

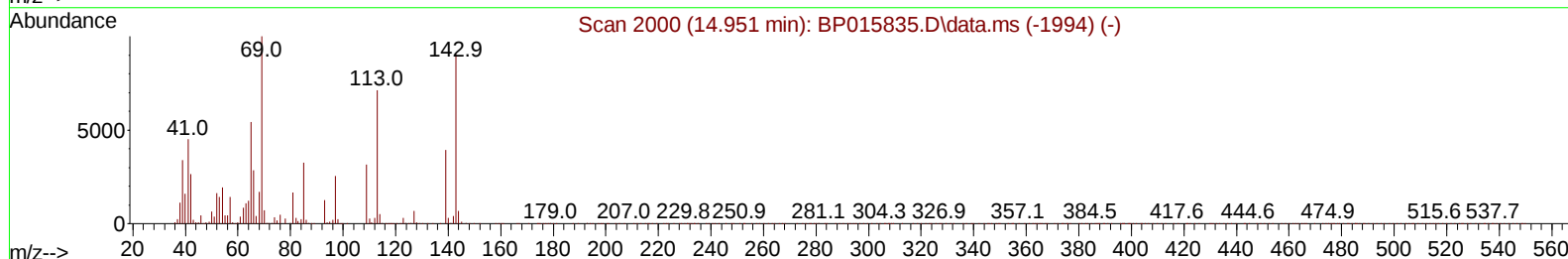
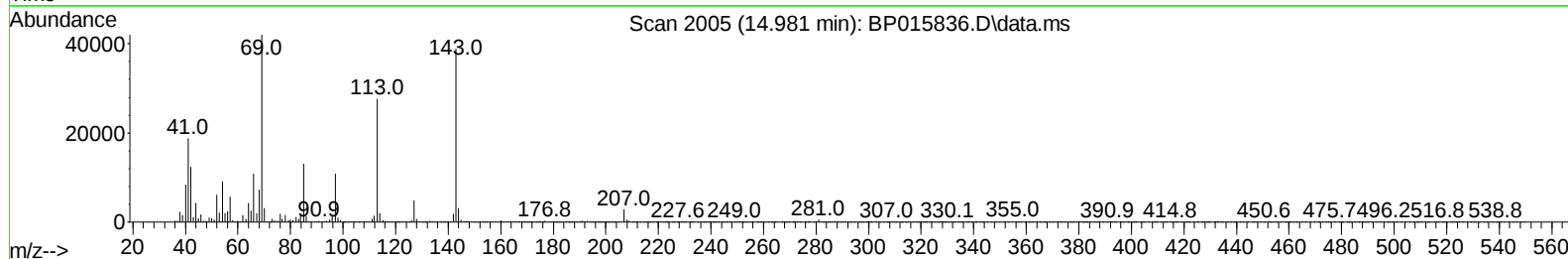
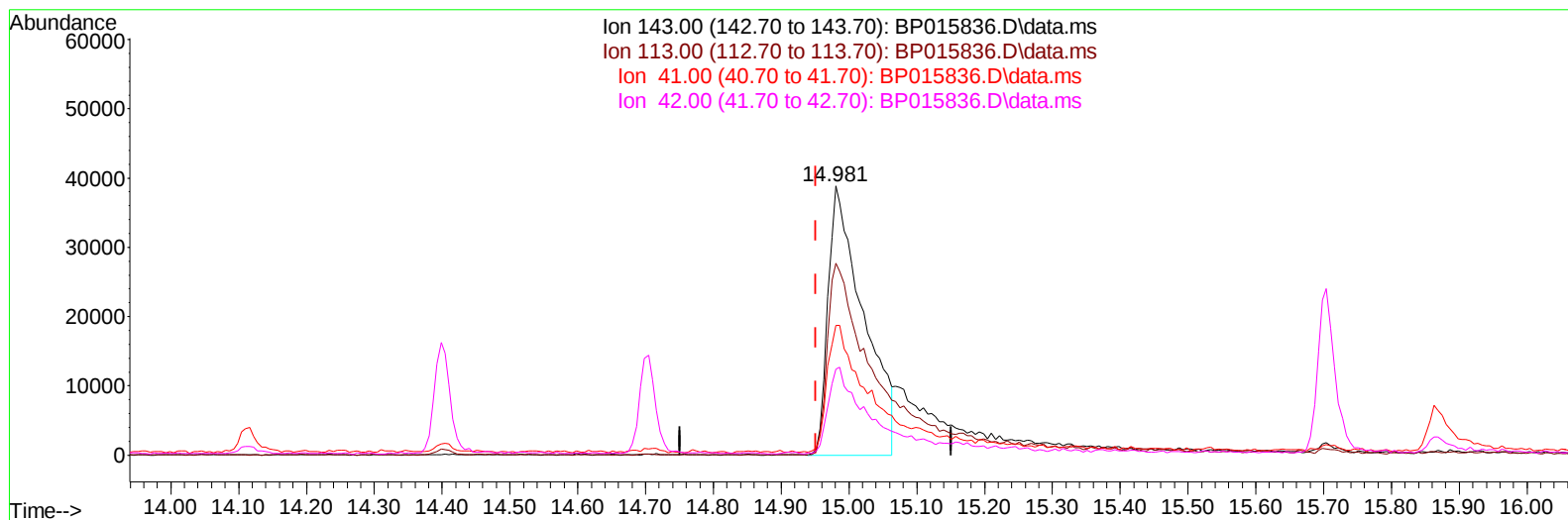
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015836.D
Acq On : 30 Jun 2023 10:34
Operator : MA/JU
Sample : PB153635BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK635

Manual IntegrationsAPPROVED

Quant Time: Jun 30 22:31:56 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jun 30 22:30:25 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/01/2023
Supervised By :mohammad ahmed 07/01/2023



TIC: BP015836.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.981min (+ 0.029) 21.31 ng/u1

response 140826

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.00	71.28#
41.00	45.70	48.33
42.00	31.20	31.87

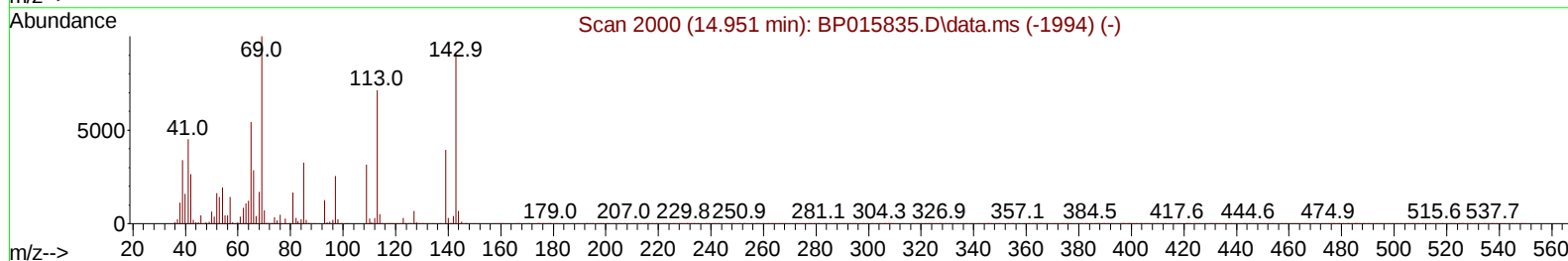
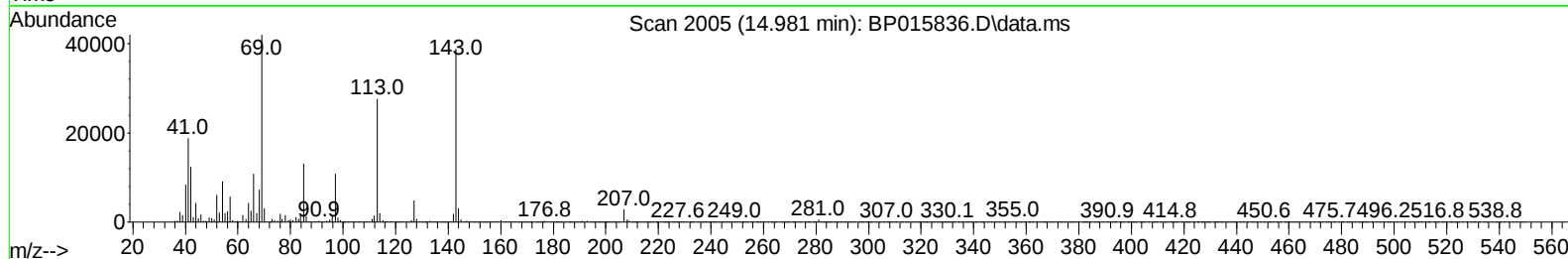
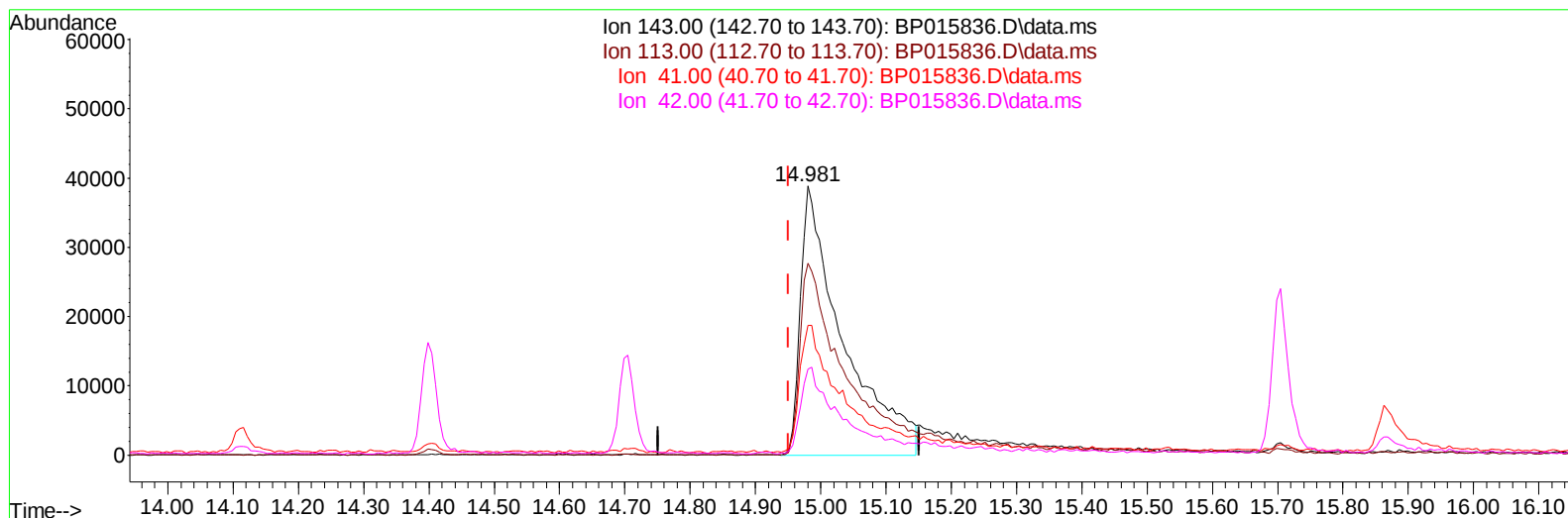
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TIC: BP015836.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.981min (+ 0.029) 26.46 ng/ul m

response 174807

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.00	71.28#
41.00	45.70	48.33
42.00	31.20	31.87

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Instrument :
 BNA_P
 ClientSampleId :
 SBLK635

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/01/2023
 Supervised By :mohammad ahmed 07/01/2023

Quant Time: Jun 30 22:42:17 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 30 22:30:25 2023
 Response via : Initial Calibration

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

Internal Standards							
1) 1,4-Dichlorobenzene-d4	8.052	152		228219	20.000	ng/ul	0.00
20) Naphthalene-d8	10.881	136		1031016	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.704	164		647755	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.463	188		1378758	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.563	240		1028974	20.000	ng/ul	# 0.00
88) Perylene-d12	24.104	264		1016585	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.375	96		37520	5.915	ng/uL	0.00
4) Pyridine-d5	3.822	84		510525	31.928	ng/ul	0.00
7) Phenol-d5	7.228	99		641926	32.874	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.381	67		431019	35.678	ng/ul	0.00
11) 2-Chlorophenol-d4	7.581	132		510827	34.655	ng/ul	0.00
15) 4-Methylphenol-d8	8.787	113		537013	34.812	ng/ul	0.00
21) Nitrobenzene-d5	9.240	128		255726	32.455	ng/ul	0.00
24) 2-Nitrophenol-d4	9.969	143		276975	30.118	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.534	165		466109	28.443	ng/ul	0.02
31) 4-Chloroaniline-d4	11.046	131		704934	33.486	ng/ul	0.01
46) Dimethylphthalate-d6	14.116	166		1627398	32.505	ng/ul	0.00
49) Acenaphthylene-d8	14.398	160		1868993	33.208	ng/ul	0.00
54) 4-Nitrophenol-d4	14.981	143		174807m	26.458	ng/ul	0.03
60) Fluorene-d10	15.704	176		1367378	33.169	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.869	200		153491	18.102	ng/ul	0.01
73) Anthracene-d10	17.569	188		2158802	34.278	ng/ul	0.00
81) Pyrene-d10	19.792	212		2380589	35.431	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.933	264		1824326	35.575	ng/ul	0.00

Target Compounds Qvalue

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

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