

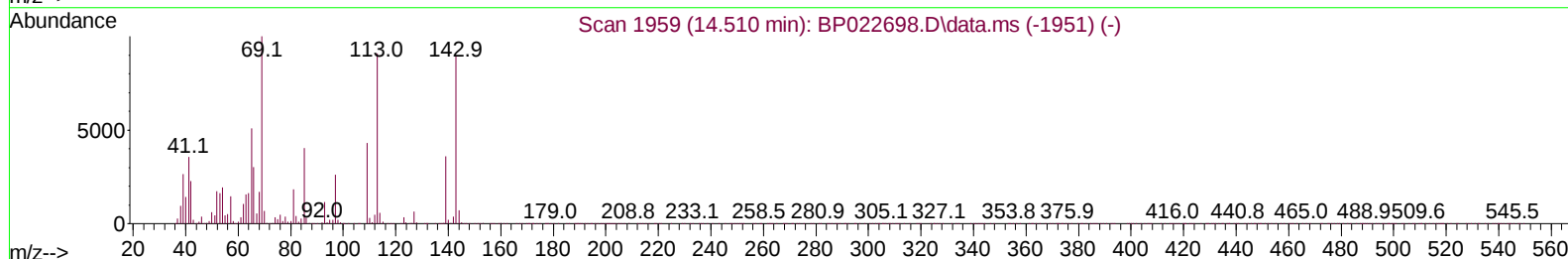
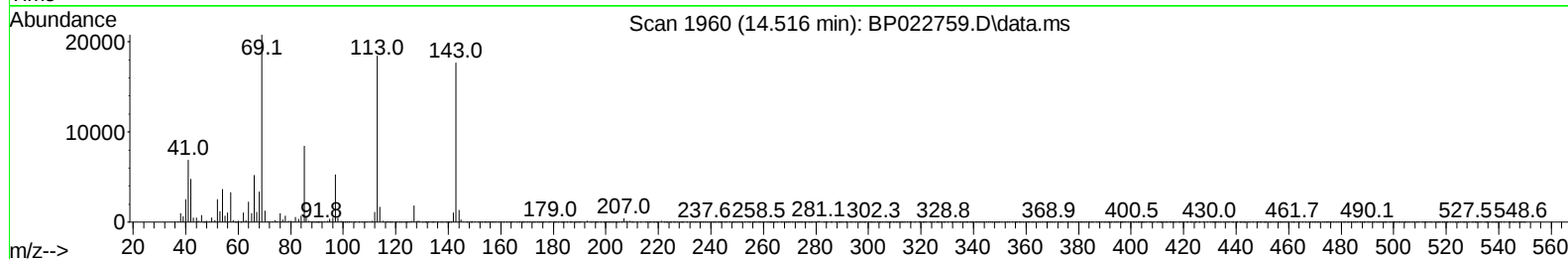
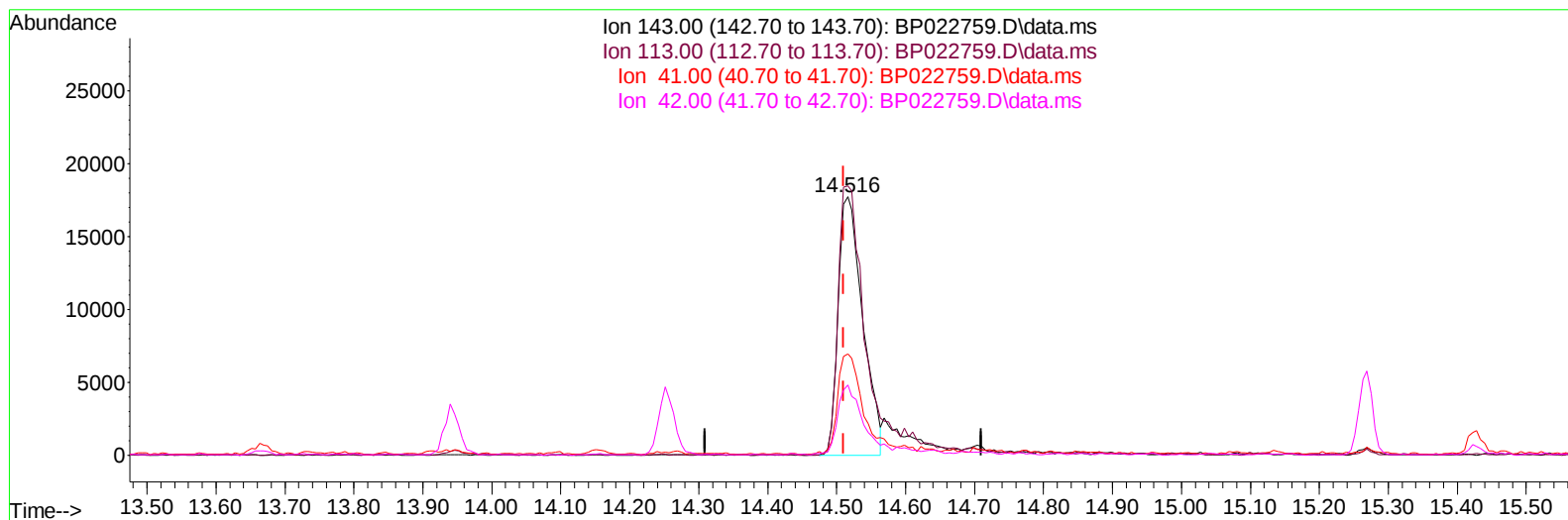
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103024\
Data File : BP022759.D
Acq On : 31 Oct 2024 07:21
Operator : RC/JU
Sample : P4592-05
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
F7H05

Manual IntegrationsAPPROVED

Quant Time: Oct 31 12:39:49 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: BP022759.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.516min (+ 0.006) 12.02 ng/u1

response 43618

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	104.15
41.00	35.90	39.11
42.00	23.70	27.08

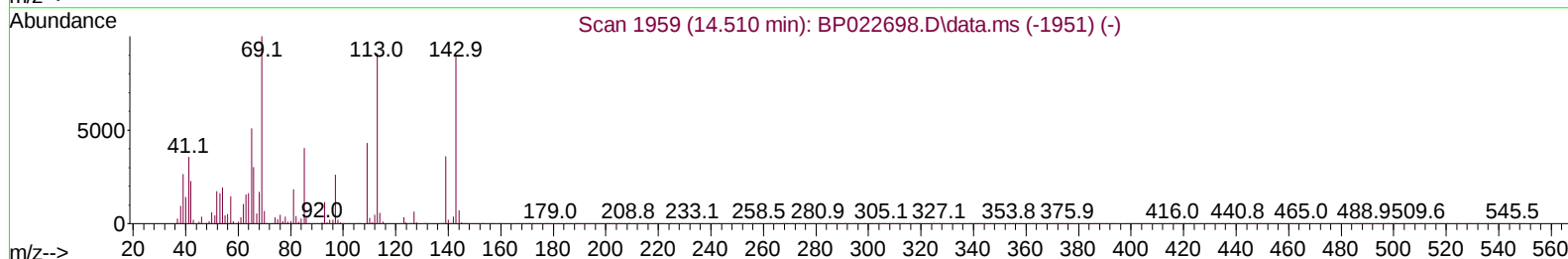
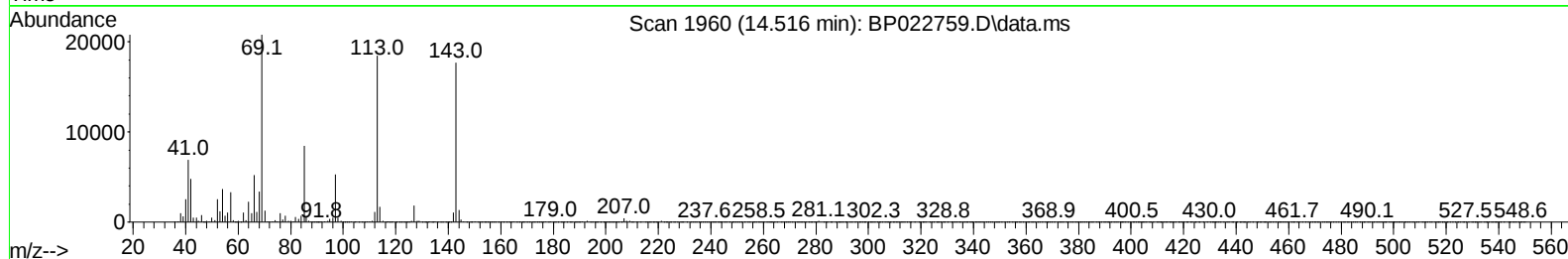
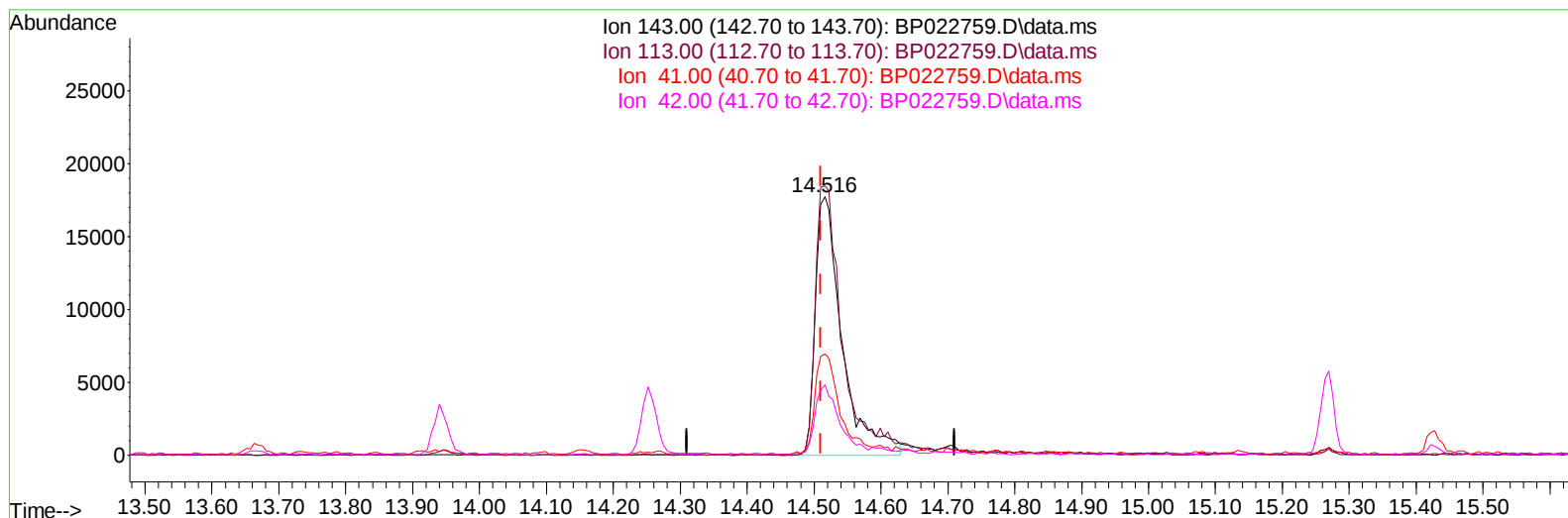
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(54) 4-Nitrophenol-d4 (S)

14.516min (+ 0.006) 13.60 ng/ul m

response 49317

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	104.15
41.00	35.90	39.11
42.00	23.70	27.08

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/04/2024
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Quant Time: Nov 01 01:34:54 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	89737	20.000	ng/ul	0.00
20) Naphthalene-d8	10.393	136	343852	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.251	164	272135	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.057	188	657358	20.000	ng/ul	0.00
79) Chrysene-d12	21.504	240	760923	20.000	ng/ul	# 0.00
88) Perylene-d12	24.757	264	871821	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	5534	2.396	ng/uL	0.00
4) Pyridine-d5	3.599	84	69585	10.977	ng/ul	0.00
7) Phenol-d5	6.834	99	115239	15.256	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.981	67	67651	15.238	ng/ul	0.00
11) 2-Chlorophenol-d4	7.181	132	89036	16.611	ng/ul	0.00
15) 4-Methylphenol-d8	8.340	113	100180	16.618	ng/ul	0.00
21) Nitrobenzene-d5	8.775	128	43648	16.509	ng/ul	0.00
24) 2-Nitrophenol-d4	9.493	143	53525	17.486	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.028	165	110416	16.972	ng/ul	0.00
31) 4-Chloroaniline-d4	10.540	131	122274	15.906	ng/ul	0.00
46) Dimethylphthalate-d6	13.669	166	404157	18.692	ng/ul	0.00
49) Acenaphthylene-d8	13.940	160	416128	18.120	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.516	143	49317m	13.596	ng/ul	0.00
60) Fluorene-d10	15.269	176	354695	18.913	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.428	200	40667	9.406	ng/ul	0.02
73) Anthracene-d10	17.163	188	613143	19.828	ng/ul	0.00
81) Pyrene-d10	19.545	212	803934	19.979	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.533	264	969314	20.819	ng/ul	0.01

Target Compounds Qvalue

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

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