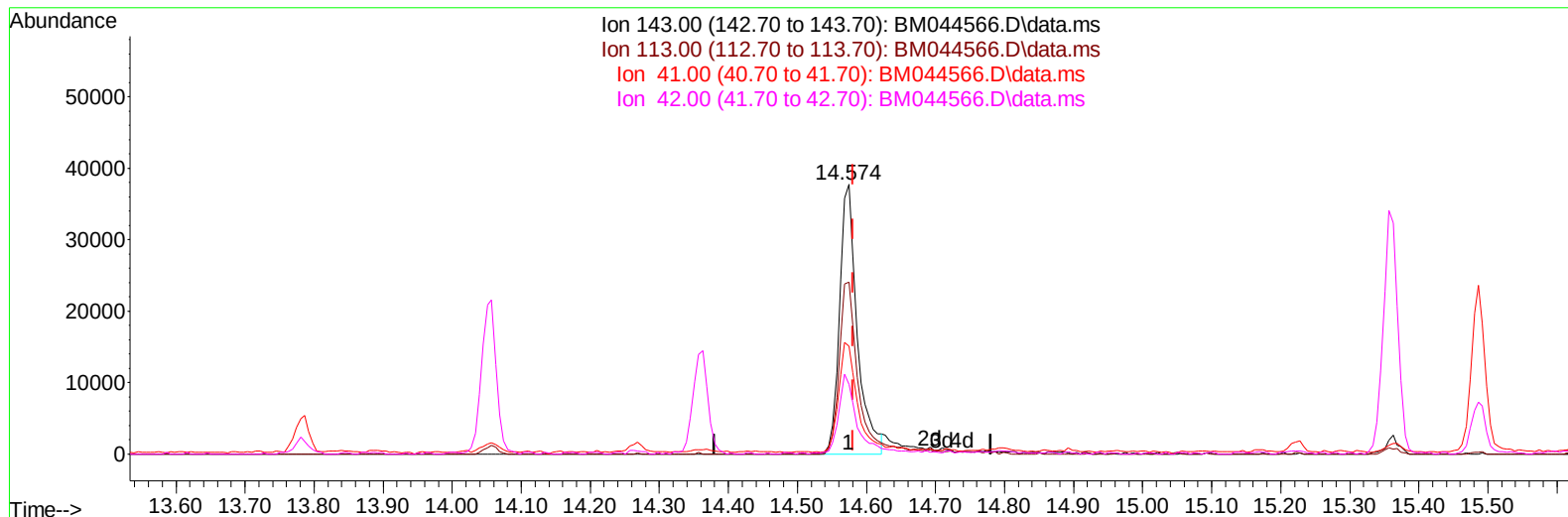
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030824\
Data File : BM044566.D
Acq On : 08 Mar 2024 15:48
Operator : MA/JU
Sample : P1703-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BH7G7

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 03/09/2024
Supervised By :mohammad ahmed 03/09/2024

Quant Time: Mar 08 16:18:32 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Mar 05 23:05:11 2024
Response via : Initial Calibration



TIC: BM044566.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.574min (-0.006) 5.82 ng/ul

response 66965

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	63.70	63.87
41.00	38.40	40.06
42.00	26.60	25.90

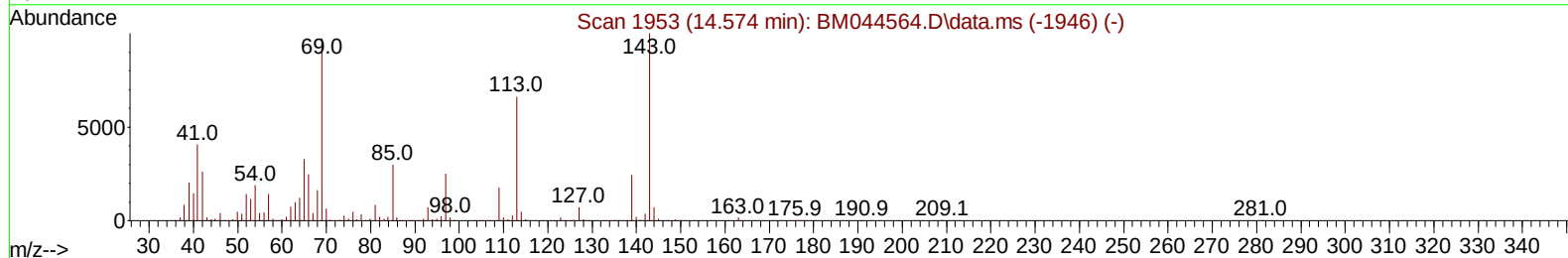
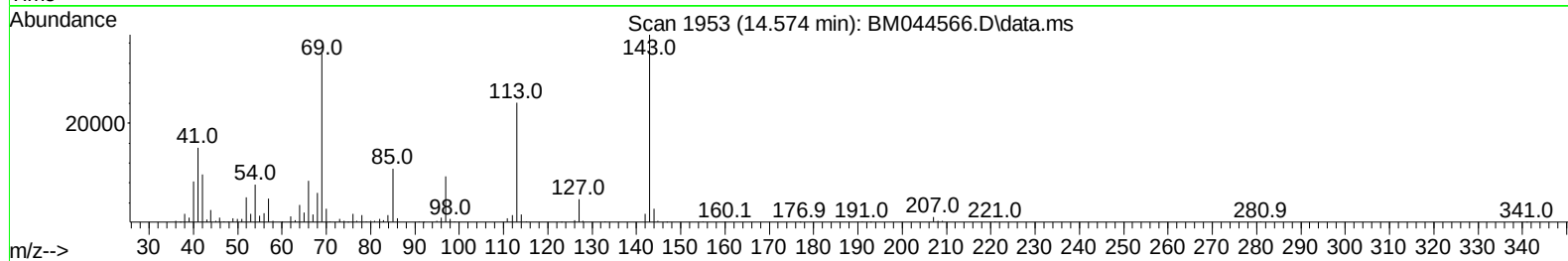
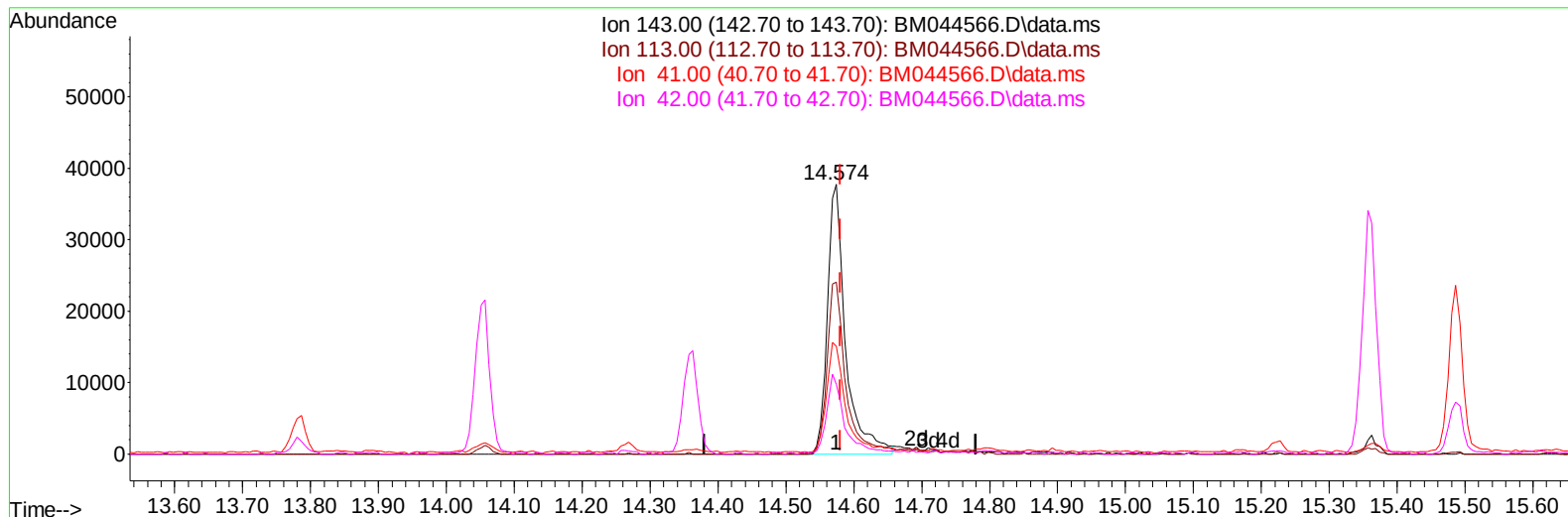
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TIC: BM044566.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.574min (-0.006) 6.11 ng/ul m

response 70402

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	63.70	63.87
41.00	38.40	40.06
42.00	26.60	25.90

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Quant Time: Mar 08 21:59:14 2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 05 23:05:11 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.716	152	242123	20.000	ng/ul	-0.02
20) Naphthalene-d8	10.498	136	1088885	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.363	164	707826	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.115	188	1438269	20.000	ng/ul	0.00
79) Chrysene-d12	21.309	240	1217439	20.000	ng/ul	0.00
88) Perylene-d12	23.562	264	1186705	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	33372	4.675	ng/uL	-0.01
4) Pyridine-d5	3.675	84	228221	10.862	ng/ul	0.00
7) Phenol-d5	6.887	99	189058	7.595	ng/ul	-0.02
9) Bis-(2-Chloroethyl)eth...	7.063	67	570233	37.086	ng/ul	-0.02
11) 2-Chlorophenol-d4	7.251	132	500322	27.350	ng/ul	-0.02
15) 4-Methylphenol-d8	8.416	113	349560	17.881	ng/ul	-0.02
21) Nitrobenzene-d5	8.881	128	357321	38.970	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.592	143	373660	37.802	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.122	165	583811	34.704	ng/ul	-0.02
31) 4-Chloroaniline-d4	10.645	131	920866	32.959	ng/ul	-0.02
46) Dimethylphthalate-d6	13.786	166	2195234	39.536	ng/ul	0.00
49) Acenaphthylene-d8	14.057	160	2499821	39.367	ng/ul	0.00
54) 4-Nitrophenol-d4	14.574	143	70402m	6.114	ng/ul	0.00
60) Fluorene-d10	15.363	176	1918336	41.113	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.486	200	337731	36.175	ng/ul	0.00
73) Anthracene-d10	17.215	188	2978831	43.330	ng/ul	0.00
81) Pyrene-d10	19.515	212	3344256	44.787	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.421	264	2706384	43.389	ng/ul	0.00

Target Compounds Qvalue

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

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