

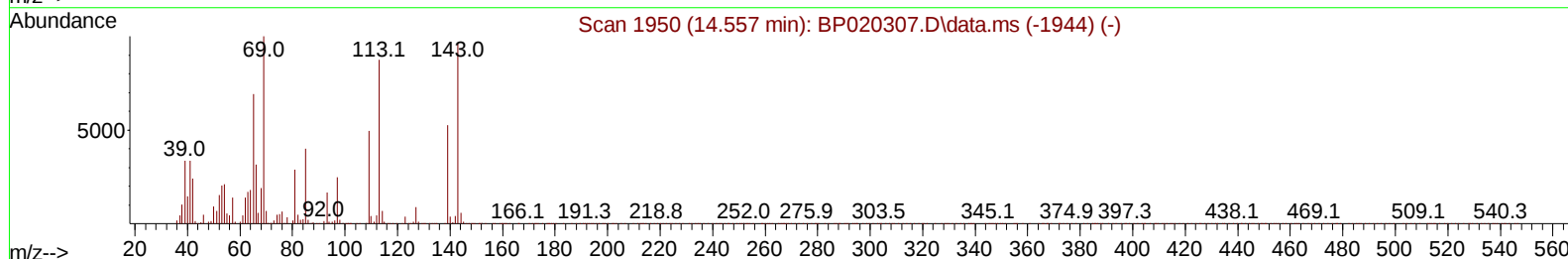
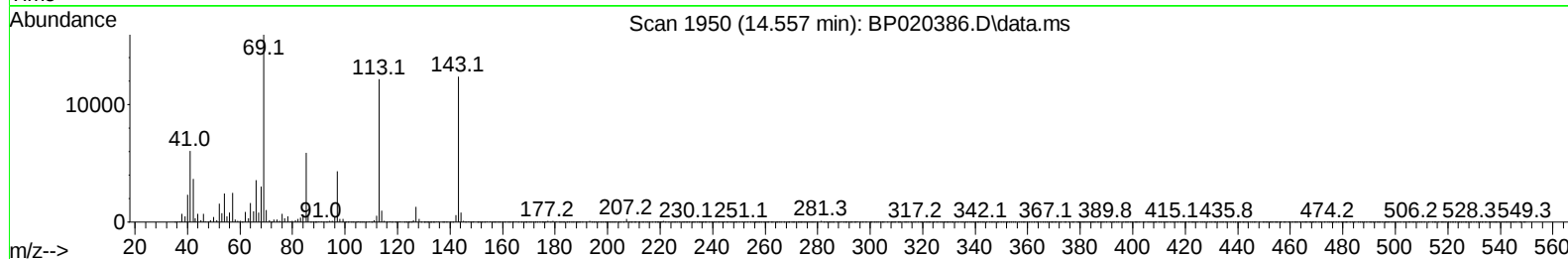
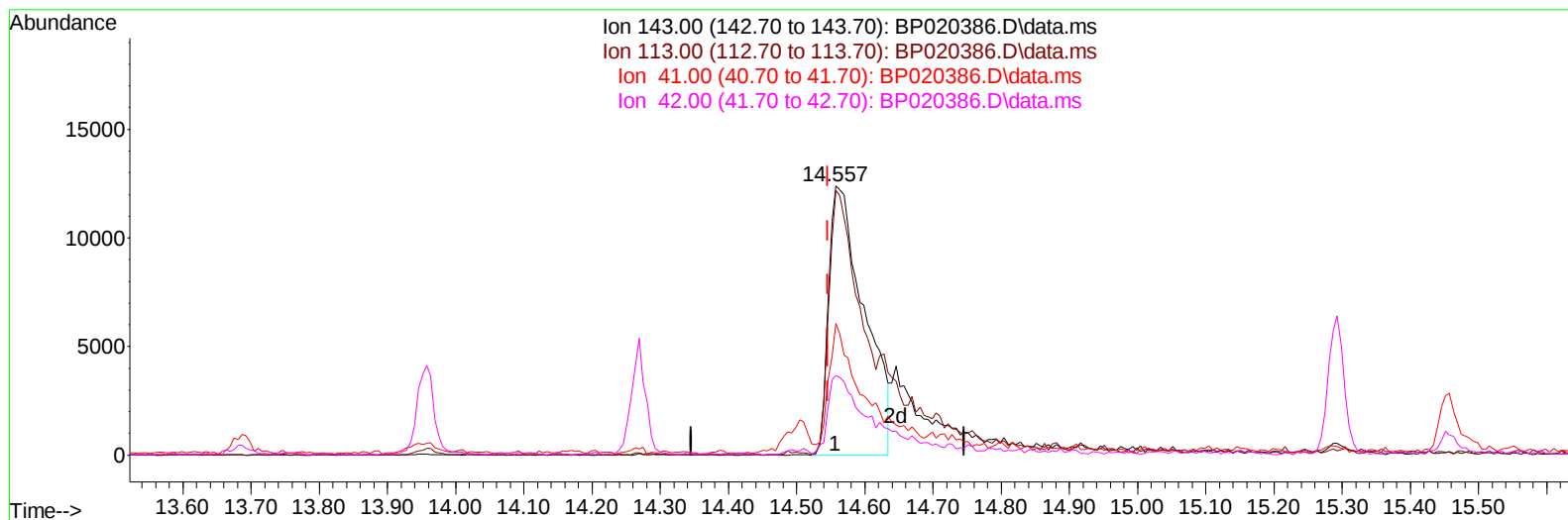
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051424\
Data File : BP020386.D
Acq On : 15 May 2024 04:35
Operator : MA/JU
Sample : P2372-16
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43W6

Manual IntegrationsAPPROVED

Quant Time: May 15 05:02:30 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 14 23:58:44 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/15/2024
Supervised By :mohammad ahmed 05/17/2024



TIC: BP020386.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.557min (+ 0.012) 17.32 ng/u1

response 45023

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	98.39
41.00	50.30	49.00
42.00	32.50	29.50

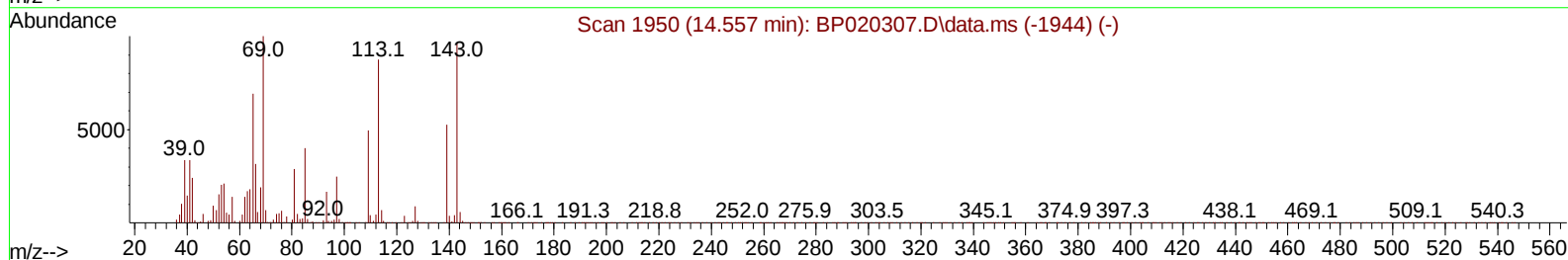
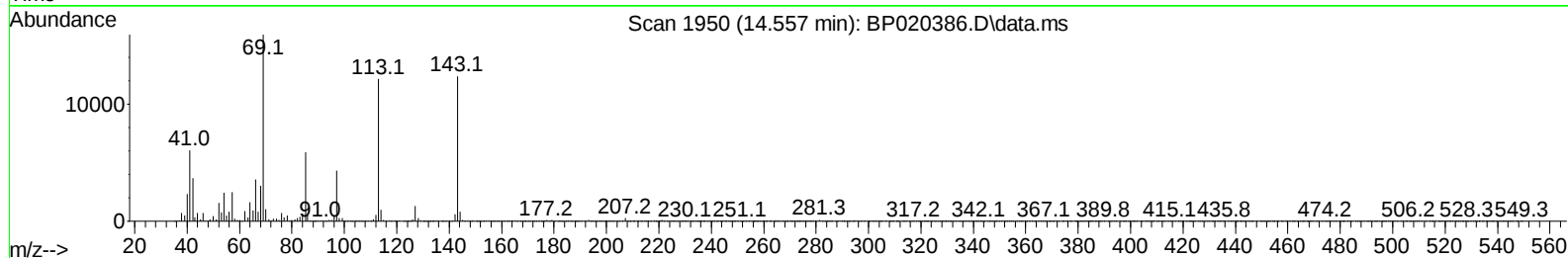
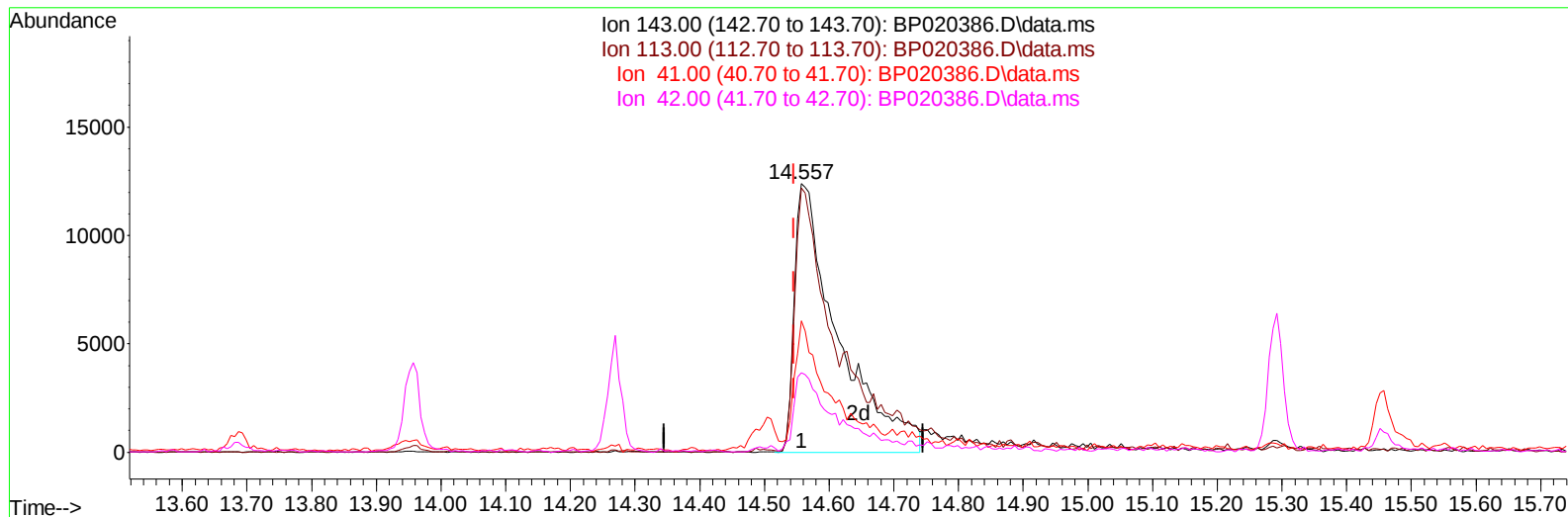
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TIC: BP020386.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.557min (+ 0.012) 22.28 ng/ul m

response 57937

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	98.39
41.00	50.30	49.00
42.00	32.50	29.50

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 Sample : P2372-16
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A43W6

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/15/2024
 Supervised By :mohammad ahmed 05/17/2024

Quant Time: May 15 05:05:52 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 14 23:58:44 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.593	152	77101	20.000	ng/ul	0.00
20) Naphthalene-d8	10.357	136	312096	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.269	164	231661	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.104	188	523897	20.000	ng/ul	0.00
79) Chrysene-d12	21.604	240	526418	20.000	ng/ul	0.00
88) Perylene-d12	24.957	264	586506	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.193	96	7516	4.117	ng/uL	0.00
4) Pyridine-d5	3.599	84	97722	18.422	ng/ul	0.00
7) Phenol-d5	6.787	99	143003	21.597	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.946	67	90040	22.385	ng/ul	0.00
11) 2-Chlorophenol-d4	7.128	132	122177	25.811	ng/ul	0.00
15) 4-Methylphenol-d8	8.305	113	118658	22.166	ng/ul	0.00
21) Nitrobenzene-d5	8.752	128	60338	25.818	ng/ul	0.00
24) 2-Nitrophenol-d4	9.469	143	68111	25.450	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.010	165	118668	24.298	ng/ul	0.00
31) 4-Chloroaniline-d4	10.534	131	161780	23.668	ng/ul	0.00
46) Dimethylphthalate-d6	13.687	166	476761	28.182	ng/ul	0.00
49) Acenaphthylene-d8	13.957	160	504673	27.005	ng/ul	0.00
54) 4-Nitrophenol-d4	14.557	143	57937m	22.283	ng/ul	0.01
60) Fluorene-d10	15.293	176	402515	28.526	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.451	200	64339	19.353	ng/ul	0.00
73) Anthracene-d10	17.204	188	635580	28.043	ng/ul	-0.01
81) Pyrene-d10	19.610	212	803916	26.125	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.727	264	786192	27.724	ng/ul	0.01

Target Compounds Qvalue

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

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