

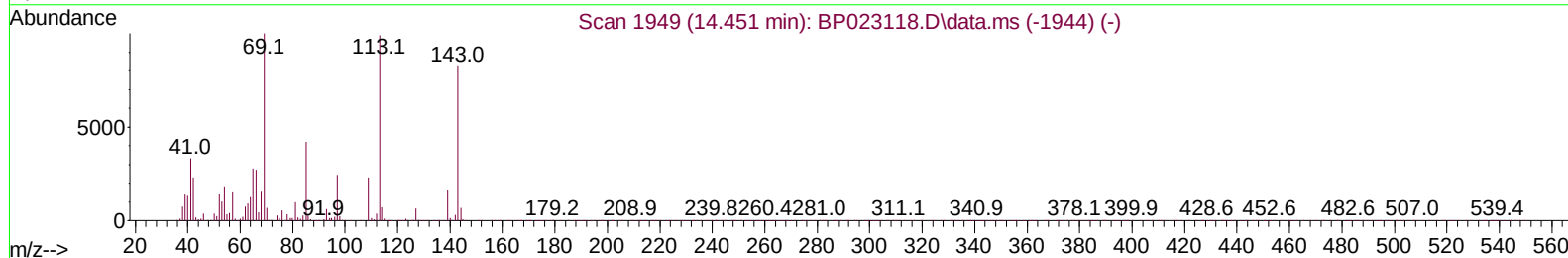
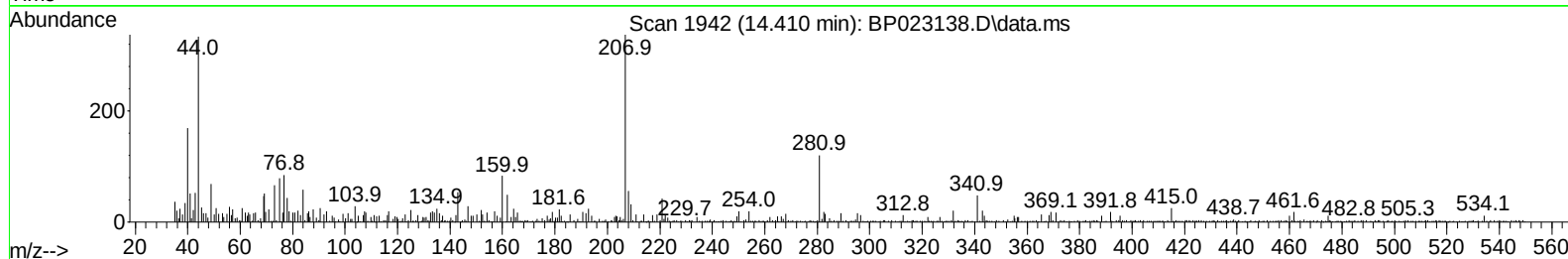
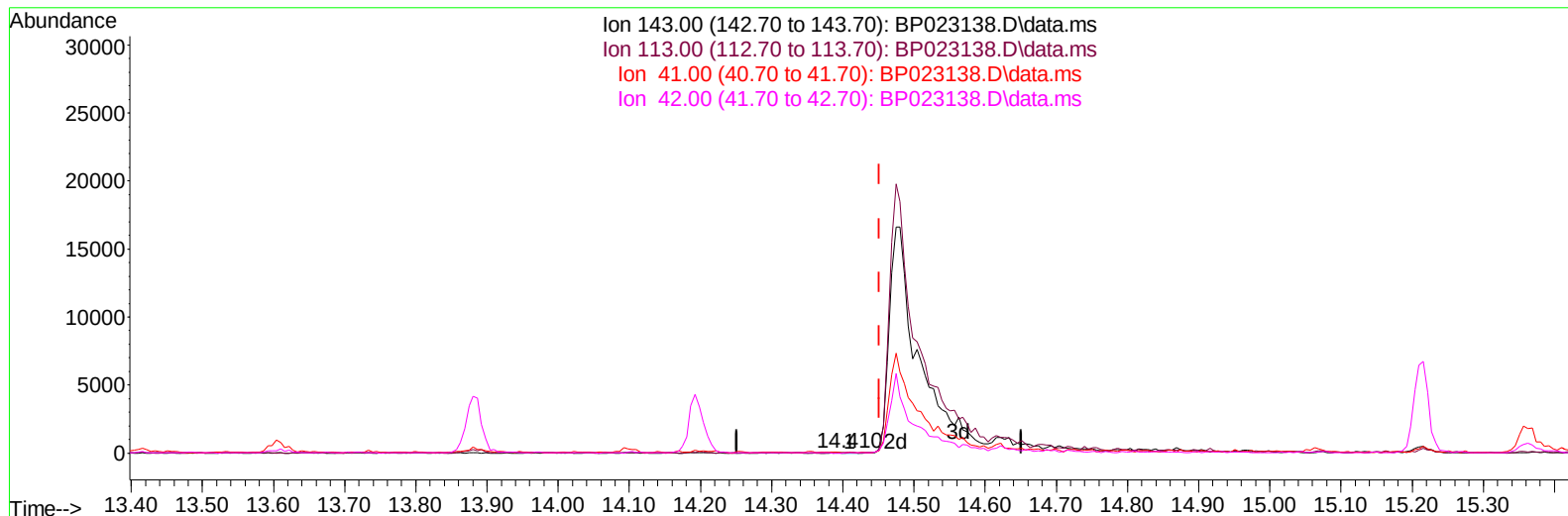
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111324\
Data File : BP023138.D
Acq On : 15 Nov 2024 14:03
Operator : RC/JU
Sample : P4724-24
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH9E4

Manual IntegrationsAPPROVED

Quant Time: Nov 15 22:52:53 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 11/16/2024
Supervised By :mohammad ahmed 11/18/2024



TIC: BP023138.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.410min (-0.042) 0.02 ng/u1

response 35

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	20.37#
41.00	39.20	107.41#
42.00	25.90	40.74#

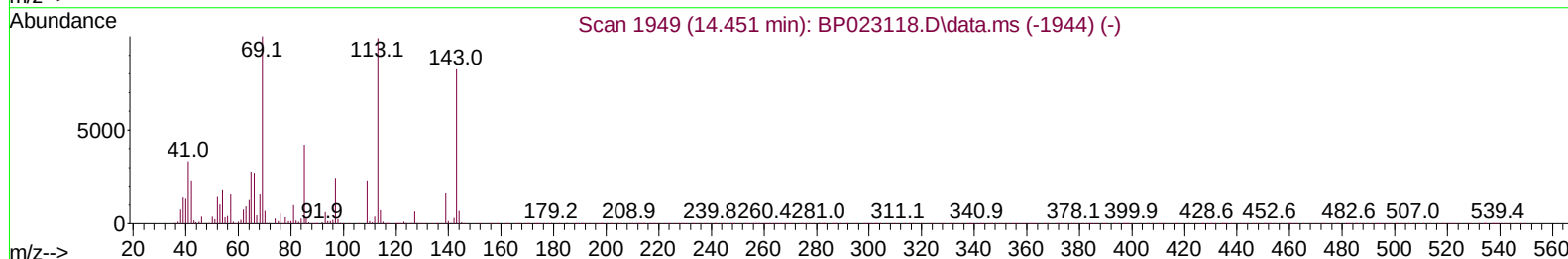
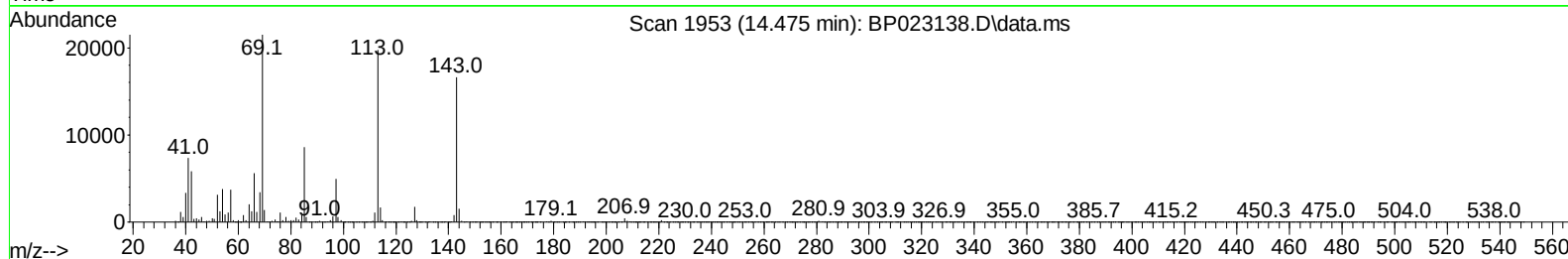
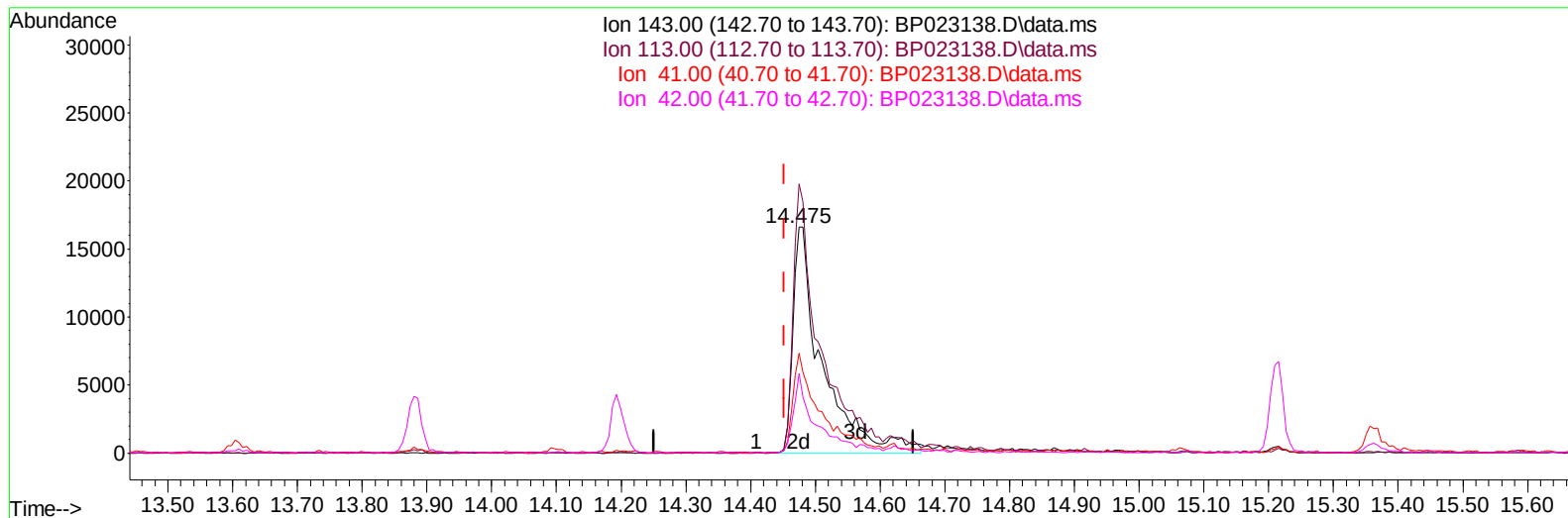
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TIC: BP023138.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.475min (+ 0.023) 25.05 ng/ul m

response 52025

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	119.03
41.00	39.20	44.22
42.00	25.90	35.21#

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 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
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 BH9E4

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/16/2024
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Quant Time: Nov 15 22:53:56 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

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	62905	20.000	ng/ul	0.00
20) Naphthalene-d8	10.334	136	237205	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.192	164	191734	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.010	188	491737	20.000	ng/ul	0.00
79) Chrysene-d12	21.445	240	598946	20.000	ng/ul	0.00
88) Perylene-d12	24.668	264	734564	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	6443	4.923	ng/uL	0.00
4) Pyridine-d5	3.557	84	45945	11.817	ng/ul	0.01
7) Phenol-d5	6.781	99	125175	26.915	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.928	67	78123	28.591	ng/ul	0.00
11) 2-Chlorophenol-d4	7.122	132	96737	28.401	ng/ul	0.00
15) 4-Methylphenol-d8	8.287	113	101965	26.643	ng/ul	0.00
21) Nitrobenzene-d5	8.722	128	47568	28.805	ng/ul	0.00
24) 2-Nitrophenol-d4	9.439	143	57893	29.407	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.981	165	115402	27.991	ng/ul	0.01
31) 4-Chloroaniline-d4	10.481	131	139646	28.392	ng/ul	0.00
46) Dimethylphthalate-d6	13.604	166	439344	31.793	ng/ul	0.00
49) Acenaphthylene-d8	13.880	160	459546	31.871	ng/ul	-0.02
54) 4-Nitrophenol-d4	14.475	143	52025m	25.052	ng/ul	0.02
60) Fluorene-d10	15.216	176	395938	32.680	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.363	200	59150	20.080	ng/ul	0.00
73) Anthracene-d10	17.104	188	683804	32.843	ng/ul	-0.01
81) Pyrene-d10	19.498	212	890217	32.268	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.450	264	1138886	32.549	ng/ul	0.00

Target Compounds Qvalue

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

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