

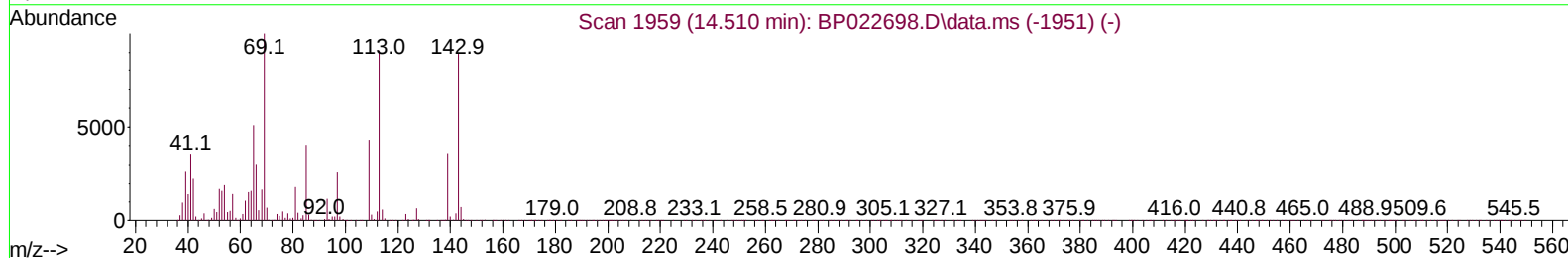
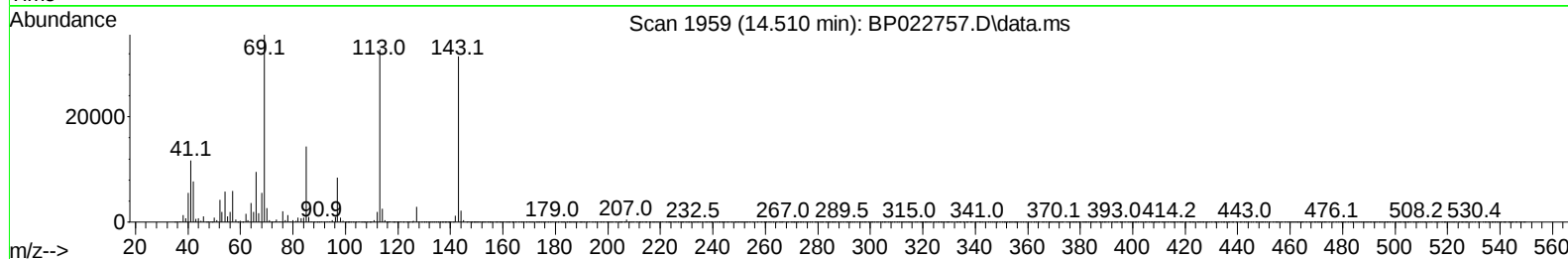
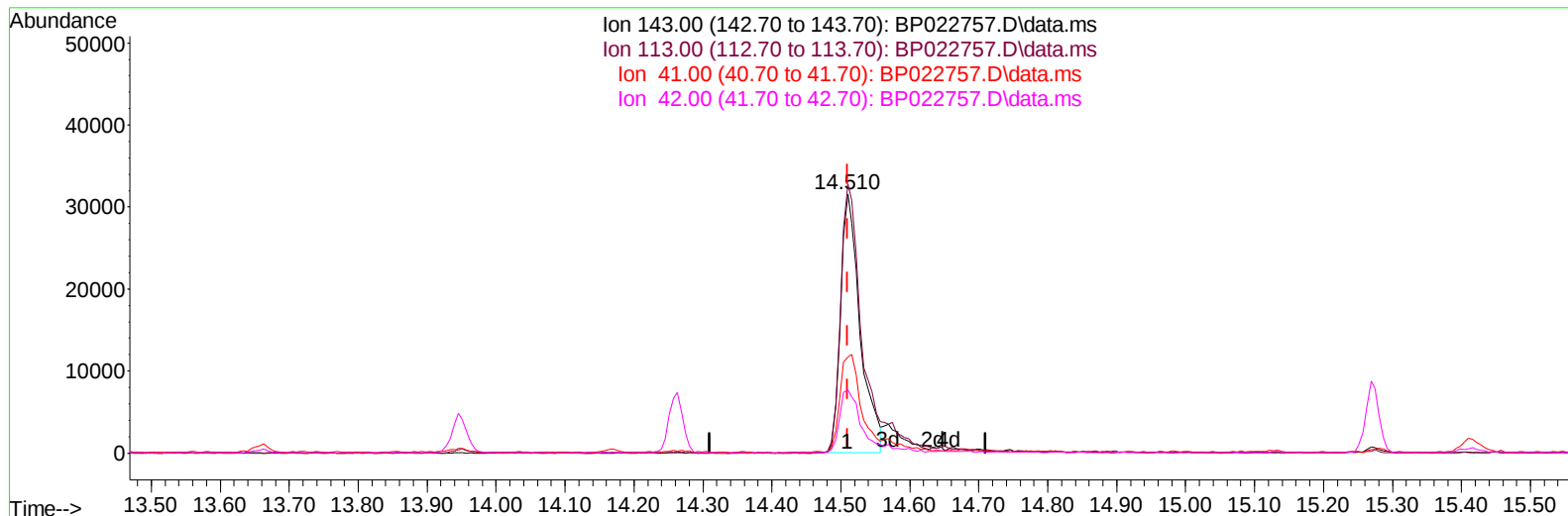
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103024\
Data File : BP022757.D
Acq On : 31 Oct 2024 06:00
Operator : RC/JU
Sample : P4592-03
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
F7H03

Manual IntegrationsAPPROVED

Quant Time: Oct 31 12:39:29 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Oct 30 10:56:39 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/04/2024
Supervised By :mohammad ahmed 11/05/2024



TIC: BP022757.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.510min (-0.000) 12.74 ng/u1

response 61835

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	103.93
41.00	35.90	37.13
42.00	23.70	24.41

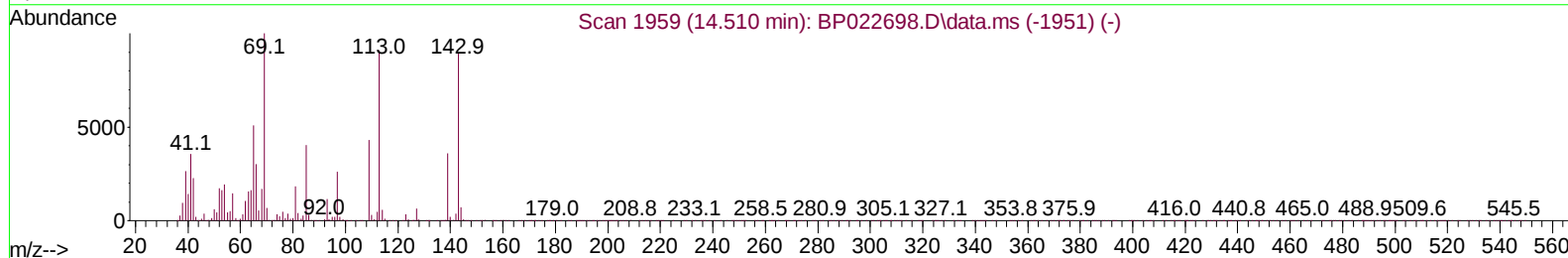
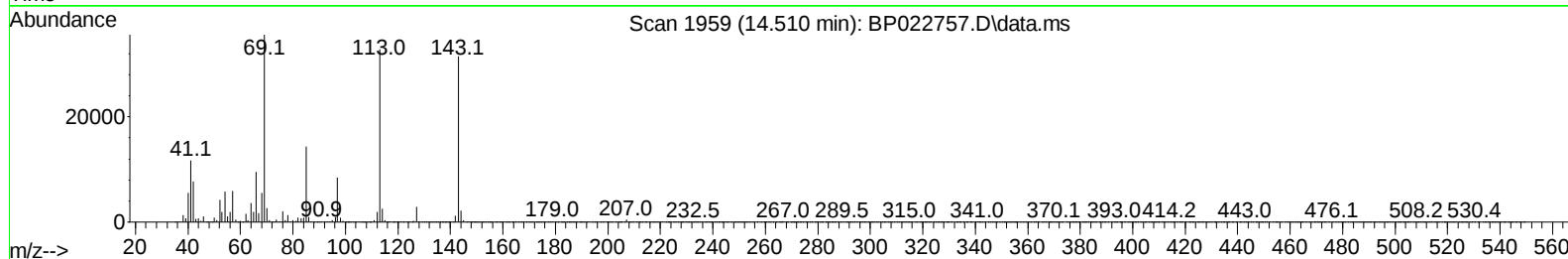
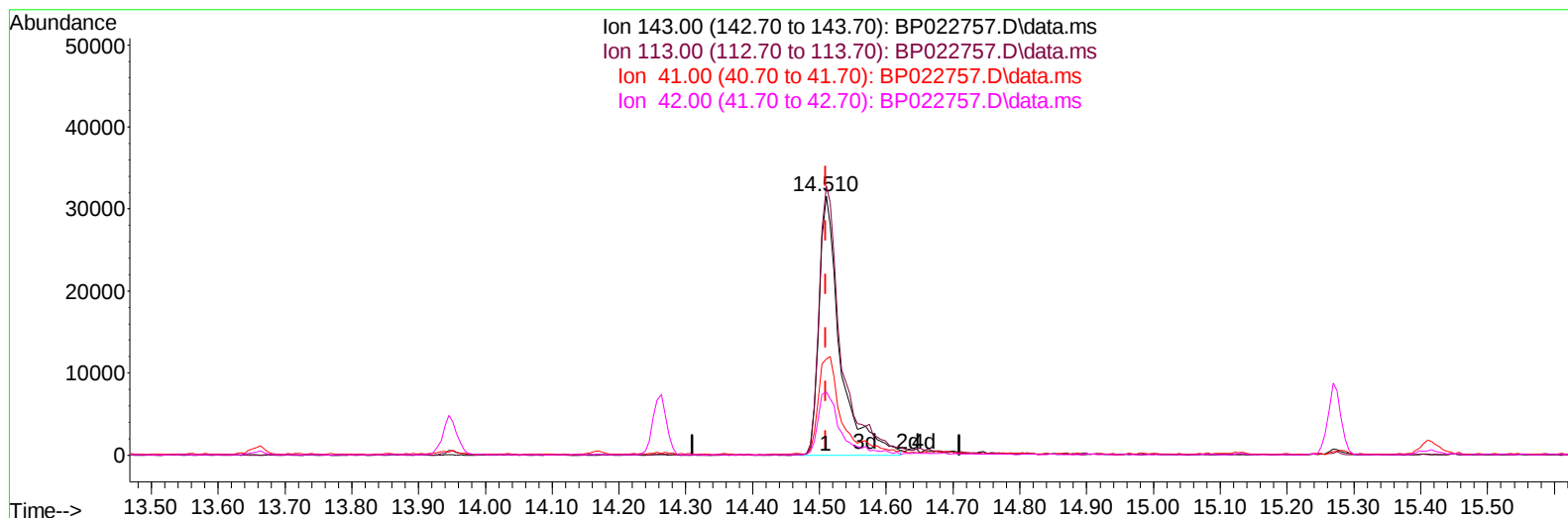
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TIC: BP022757.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.510min (-0.000) 14.28 ng/ul m

response 69267

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	103.93
41.00	35.90	37.13
42.00	23.70	24.41

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/04/2024
 Supervised By :mohammad ahmed 11/05/2024

Quant Time: Nov 01 01:19:08 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 30 10:56:39 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.634	152	117466	20.000	ng/ul	0.00
20) Naphthalene-d8	10.393	136	453663	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.263	164	363996	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.063	188	899099	20.000	ng/ul	0.00
79) Chrysene-d12	21.492	240	954769	20.000	ng/ul	0.00
88) Perylene-d12	24.751	264	1097586	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	7296	2.414	ng/uL	0.00
4) Pyridine-d5	3.593	84	102113	12.305	ng/ul	0.00
7) Phenol-d5	6.834	99	161568	16.340	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.981	67	92792	15.967	ng/ul	0.00
11) 2-Chlorophenol-d4	7.181	132	118927	16.950	ng/ul	0.00
15) 4-Methylphenol-d8	8.340	113	138764	17.585	ng/ul	0.00
21) Nitrobenzene-d5	8.781	128	61391	17.600	ng/ul	0.00
24) 2-Nitrophenol-d4	9.487	143	77103	19.092	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.028	165	154311	17.978	ng/ul	0.00
31) 4-Chloroaniline-d4	10.534	131	171922	16.951	ng/ul	0.00
46) Dimethylphthalate-d6	13.657	166	561319	19.409	ng/ul	0.00
49) Acenaphthylene-d8	13.945	160	577053	18.786	ng/ul	0.00
54) 4-Nitrophenol-d4	14.510	143	69267m	14.277	ng/ul	0.00
60) Fluorene-d10	15.275	176	498222	19.862	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.410	200	56966	9.634	ng/ul	0.00
73) Anthracene-d10	17.163	188	857858	20.282	ng/ul	0.00
81) Pyrene-d10	19.551	212	1068841	21.169	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.533	264	1252428	21.366	ng/ul	0.01

Target Compounds Qvalue

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

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