

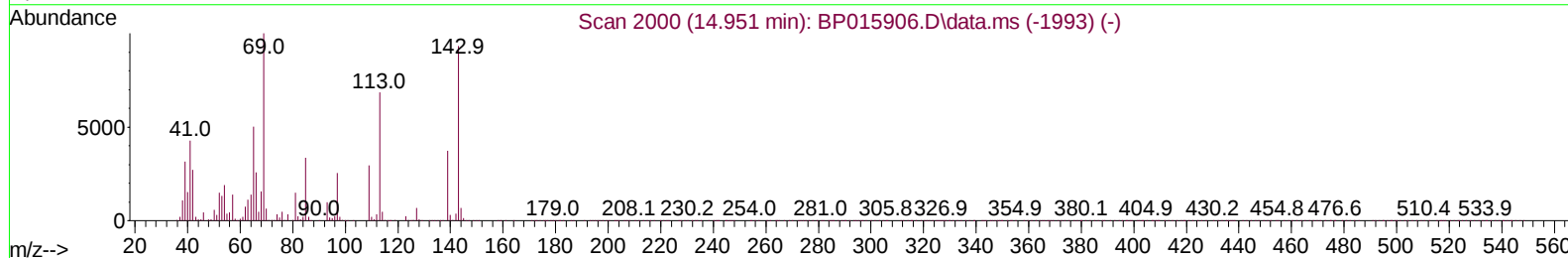
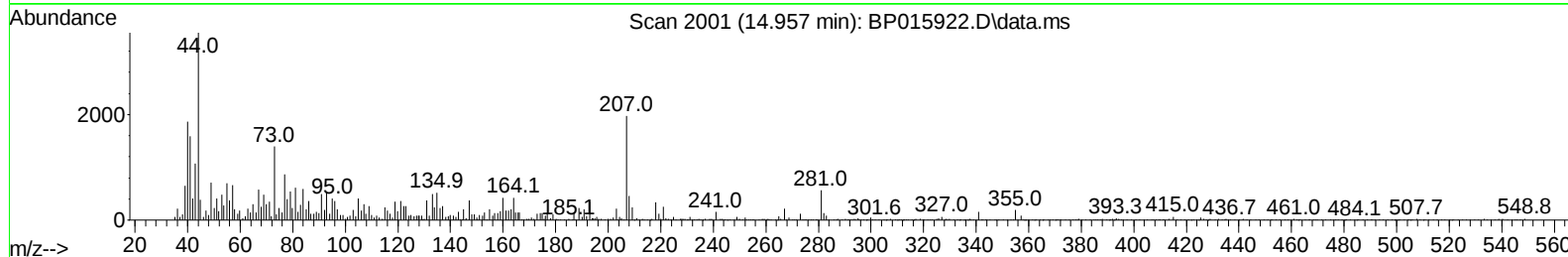
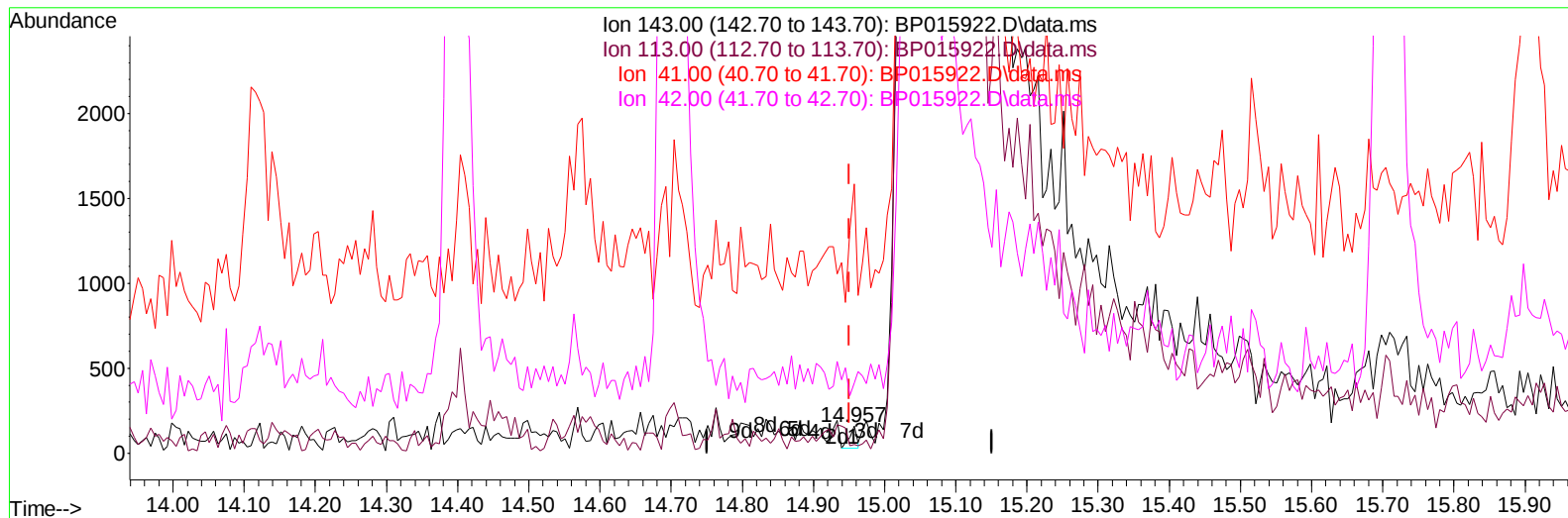
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070223\
Data File : BP015922.D
Acq On : 02 Jul 2023 13:27
Operator : MA/JU
Sample : 03243-18
Misc :
ALS Vial : 89 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGD8

Manual IntegrationsAPPROVED

Quant Time: Jul 02 14:28:30 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sun Jul 02 06:12:16 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/03/2023
Supervised By :mohammad ahmed 07/03/2023



TIC: BP015922.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.957min (+ 0.006) 0.01 ng/u1

response 93

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.00	30.54#
41.00	45.70	948.50#
42.00	31.20	246.71#

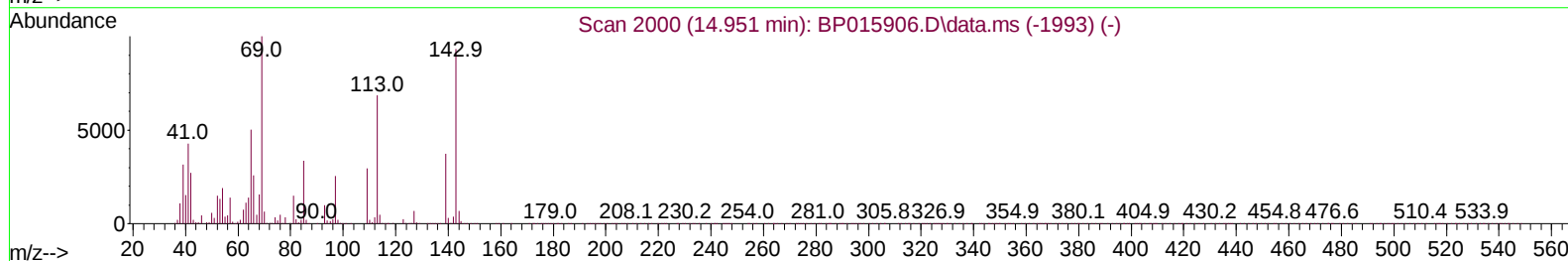
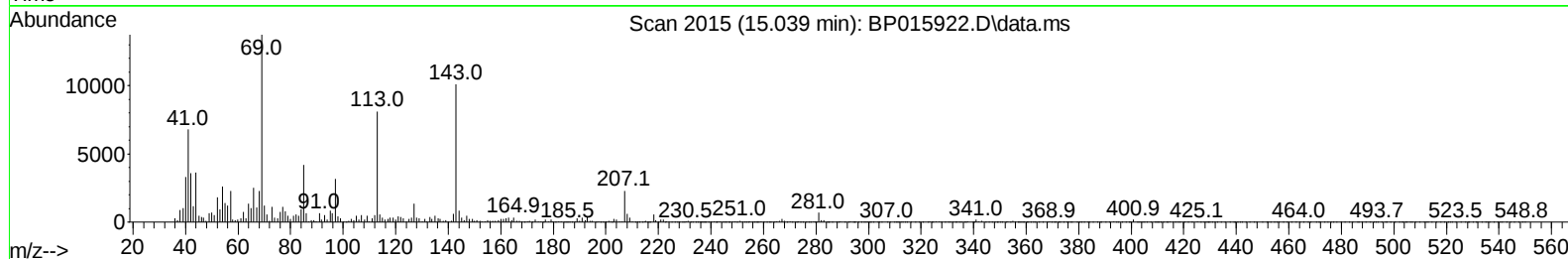
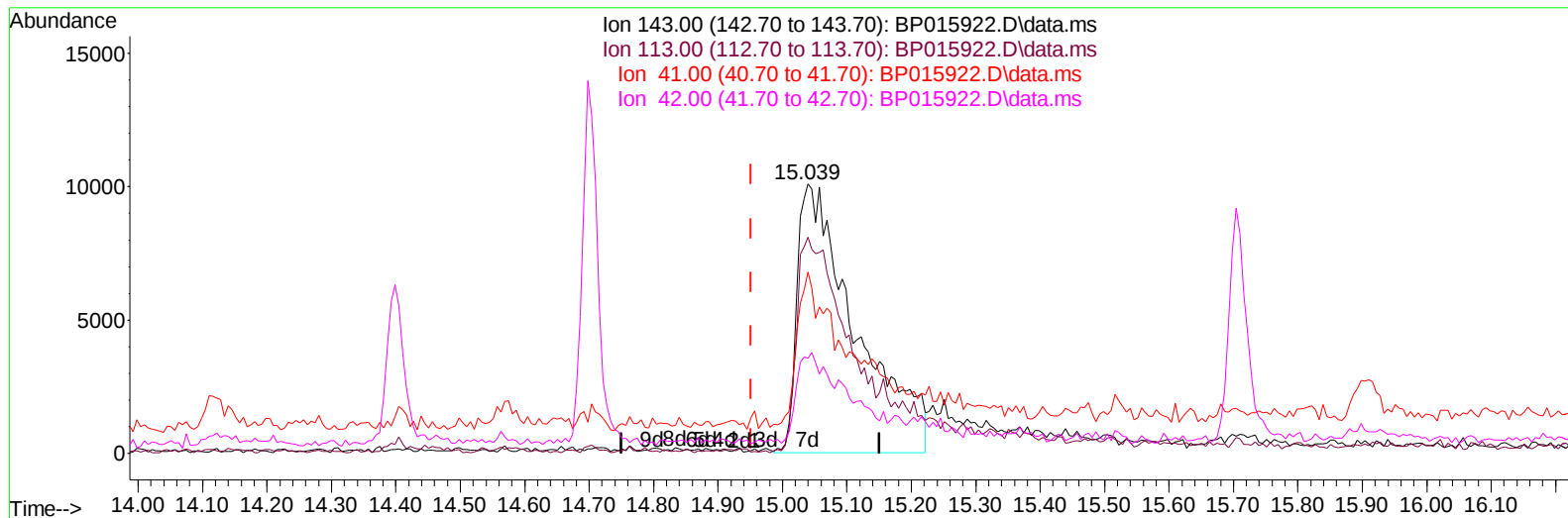
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TIC: BP015922.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.039min (+ 0.088) 10.26 ng/ul m

response 63807

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.00	80.27
41.00	45.70	67.22#
42.00	31.20	35.34

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 QLast Update : Sun Jul 02 06:12:16 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.051	152	235748	20.000	ng/ul	0.00
20) Naphthalene-d8	10.881	136	1054287	20.000	ng/ul	# 0.01
38) Acenaphthene-d10	14.698	164	609438	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.463	188	1092238	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.557	240	781268	20.000	ng/ul	# 0.00
88) Perylene-d12	24.098	264	915264	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.375	96	14387	2.196	ng/uL	0.00
4) Pyridine-d5	3.834	84	95719	5.795	ng/ul	0.02
7) Phenol-d5	7.251	99	258479	12.814	ng/ul	0.03
9) Bis-(2-Chloroethyl)eth...	7.381	67	179967	14.421	ng/ul	0.01
11) 2-Chlorophenol-d4	7.587	132	216742	14.235	ng/ul	0.01
15) 4-Methylphenol-d8	8.810	113	176647	11.086	ng/ul	0.03
21) Nitrobenzene-d5	9.251	128	103975	12.905	ng/ul	0.02
24) 2-Nitrophenol-d4	9.981	143	112542	11.968	ng/ul	0.02
28) 2,4-Dichlorophenol-d3	10.569	165	184166	10.990	ng/ul	0.06
31) 4-Chloroaniline-d4	11.075	131	206023	9.571	ng/ul	0.05
46) Dimethylphthalate-d6	14.116	166	665157	14.121	ng/ul	0.01
49) Acenaphthylene-d8	14.398	160	739891	13.973	ng/ul	0.00
54) 4-Nitrophenol-d4	15.039	143	63807m	10.265	ng/ul	0.09
60) Fluorene-d10	15.704	176	544270	14.033	ng/ul	0.01
65) 4,6-Dinitro-2-methylph...	15.898	200	36418	5.422	ng/ul	0.04
73) Anthracene-d10	17.569	188	752596	15.085	ng/ul	0.01
81) Pyrene-d10	19.792	212	829667	16.263	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.933	264	763800	16.543	ng/ul	0.00

Target Compounds Qvalue

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

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