

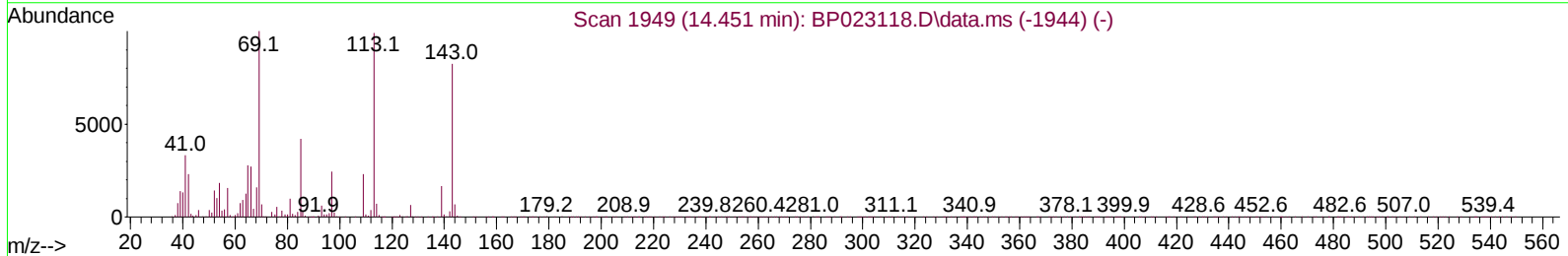
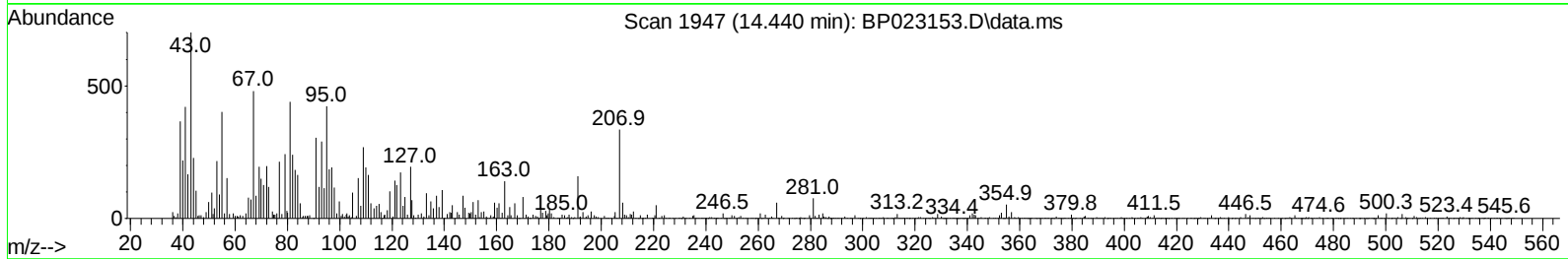
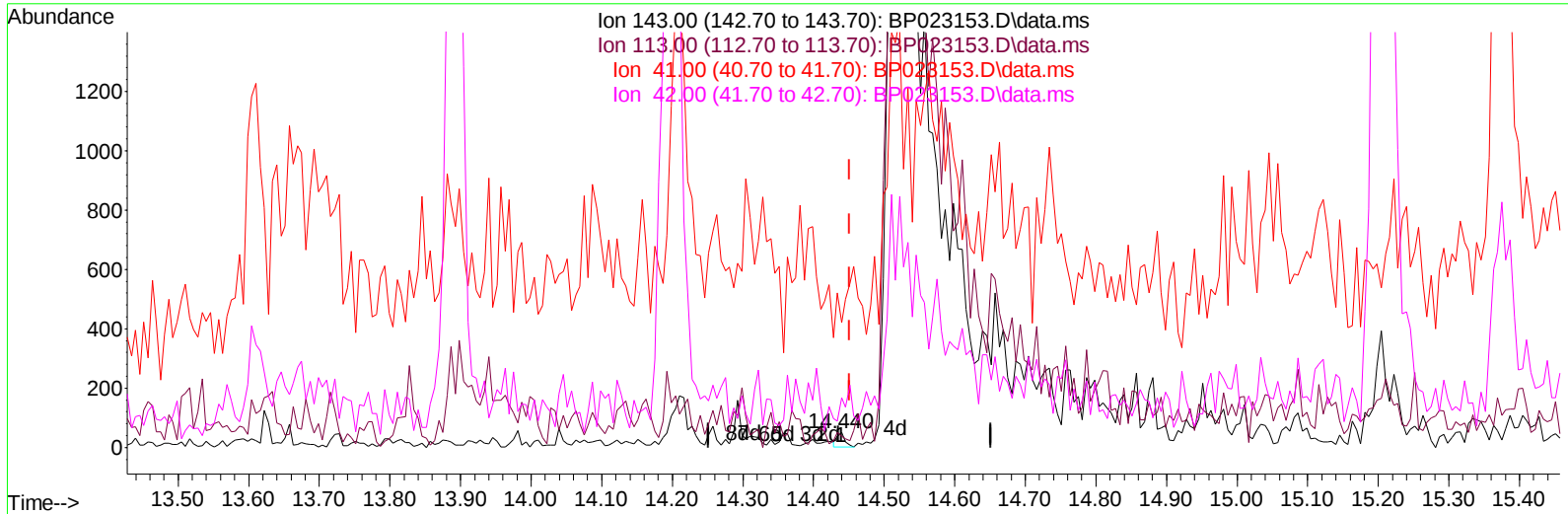
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111324\
Data File : BP023153.D
Acq On : 16 Nov 2024 00:57
Operator : RC/JU
Sample : P4803-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A0AP6

Manual IntegrationsAPPROVED

Quant Time: Nov 16 01:35:53 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 13 04:49:01 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/18/2024
Supervised By :mohammad ahmed 11/18/2024



TIC: BP023153.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.440min (-0.012) 0.03 ng/u1

response 38

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	76.00#
41.00	39.20	842.00#
42.00	25.90	334.00#

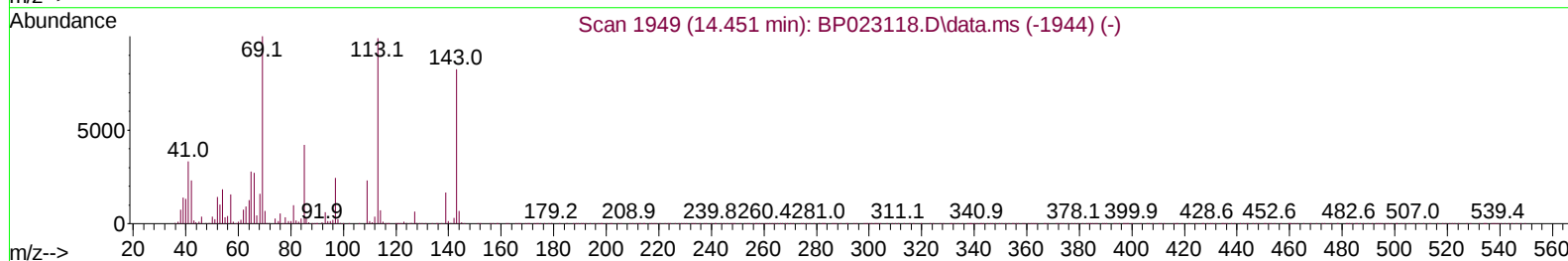
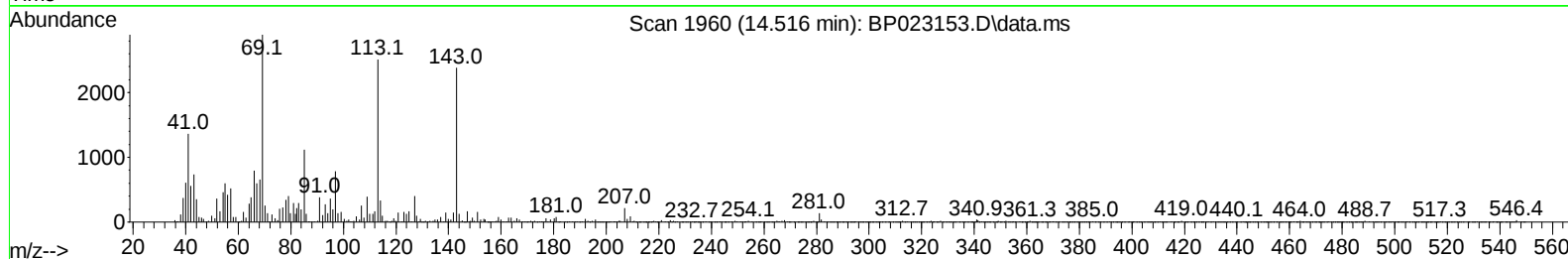
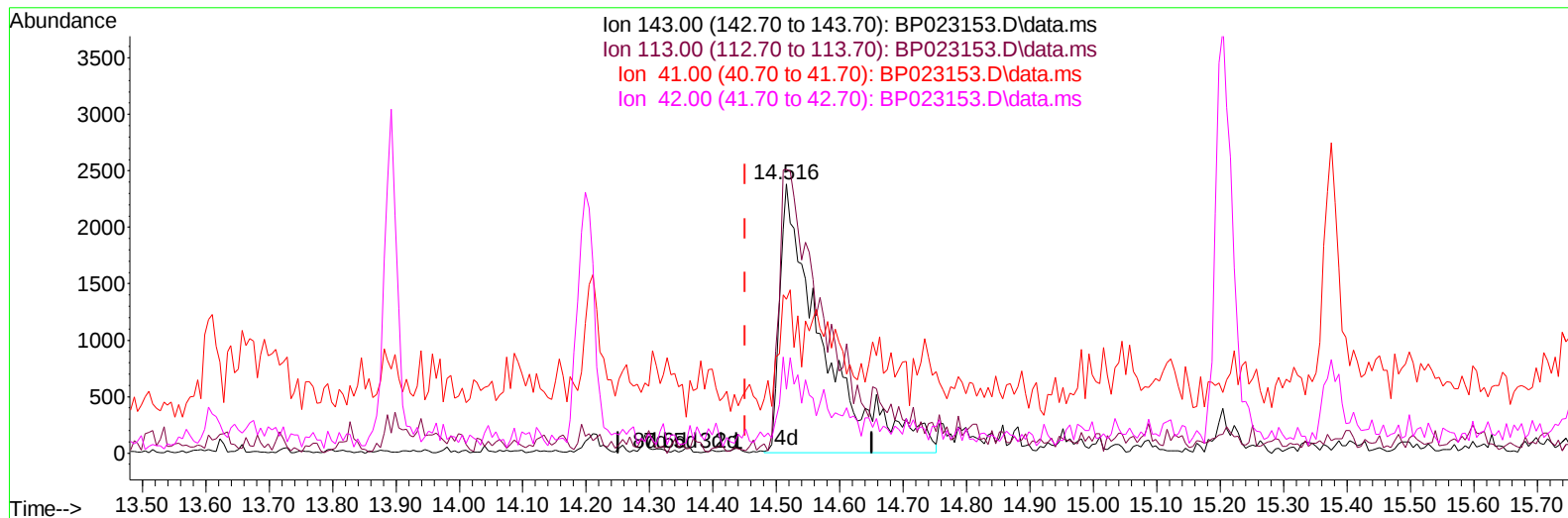
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TIC: BP023153.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.516min (+ 0.065) 9.34 ng/ul m

response 11235

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	105.46
41.00	39.20	57.26#
42.00	25.90	23.64

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 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A0AP6

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/18/2024
 Supervised By :mohammad ahmed 11/18/2024

Quant Time: Nov 16 01:37:06 2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 13 04:49:01 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.581	152	36776	20.000	ng/ul	0.00
20) Naphthalene-d8	10.328	136	141454	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.198	164	111076	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.004	188	301450	20.000	ng/ul	-0.01
79) Chrysene-d12	21.451	240	418454	20.000	ng/ul	0.00
88) Perylene-d12	24.669	264	493648	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.134	96	4575	5.980	ng/uL	0.00
4) Pyridine-d5	3.552	84	46156	20.305	ng/ul	0.00
7) Phenol-d5	6.787	99	35549	13.075	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.928	67	48514	30.370	ng/ul	0.00
11) 2-Chlorophenol-d4	7.122	132	58366	29.311	ng/ul	0.00
15) 4-Methylphenol-d8	8.293	113	60468	27.025	ng/ul	0.00
21) Nitrobenzene-d5	8.722	128	30385	30.855	ng/ul	0.00
24) 2-Nitrophenol-d4	9.434	143	36306	30.925	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.981	165	79637	32.391	ng/ul	0.01
31) 4-Chloroaniline-d4	10.481	131	83303	28.401	ng/ul	0.00
46) Dimethylphthalate-d6	13.610	166	286858	35.832	ng/ul	0.00
49) Acenaphthylene-d8	13.893	160	271078	32.452	ng/ul	0.00
54) 4-Nitrophenol-d4	14.516	143	11235m	9.338	ng/ul	0.06
60) Fluorene-d10	15.204	176	251739	35.866	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.375	200	46063	25.508	ng/ul	0.01
73) Anthracene-d10	17.110	188	496508	38.900	ng/ul	0.00
81) Pyrene-d10	19.492	212	725294	37.630	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.445	264	934476	39.741	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.164	88	3555	4.021	ng/uL	90

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

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