

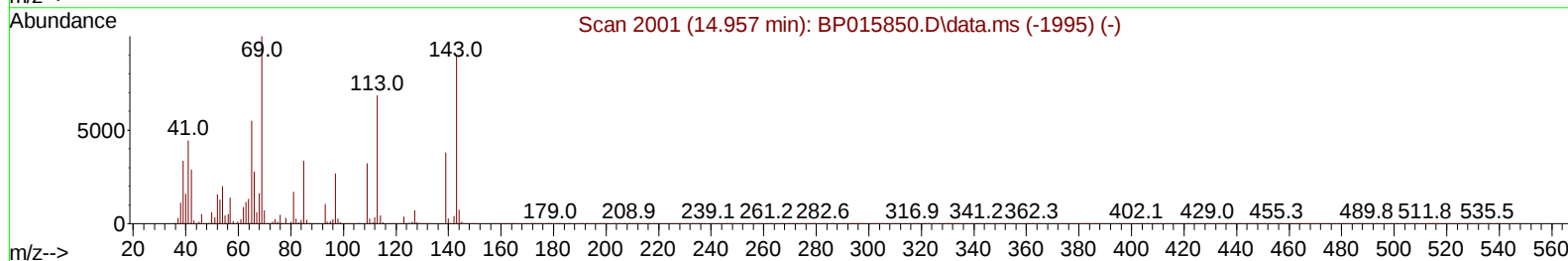
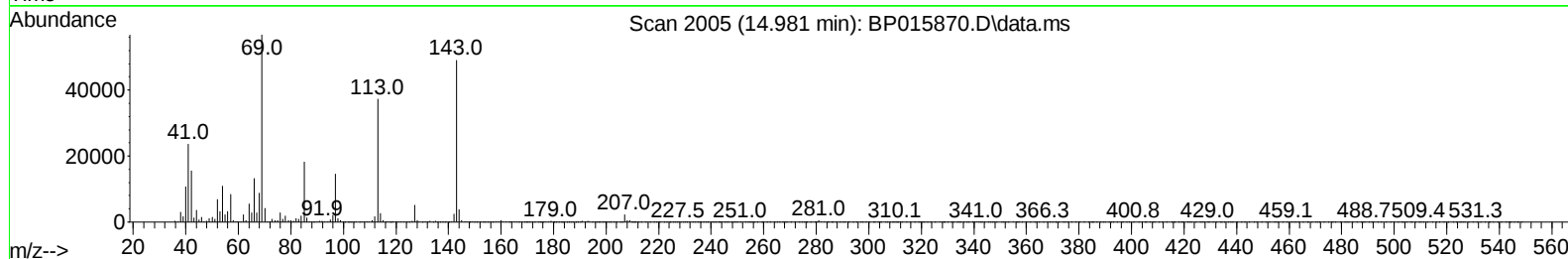
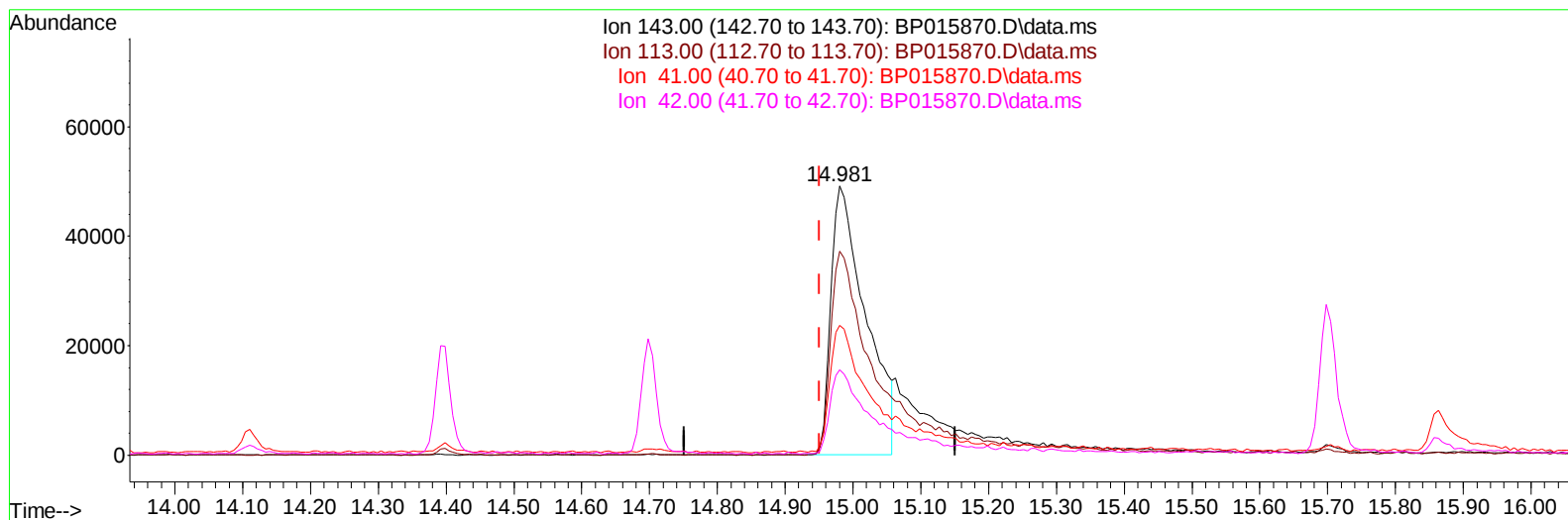
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015870.D
Acq On : 01 Jul 2023 06:24
Operator : MA/JU
Sample : PB153698BL
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK698

Manual IntegrationsAPPROVED

Quant Time: Jul 01 06:59:47 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: BP015870.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.981min (+ 0.029) 19.60 ng/u1

response 175028

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.00	75.92
41.00	45.70	48.17
42.00	31.20	31.63

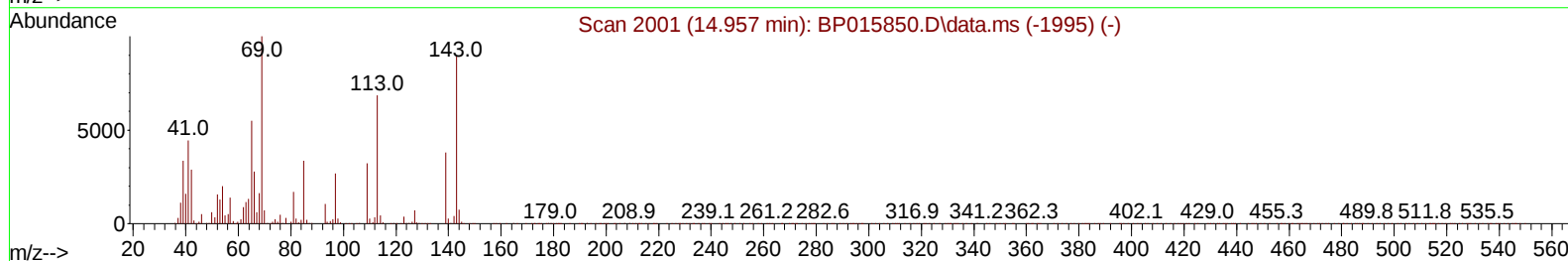
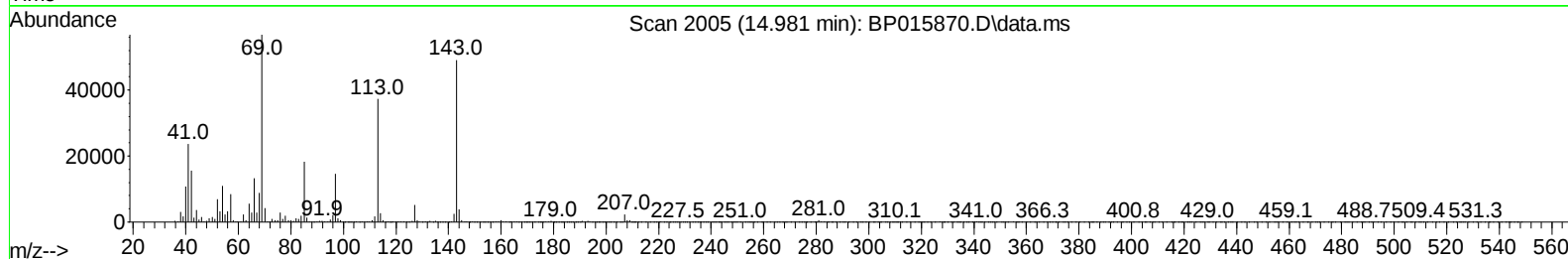
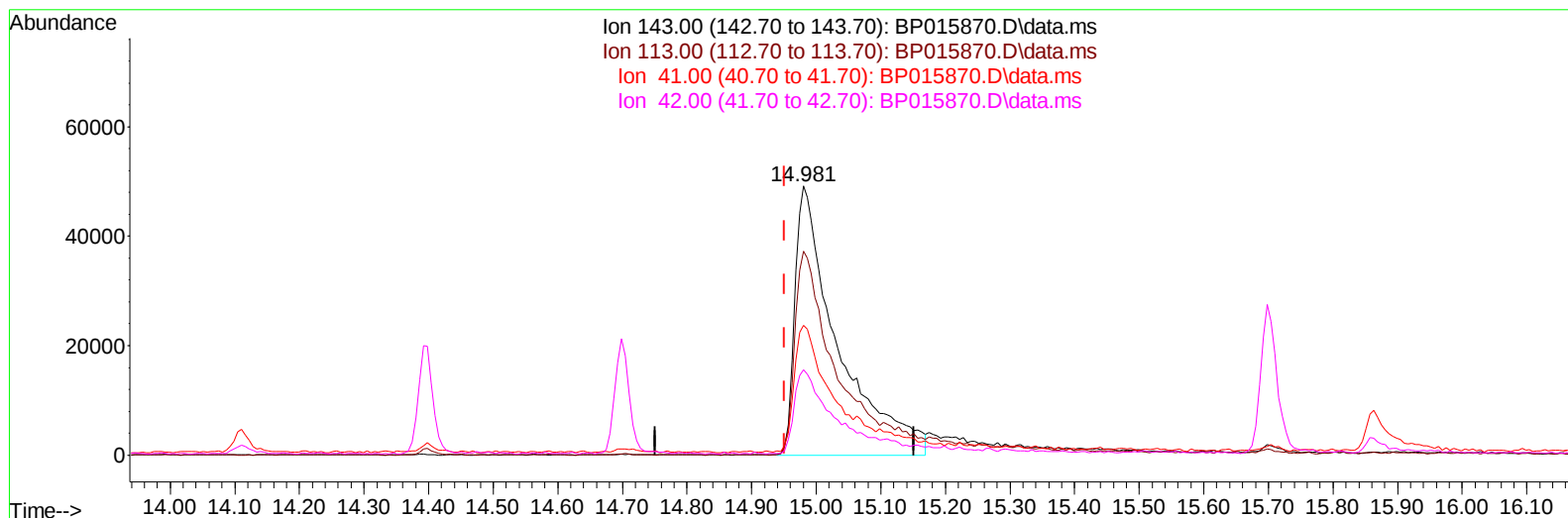
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TIC: BP015870.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.981min (+ 0.029) 25.12 ng/ul m

response 224333

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.00	75.92
41.00	45.70	48.17
42.00	31.20	31.63

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 ALS Vial : 37 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

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

Quant Time: Jul 01 07:00:29 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		306013	20.000	ng/ul	0.00
20) Naphthalene-d8	10.875	136		1393861	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.698	164		875464	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.463	188		1859325	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.557	240		1353530	20.000	ng/ul	# 0.00
88) Perylene-d12	24.098	264		1401662	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.375	96		40315	4.740	ng/uL	0.00
4) Pyridine-d5	3.817	84		591263	27.577	ng/ul	0.00
7) Phenol-d5	7.228	99		732838	27.989	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.375	67		497277	30.698	ng/ul	0.00
11) 2-Chlorophenol-d4	7.581	132		588484	29.774	ng/ul	0.00
15) 4-Methylphenol-d8	8.781	113		616530	29.807	ng/ul	0.00
21) Nitrobenzene-d5	9.240	128		298741	28.045	ng/ul	0.00
24) 2-Nitrophenol-d4	9.969	143		323587	26.027	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.528	165		507086	22.888	ng/ul	0.01
31) 4-Chloroaniline-d4	11.040	131		817621	28.729	ng/ul	0.00
46) Dimethylphthalate-d6	14.110	166		1868803	27.618	ng/ul	0.00
49) Acenaphthylene-d8	14.398	160		2131316	28.019	ng/ul	0.00
54) 4-Nitrophenol-d4	14.981	143		224333m	25.122	ng/ul	0.03
60) Fluorene-d10	15.698	176		1574289	28.255	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.863	200		185142	16.191	ng/ul	0.00
73) Anthracene-d10	17.563	188		2494151	29.367	ng/ul	0.00
81) Pyrene-d10	19.792	212		2644839	29.925	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.927	264		2054402	29.055	ng/ul	0.00

Target Compounds Qvalue

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

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