

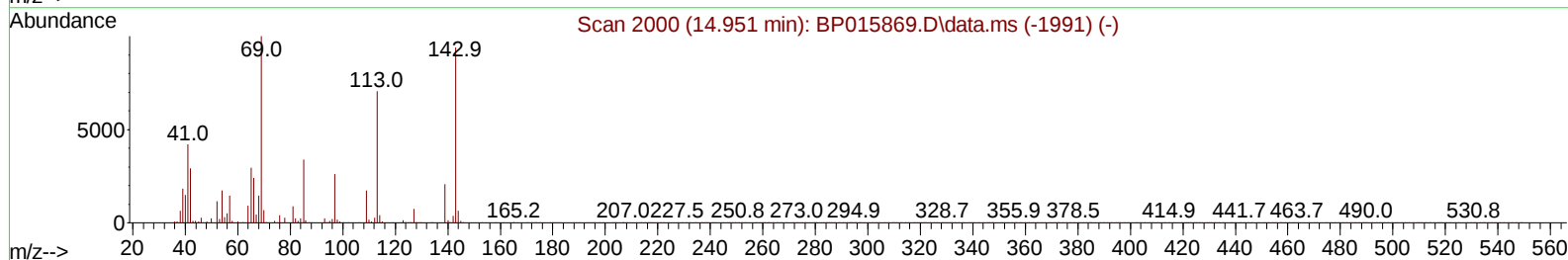
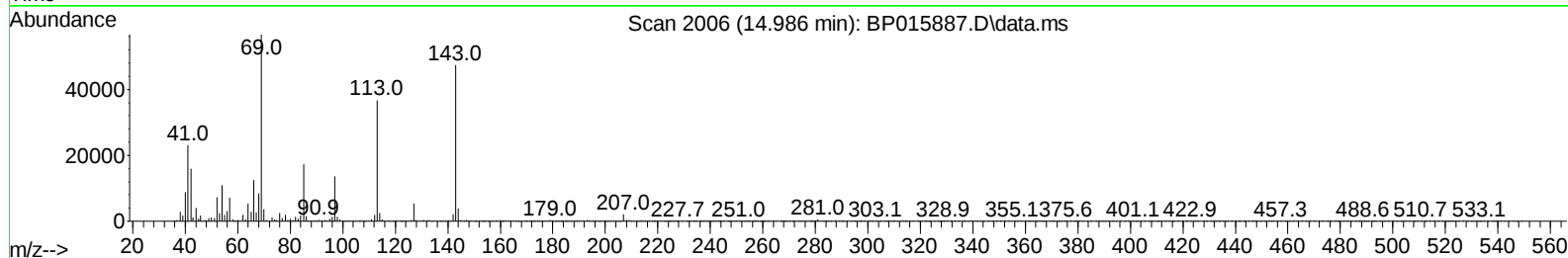
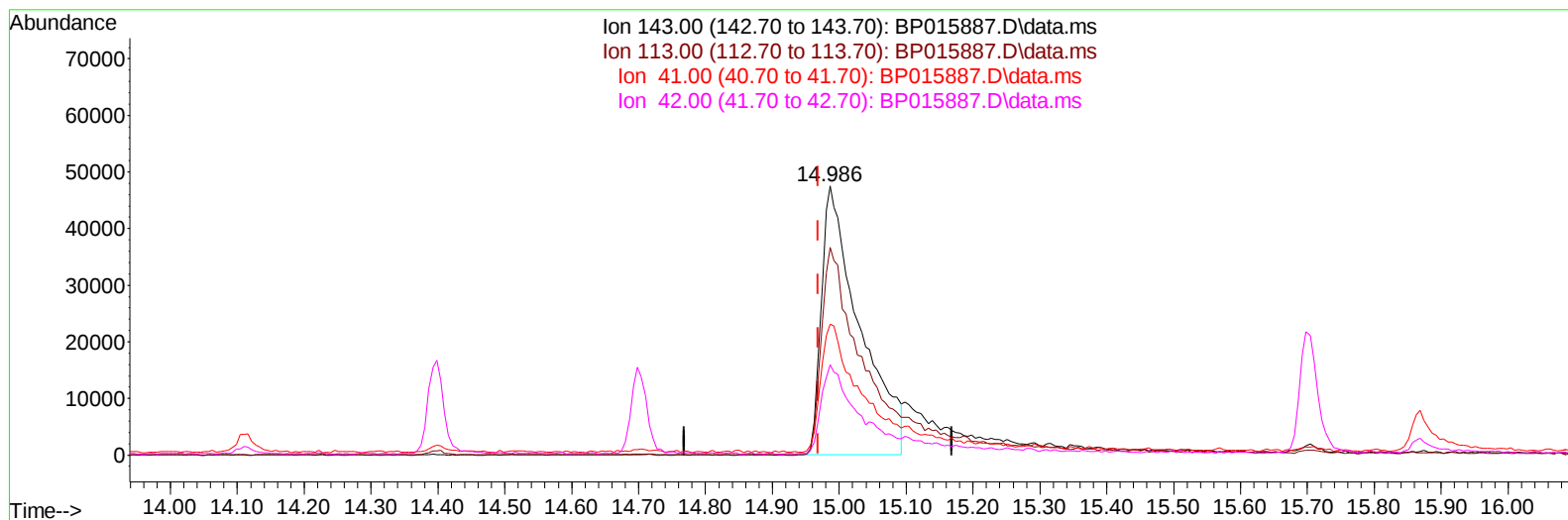
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015887.D
Acq On : 01 Jul 2023 17:05
Operator : MA/JU
Sample : PB153707BL
Misc :
ALS Vial : 56 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK707

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/02/2023
Supervised By :Sohil Jodhani 07/02/2023

Quant Time: Jul 02 02:42:42 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sun Jul 02 02:29:46 2023
Response via : Initial Calibration



TIC: BP015887.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.986min (+ 0.017) 27.23 ng/u1

response 188308

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.00	77.26
41.00	45.70	48.63
42.00	31.20	33.65

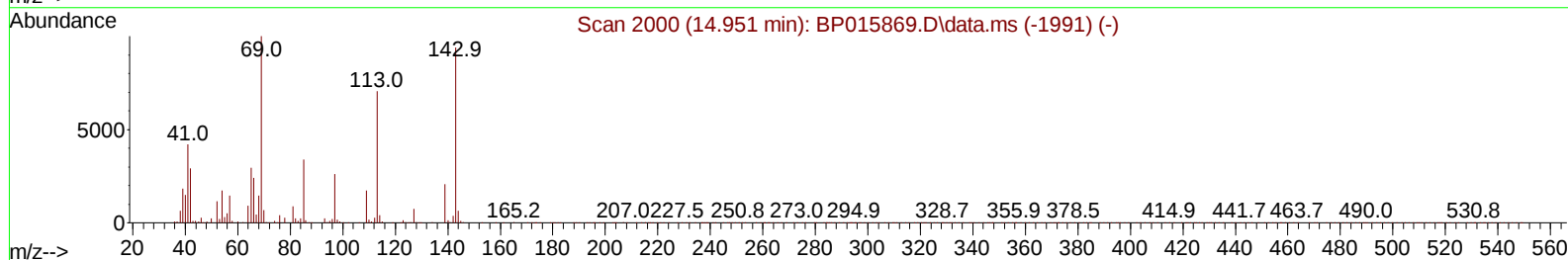
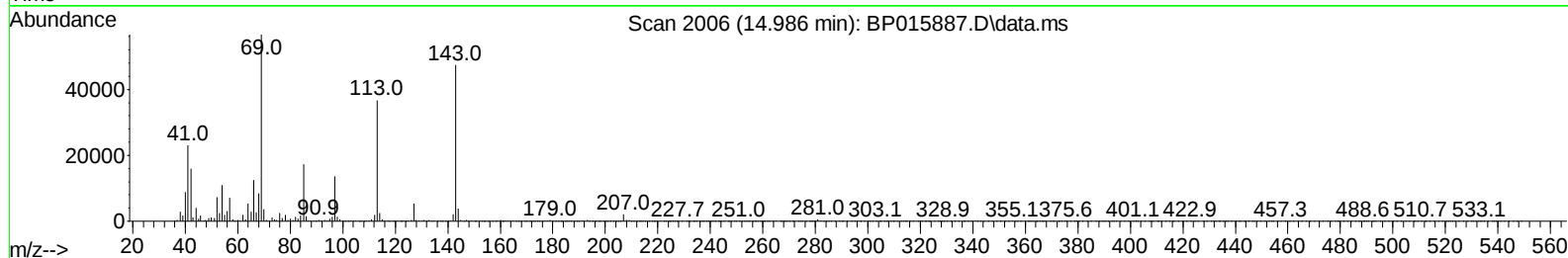
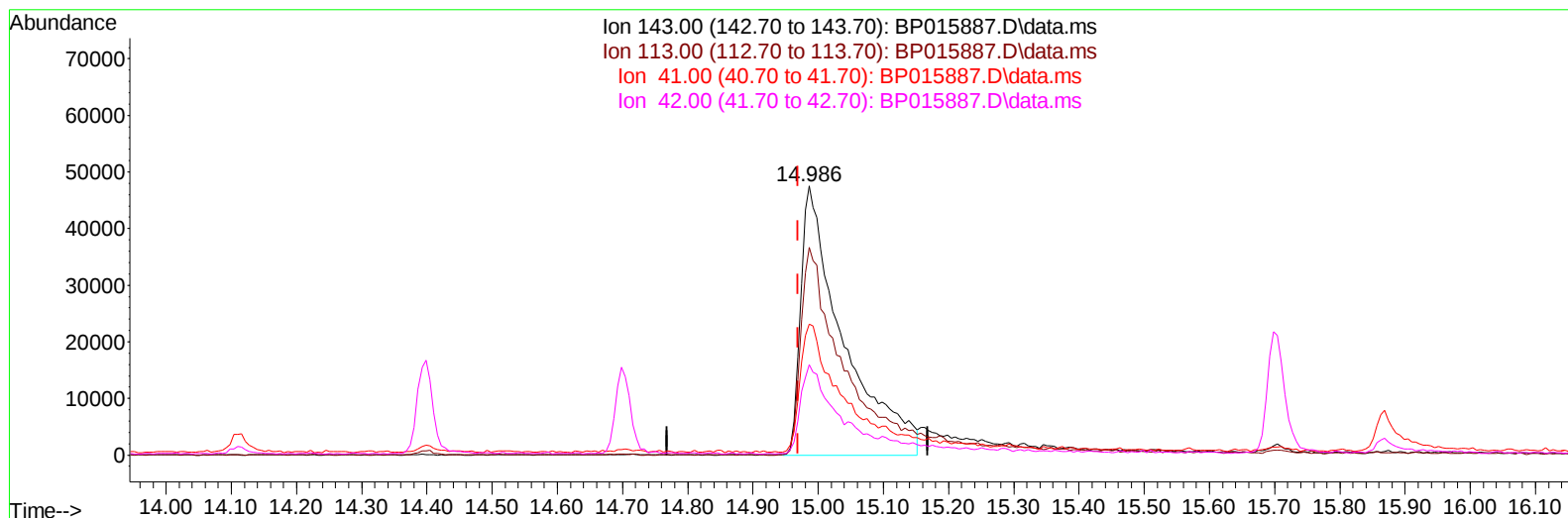
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TIC: BP015887.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.986min (+ 0.017) 30.77 ng/ul m

response 212821

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.00	77.26
41.00	45.70	48.63
42.00	31.20	33.65

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 Operator : MA/JU
 Sample : PB153707BL
 Misc :
 ALS Vial : 56 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/02/2023
 Supervised By :Sohil Jodhani 07/02/2023

Quant Time: Jul 02 02:51:27 2023
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 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Jul 02 02:29:46 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.051	152	279687	20.000	ng/ul	0.00
20) Naphthalene-d8	10.881	136	1153586	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.704	164	678062	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.463	188	1479686	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.557	240	1350722	20.000	ng/ul	# 0.00
88) Perylene-d12	24.098	264	1552394	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.375	96	43907	5.648	ng/uL	0.00
4) Pyridine-d5	3.822	84	584635	29.835	ng/ul	0.00
7) Phenol-d5	7.228	99	727494	30.400	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.375	67	490365	33.121	ng/ul	0.00
11) 2-Chlorophenol-d4	7.581	132	590939	32.713	ng/ul	0.00
15) 4-Methylphenol-d8	8.787	113	577330	30.539	ng/ul	0.00
21) Nitrobenzene-d5	9.240	128	277200	31.442	ng/ul	0.00
24) 2-Nitrophenol-d4	9.969	143	301358	29.288	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.534	165	486571	26.537	ng/ul	0.01
31) 4-Chloroaniline-d4	11.045	131	724057	30.740	ng/ul	0.00
46) Dimethylphthalate-d6	14.110	166	1607764	30.677	ng/ul	0.00
49) Acenaphthylene-d8	14.398	160	1851088	31.420	ng/ul	0.00
54) 4-Nitrophenol-d4	14.986	143	212821m	30.772	ng/ul	0.02
60) Fluorene-d10	15.698	176	1357974	31.468	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.869	200	170794	18.769	ng/ul	0.00
73) Anthracene-d10	17.569	188	2244863	33.213	ng/ul	0.00
81) Pyrene-d10	19.792	212	2652762	30.077	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.927	264	2615416	33.398	ng/ul	-0.01

Target Compounds Qvalue

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

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