

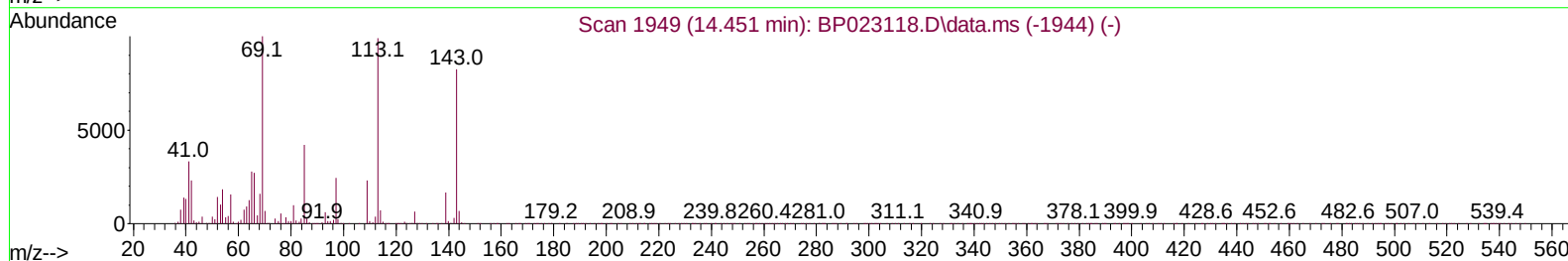
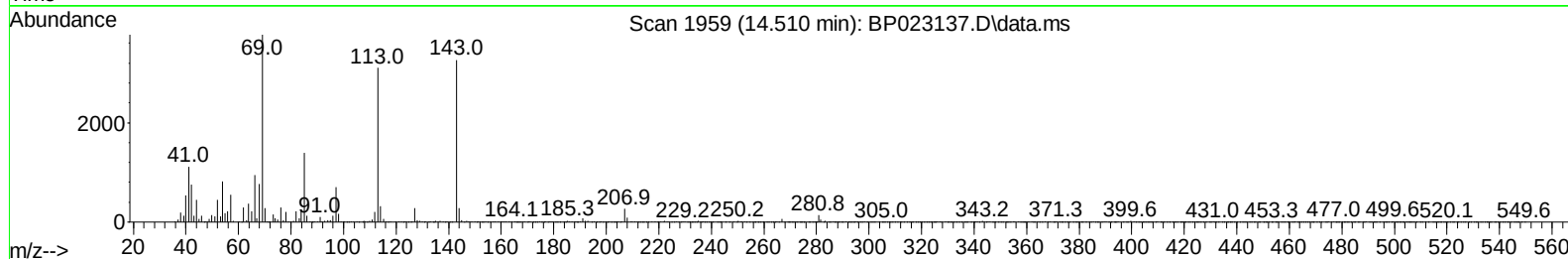
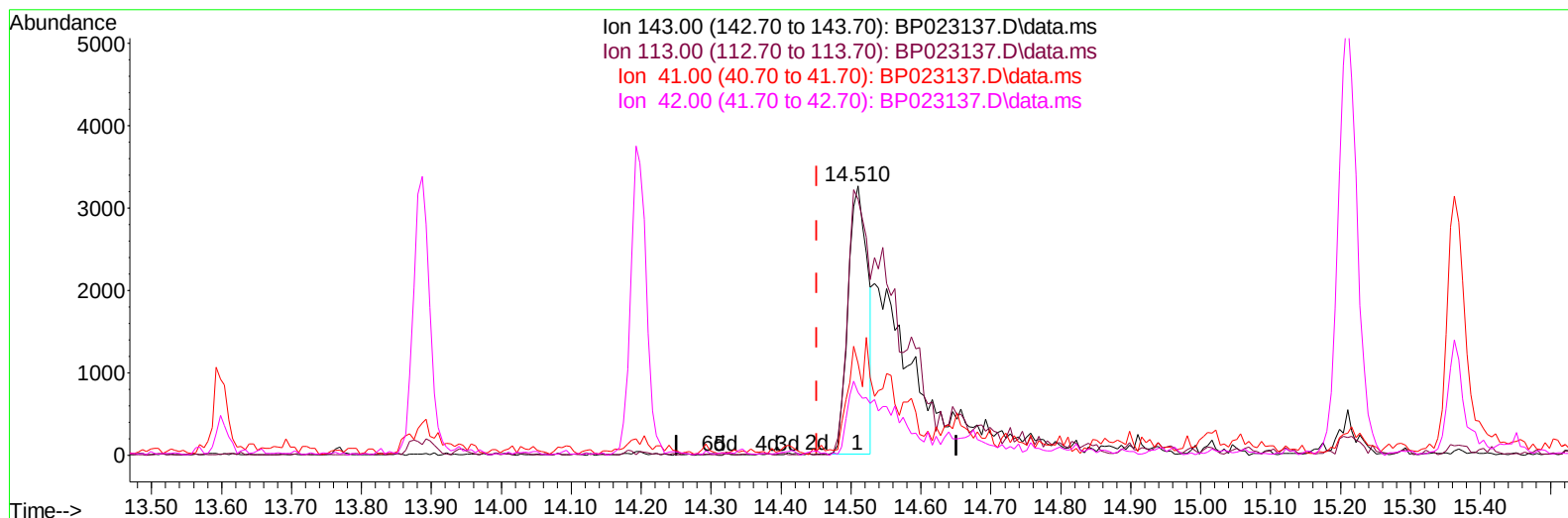
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111324\
Data File : BP023137.D
Acq On : 15 Nov 2024 13:23
Operator : RC/JU
Sample : P4740-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0CC3

Manual IntegrationsAPPROVED

Quant Time: Nov 15 22:49:41 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: BP023137.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.510min (+ 0.059) 3.16 ng/ul

response 6319

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	95.04
41.00	39.20	34.14
42.00	25.90	23.12

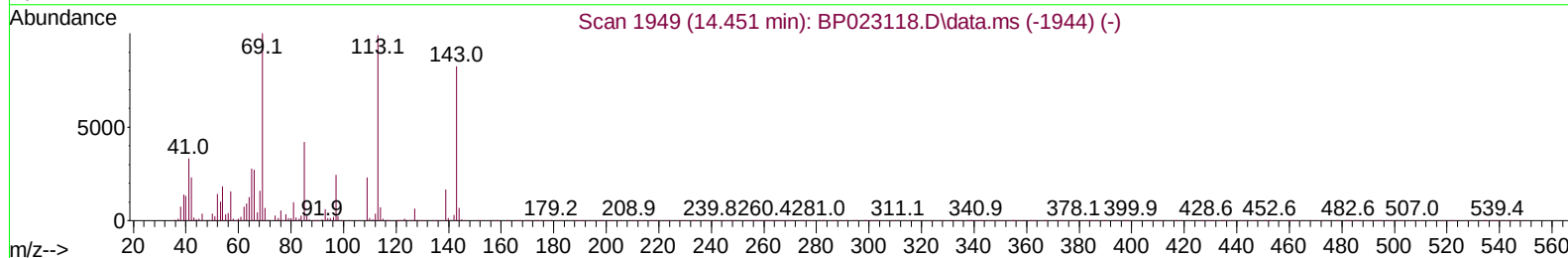
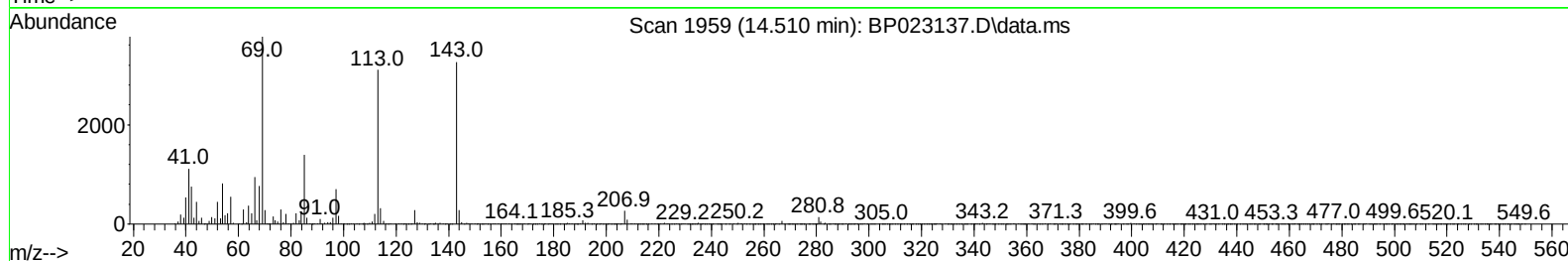
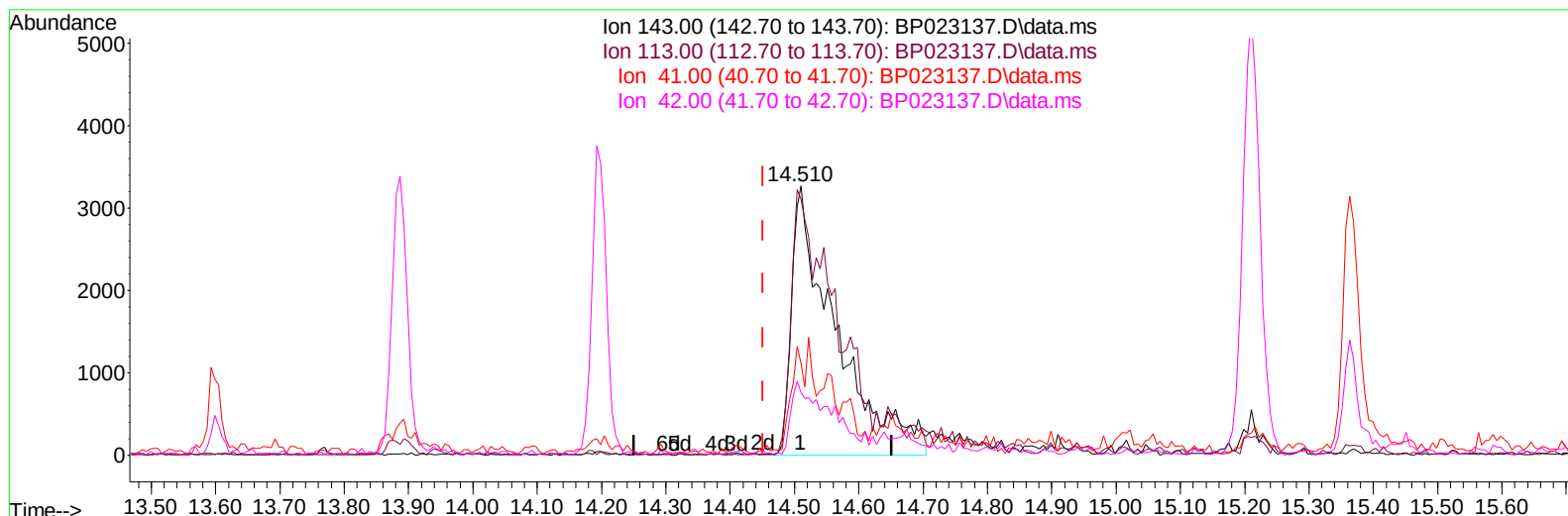
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TIC: BP023137.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.510min (+ 0.059) 7.72 ng/ul m

response 15450

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	95.04
41.00	39.20	34.14
42.00	25.90	23.12

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/16/2024
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Quant Time: Nov 15 22:51:28 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	60717	20.000	ng/ul	0.00
20) Naphthalene-d8	10.328	136	232079	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.198	164	184679	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.004	188	449123	20.000	ng/ul	-0.01
79) Chrysene-d12	21.445	240	545451	20.000	ng/ul	0.00
88) Perylene-d12	24.662	264	647157	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.134	96	6169	4.884	ng/uL	0.00
4) Pyridine-d5	3.564	84	19024	5.069	ng/ul	0.02
7) Phenol-d5	6.787	99	31972	7.122	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.928	67	72312	27.418	ng/ul	0.00
11) 2-Chlorophenol-d4	7.128	132	82470	25.085	ng/ul	0.00
15) 4-Methylphenol-d8	8.293	113	48982	13.260	ng/ul	0.00
21) Nitrobenzene-d5	8.722	128	45576	28.209	ng/ul	0.00
24) 2-Nitrophenol-d4	9.434	143	54335	28.209	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.981	165	111338	27.602	ng/ul	0.01
31) 4-Chloroaniline-d4	10.487	131	77946	16.197	ng/ul	0.00
46) Dimethylphthalate-d6	13.598	166	448334	33.683	ng/ul	-0.01
49) Acenaphthylene-d8	13.887	160	422125	30.394	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.510	143	15450m	7.724	ng/ul	0.06
60) Fluorene-d10	15.210	176	393324	33.704	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.363	200	84599	31.444	ng/ul	0.00
73) Anthracene-d10	17.110	188	745007	39.177	ng/ul	0.00
81) Pyrene-d10	19.492	212	974780	38.799	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.445	264	1145468	37.158	ng/ul	0.00

Target Compounds Qvalue

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

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