

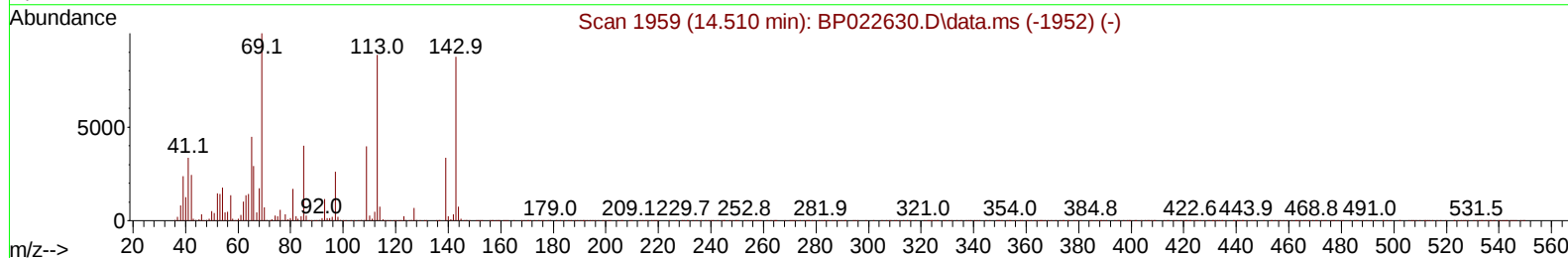
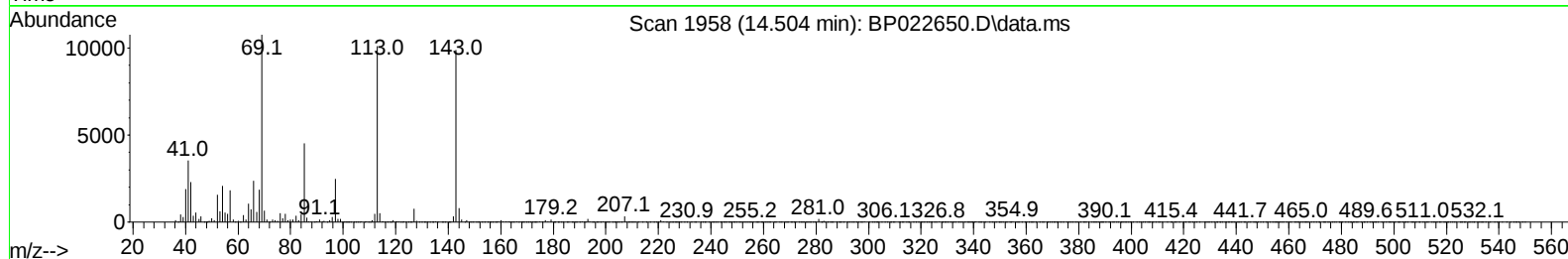
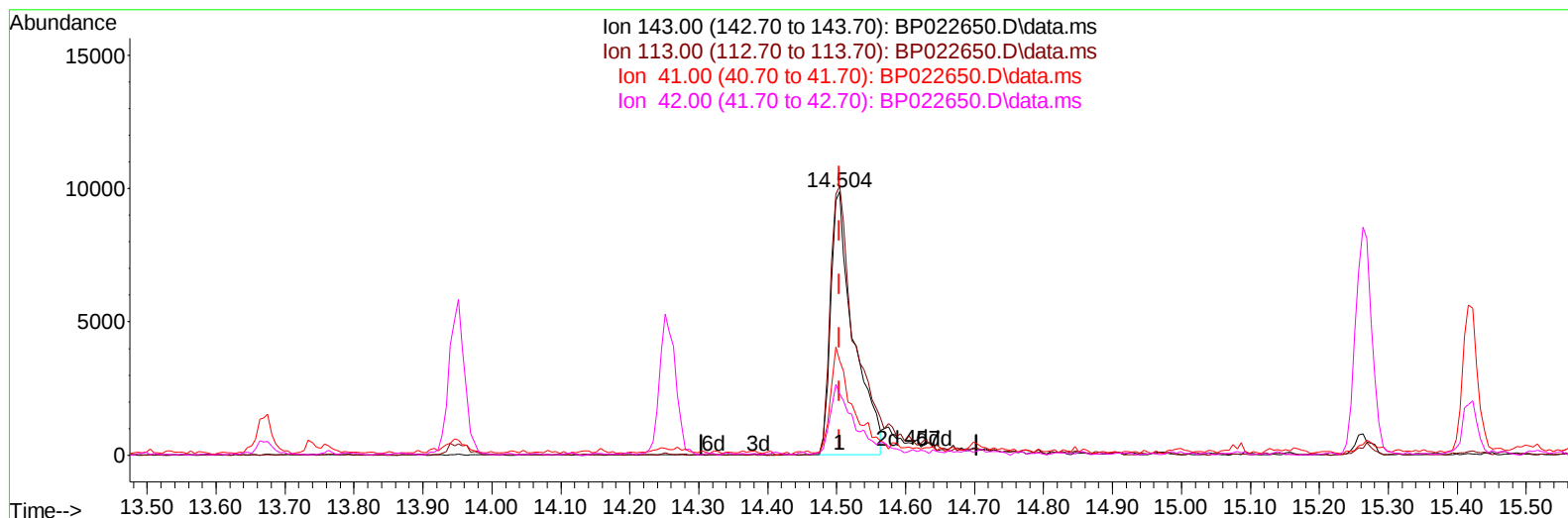
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102524\
Data File : BP022650.D
Acq On : 26 Oct 2024 20:03
Operator : RC/JU
Sample : P4532-10
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCNX1

Manual IntegrationsAPPROVED

Quant Time: Oct 28 00:27:05 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Oct 26 05:55:04 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/28/2024
Supervised By :mohammad ahmed 10/29/2024



TIC: BP022650.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.504min (+ 0.000) 5.84 ng/u1

response 22510

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	102.11
41.00	35.90	35.77
42.00	23.70	23.26

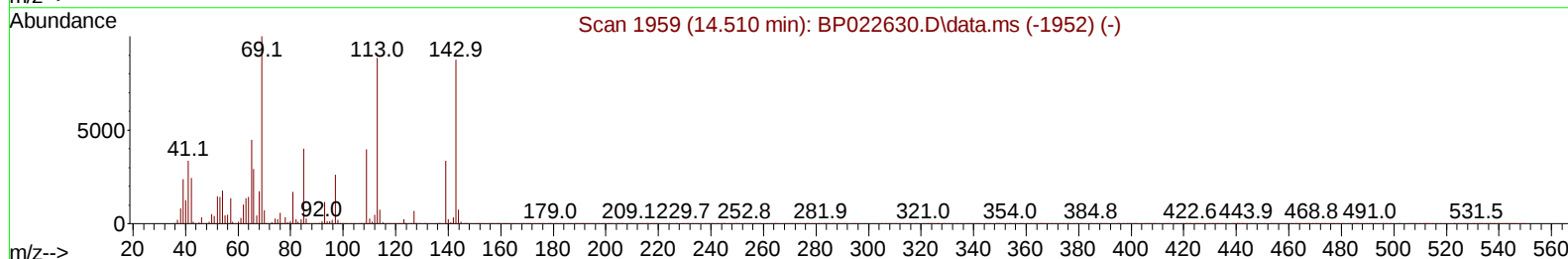
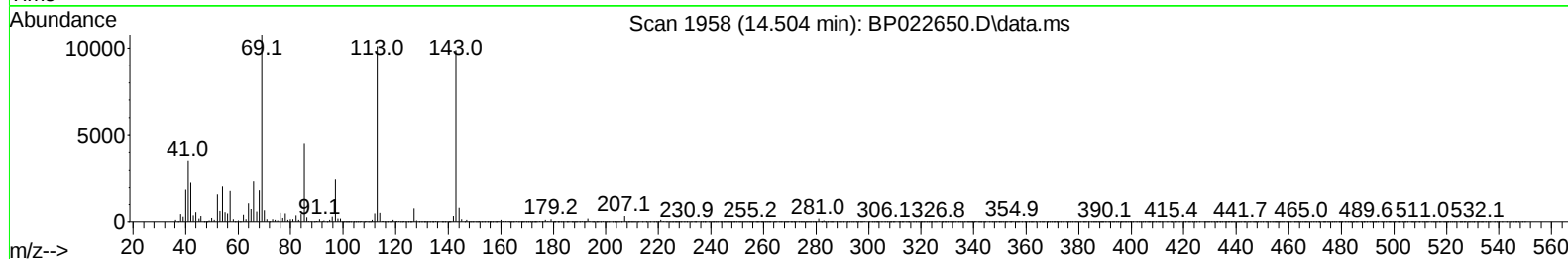
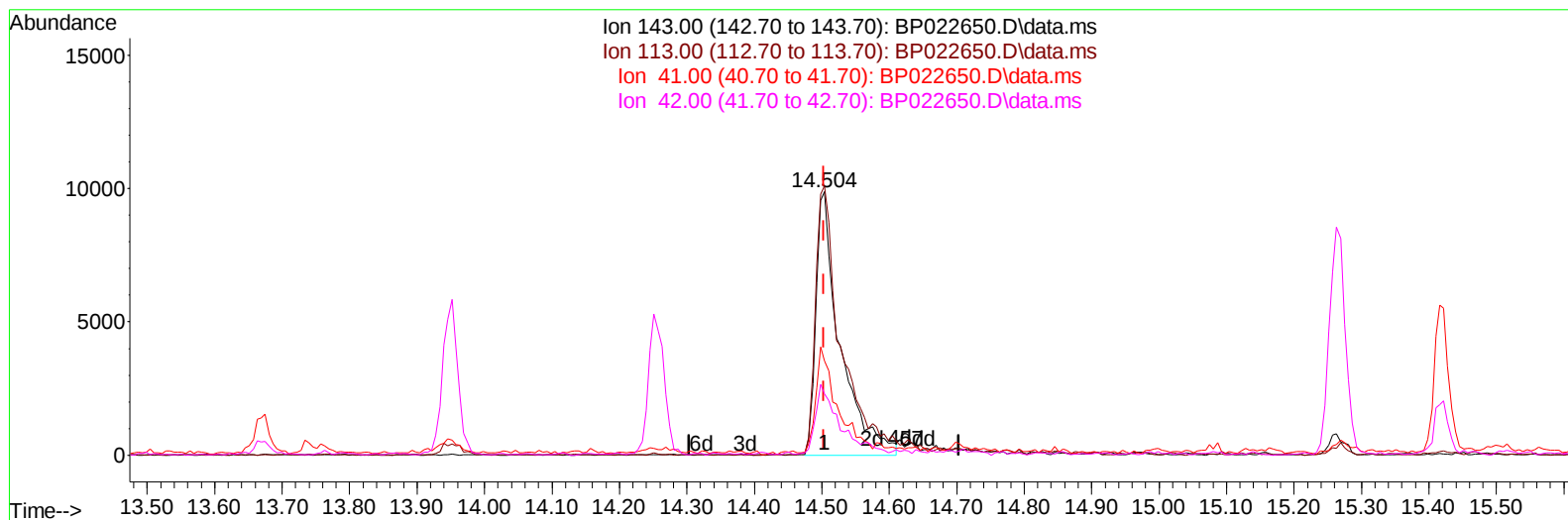
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TIC: BP022650.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.504min (+ 0.000) 6.36 ng/ul m

response 24498

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	102.11
41.00	35.90	35.77
42.00	23.70	23.26

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/28/2024
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Quant Time: Oct 28 00:27:37 2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 26 05:55:04 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.646	152	99143	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.399	136	377998	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.257	164	288972	20.000	ng/ul	-0.02
64) Phenanthrene-d10	17.063	188	687257	20.000	ng/ul	-0.02
79) Chrysene-d12	21.504	240	758210	20.000	ng/ul	# 0.00
88) Perylene-d12	24.751	264	880060	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.193	96	10067	3.946	ng/uL	0.00
4) Pyridine-d5	3.611	84	41502	5.926	ng/ul	0.00
7) Phenol-d5	6.840	99	57092	6.841	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	6.987	67	130776	26.661	ng/ul	-0.02
11) 2-Chlorophenol-d4	7.193	132	149062	25.172	ng/ul	-0.01
15) 4-Methylphenol-d8	8.352	113	115820	17.390	ng/ul	-0.01
21) Nitrobenzene-d5	8.787	128	83909	28.870	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.499	143	102270	30.393	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.040	165	206333	28.851	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.534	131	179745	21.270	ng/ul	-0.02
46) Dimethylphthalate-d6	13.669	166	752144	32.760	ng/ul	-0.01
49) Acenaphthylene-d8	13.951	160	745216	30.560	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.504	143	24498m	6.360	ng/ul	0.00
60) Fluorene-d10	15.263	176	643281	32.302	ng/ul	-0.03
65) 4,6-Dinitro-2-methylph...	15.422	200	135030	29.874	ng/ul	0.00
73) Anthracene-d10	17.169	188	1122377	34.716	ng/ul	0.00
81) Pyrene-d10	19.545	212	1411625	35.206	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.521	264	1662434	35.371	ng/ul	0.00

Target Compounds Qvalue

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

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