

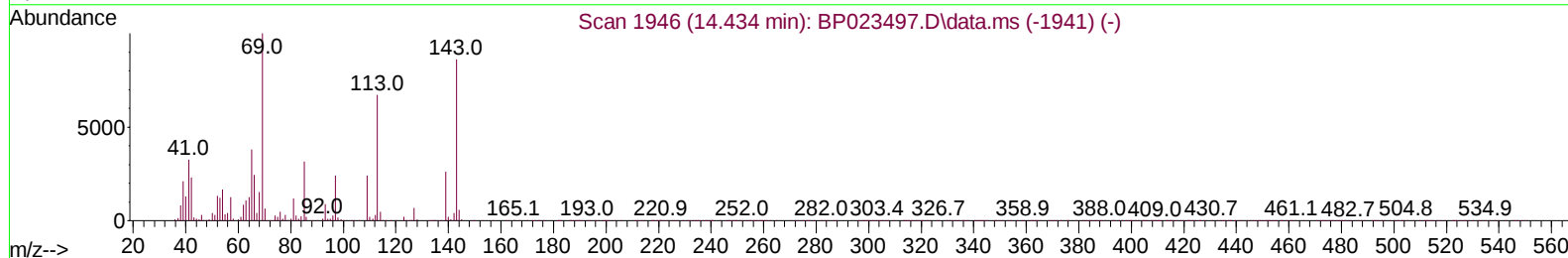
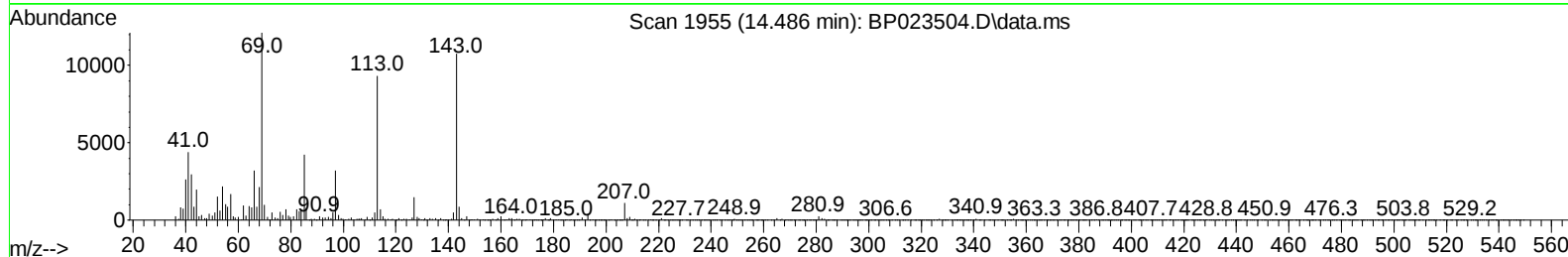
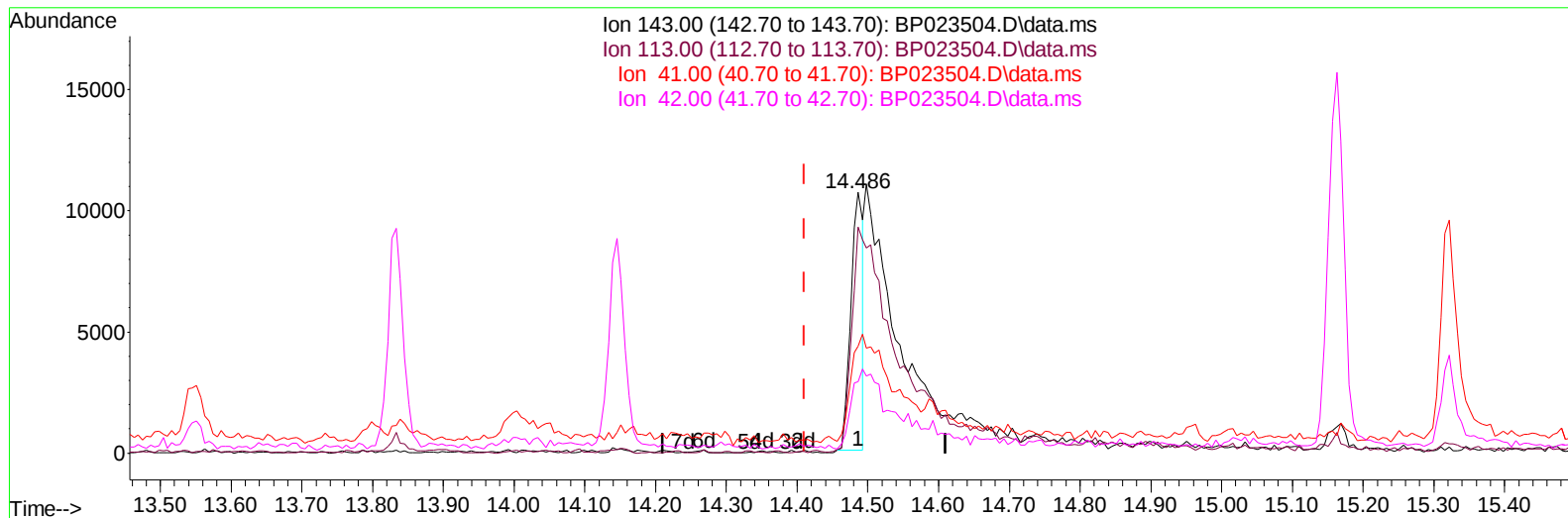
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
Data File : BP023504.D
Acq On : 18 Dec 2024 19:02
Operator : RC/JU
Sample : P5235-05 10X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AT3

Manual IntegrationsAPPROVED

Quant Time: Dec 19 01:08:31 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Dec 11 00:30:30 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 12/21/2024
Supervised By :mohammad ahmed 12/21/2024



TIC: BP023504.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.486min (+ 0.076) 2.61 ng/ul

response 12791

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	106.10	86.63
41.00	45.60	40.94
42.00	31.70	27.30

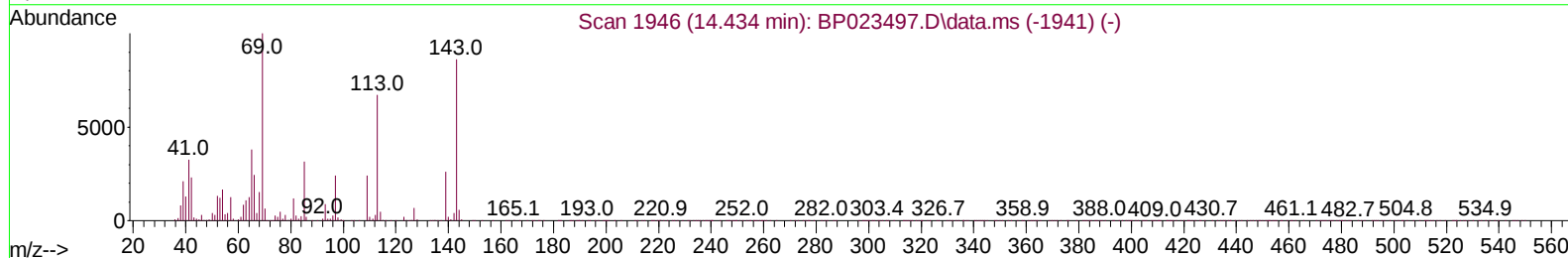
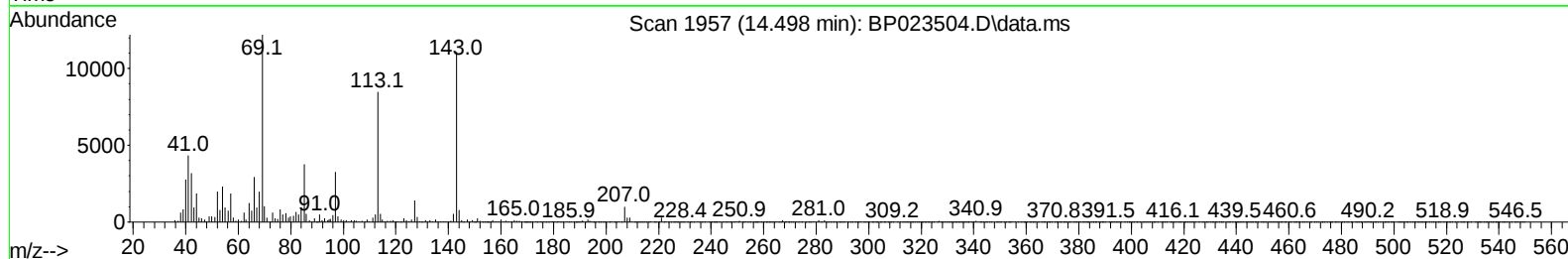
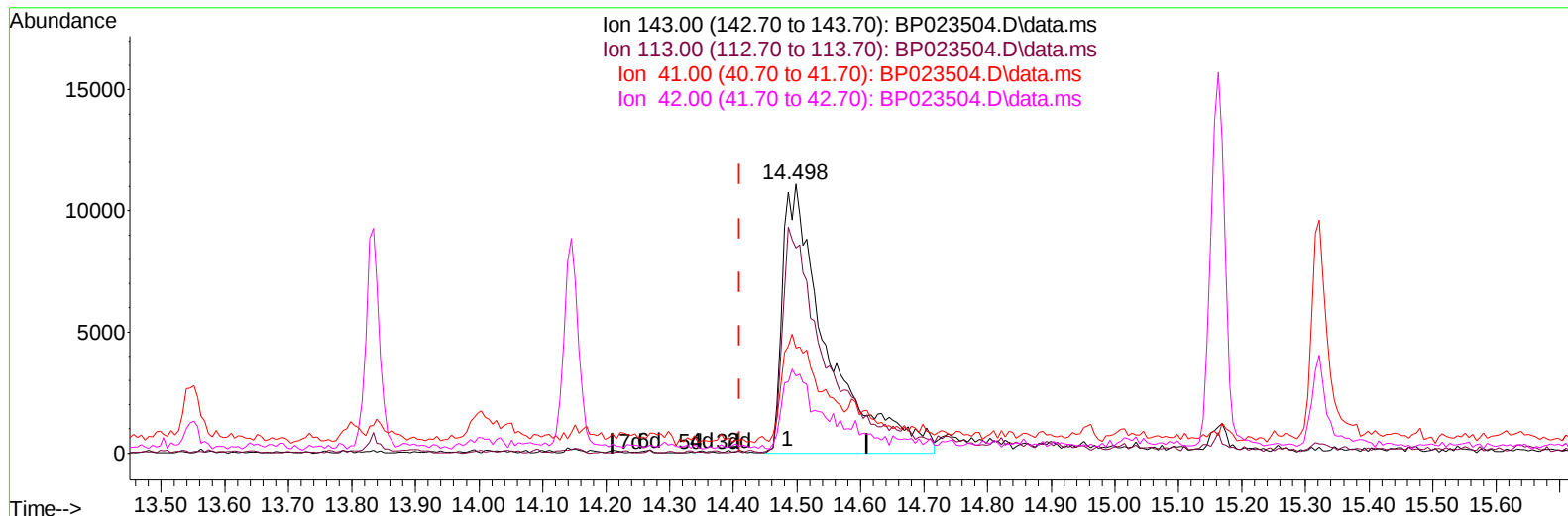
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TIC: BP023504.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.088) 11.07 ng/ul m

response 54274

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	106.10	76.30#
41.00	45.60	39.04
42.00	31.70	28.66

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 Sample : P5235-05 10X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AT3

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 12/21/2024
 Supervised By :mohammad ahmed 12/21/2024

Quant Time: Dec 19 03:55:12 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 11 00:30:30 2024
 Response via : Initial Calibration

Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.528	152		213319	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.269	136		770673	20.000	ng/ul	#-0.01
38) Acenaphthene-d10	14.145	164		515406	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.957	188		1120601	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.386	240		1360018	20.000	ng/ul	# 0.00
88) Perylene-d12	24.545	264		1527563	20.000	ng/ul	-0.03
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.093	96		23671	4.085	ng/uL	-0.01
4) Pyridine-d5	3.522	84		115291	7.738	ng/ul	0.00
7) Phenol-d5	6.751	99		165344	9.879	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.881	67		380498	36.032	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.081	132		374096	30.283	ng/ul	0.00
15) 4-Methylphenol-d8	8.251	113		287336	21.270	ng/ul	0.00
21) Nitrobenzene-d5	8.675	128		189502	32.929	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.381	143		228575	34.462	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	9.922	165		413308	29.525	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.422	131		412914	26.331	ng/ul	-0.02
46) Dimethylphthalate-d6	13.551	166		1348813	33.711	ng/ul	0.00
49) Acenaphthylene-d8	13.833	160		1255086	29.739	ng/ul	0.00
54) 4-Nitrophenol-d4	14.498	143		54274m	11.069	ng/ul	0.09
60) Fluorene-d10	15.163	176		1080289	30.696	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.322	200		219907	32.471	ng/ul	0.01
73) Anthracene-d10	17.051	188		1995794	39.024	ng/ul	0.00
81) Pyrene-d10	19.439	212		2548090	35.861	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.333	264		3056329	37.375	ng/ul	-0.01

Target Compounds Qvalue

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

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