

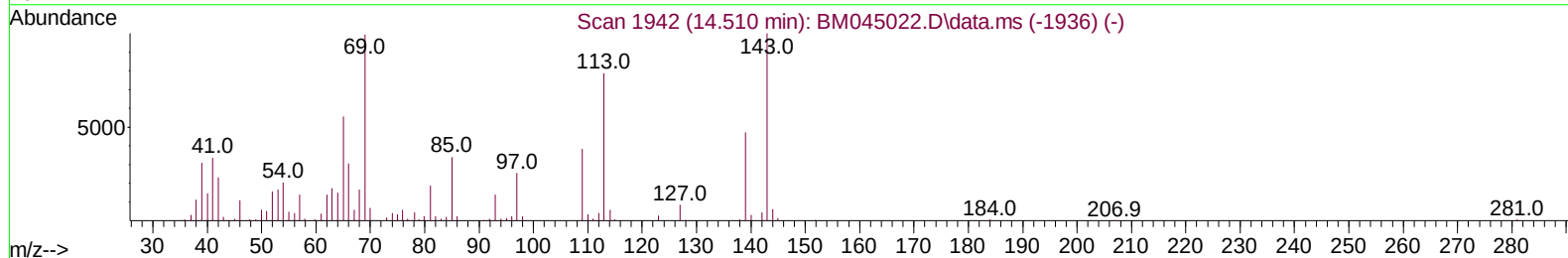
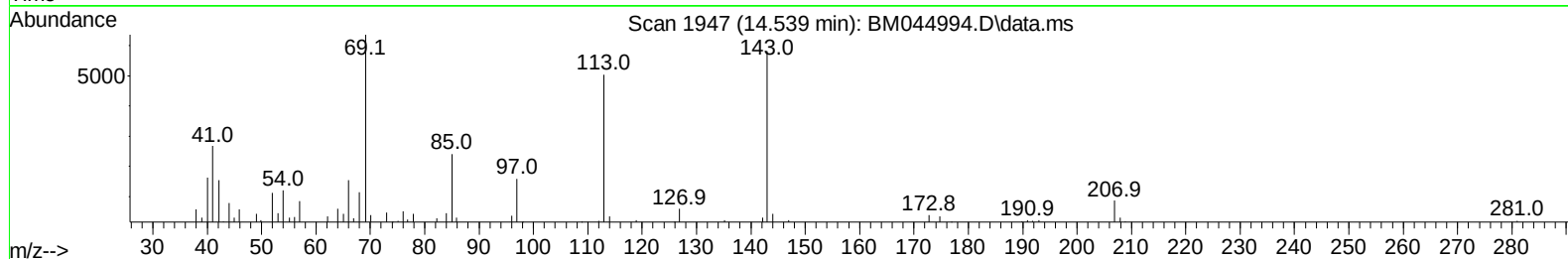
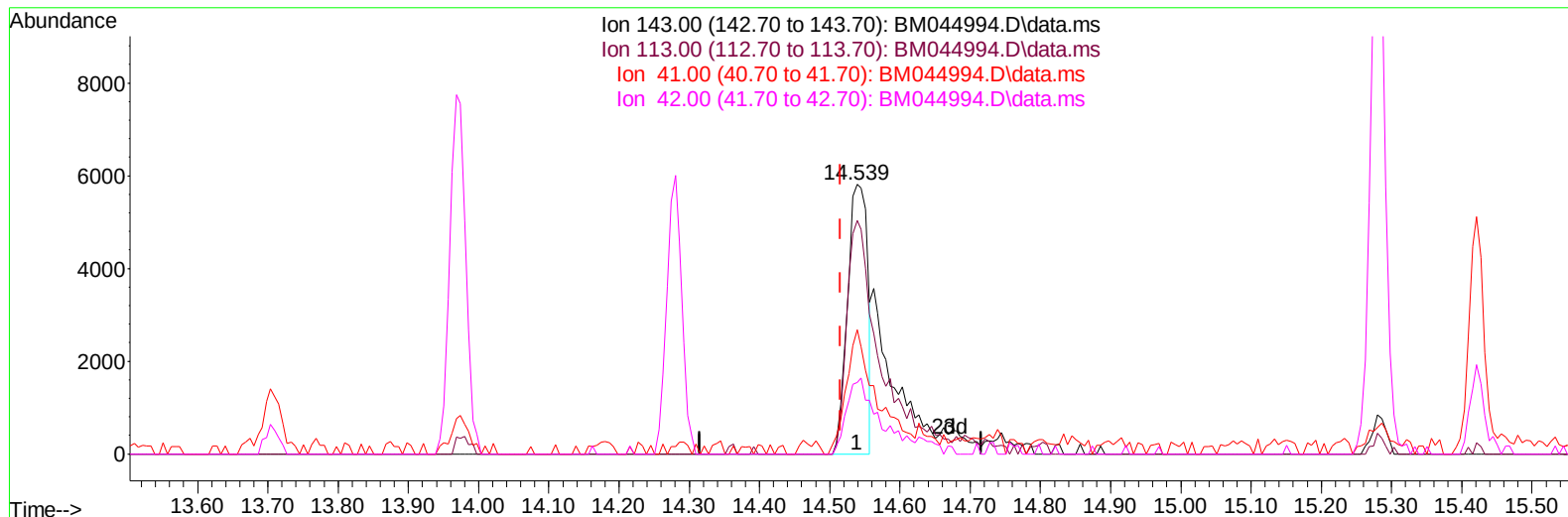
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
Data File : BM044994.D
Acq On : 07 Apr 2024 05:49
Operator : MA/JU
Sample : P2013-10
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EOMB6

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 04/09/2024
Supervised By :mohammad ahmed 04/09/2024

Quant Time: Apr 08 05:07:35 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Apr 08 05:01:54 2024
Response via : Initial Calibration



TIC: BM044994.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.539min (+ 0.023) 3.63 ng/ul

response 11618

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.80	86.75
41.00	36.50	46.05#
42.00	23.30	26.50

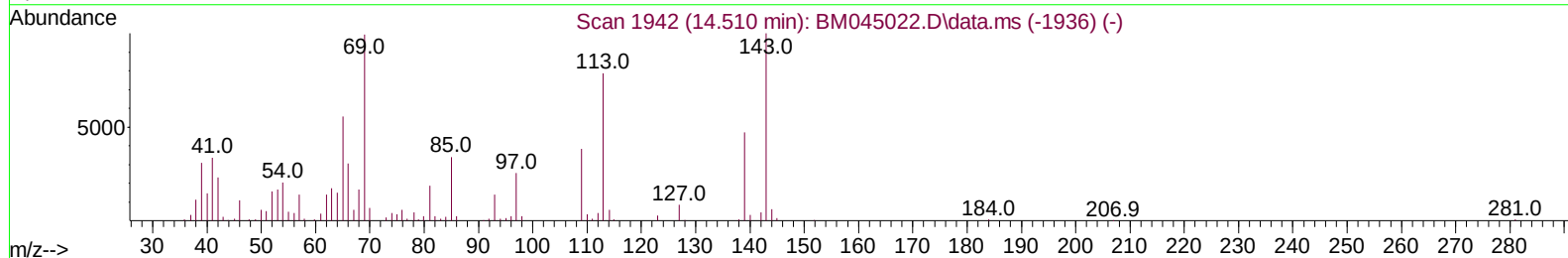
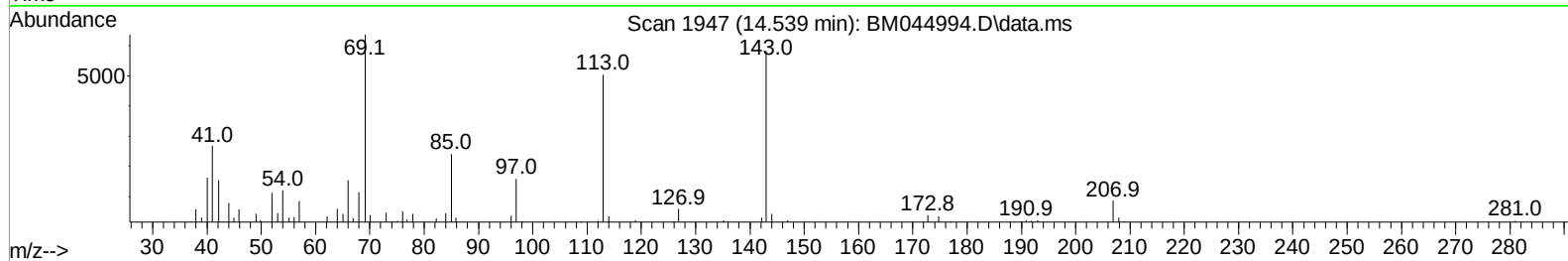
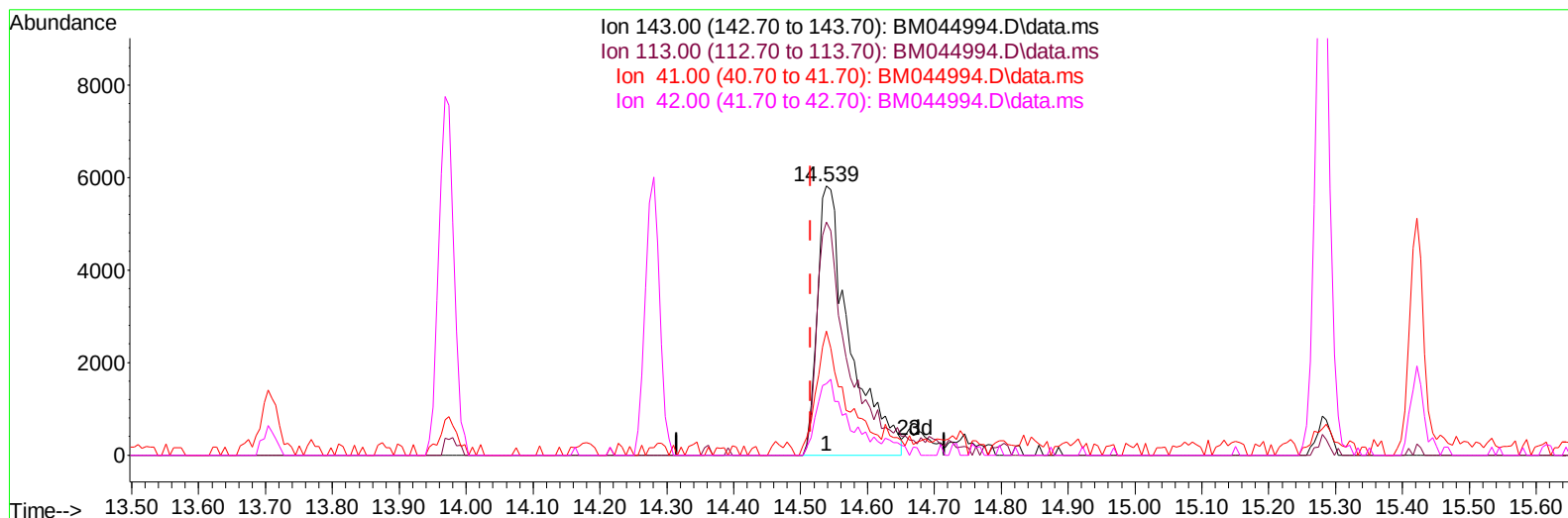
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TIC: BM044994.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.539min (+ 0.023) 6.09 ng/ul m

response 19500

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.80	86.75
41.00	36.50	46.05#
42.00	23.30	26.50

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Instrument :
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 E0MB6

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 04/09/2024
 Supervised By :mohammad ahmed 04/09/2024

Quant Time: Apr 08 05:44:22 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Apr 08 05:01:54 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.634	152	80228	20.000	ng/ul	0.00
20) Naphthalene-d8	10.404	136	350439	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.280	164	220274	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.039	188	456085	20.000	ng/ul	0.00
79) Chrysene-d12	21.244	240	429610	20.000	ng/ul	0.00
88) Perylene-d12	23.468	264	515407	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.210	96	11918	5.547	ng/uL	0.00
4) Pyridine-d5	3.616	84	61605	10.544	ng/ul	0.00
7) Phenol-d5	6.816	99	57341	8.431	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.987	67	164262	40.537	ng/ul	0.00
11) 2-Chlorophenol-d4	7.169	132	174646	32.776	ng/ul	0.00
15) 4-Methylphenol-d8	8.345	113	109002	20.475	ng/ul	0.00
21) Nitrobenzene-d5	8.792	128	111969	41.596	ng/ul	0.00
24) 2-Nitrophenol-d4	9.504	143	118268	41.626	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.039	165	180161	34.585	ng/ul	0.00
31) 4-Chloroaniline-d4	10.563	131	231355	27.896	ng/ul	0.00
46) Dimethylphthalate-d6	13.704	166	636829	41.522	ng/ul	0.00
49) Acenaphthylene-d8	13.974	160	755417	41.692	ng/ul	0.00
54) 4-Nitrophenol-d4	14.539	143	19500m	6.092	ng/ul	0.02
60) Fluorene-d10	15.280	176	576904	42.192	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.421	200	87029	32.183	ng/ul	0.00
73) Anthracene-d10	17.139	188	962581	46.916	ng/ul	0.00
81) Pyrene-d10	19.445	212	1117731	52.838	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.333	264	1217263	48.414	ng/ul	0.00

Target Compounds Qvalue

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

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