

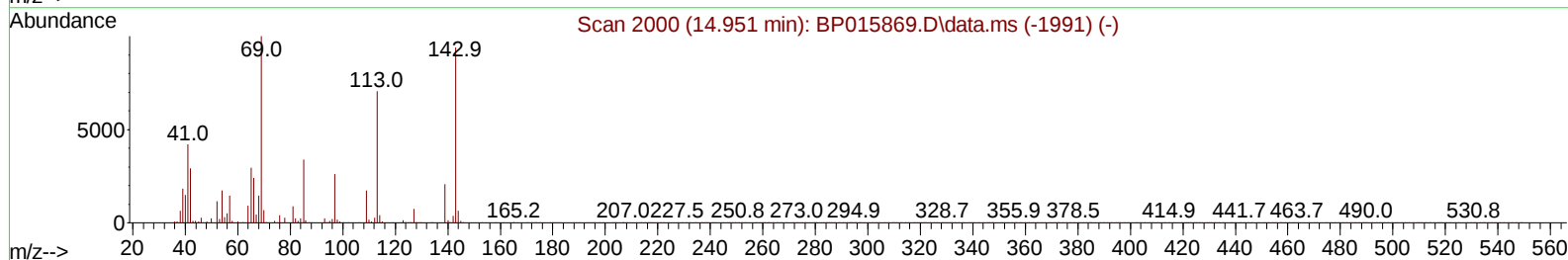
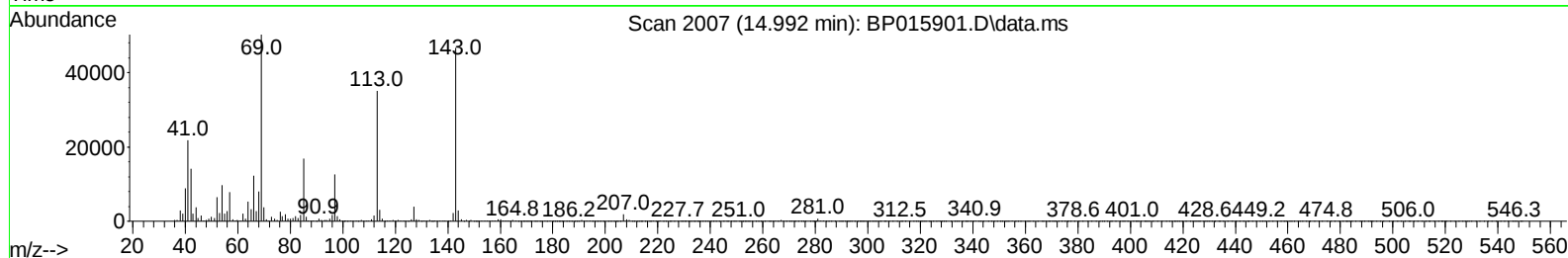
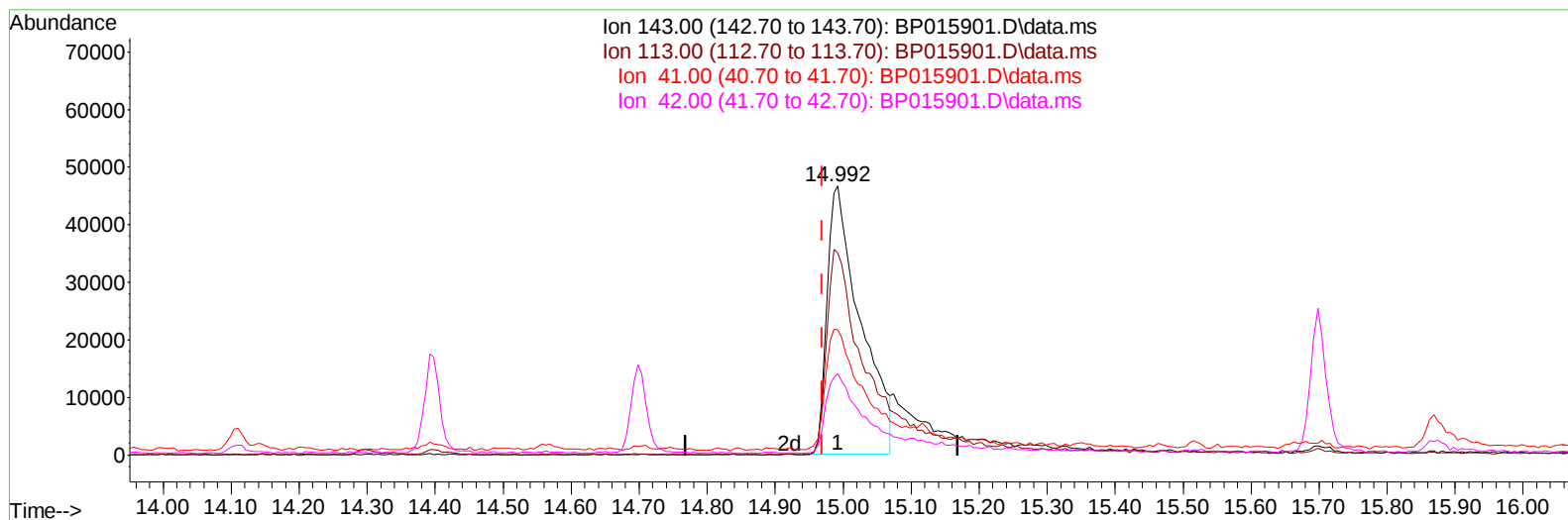
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
 Data File : BP015901.D
 Acq On : 02 Jul 2023 01:01
 Operator : MA/JU
 Sample : 03242-20
 Misc :
 ALS Vial : 69 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCGC1

Manual IntegrationsAPPROVED

Quant Time: Jul 02 02:48:00 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

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



TIC: BP015901.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.992min (+ 0.023) 23.89 ng/u1

response 158649

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.00	75.16#
41.00	45.70	46.54
42.00	31.20	30.29

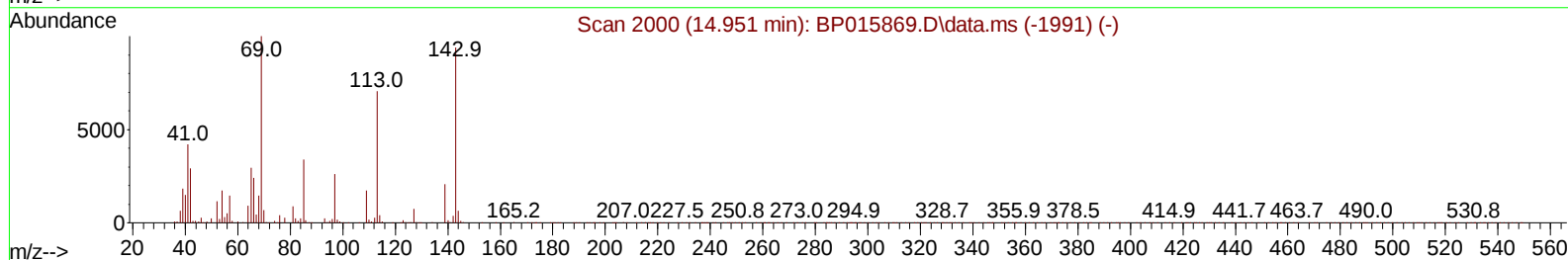
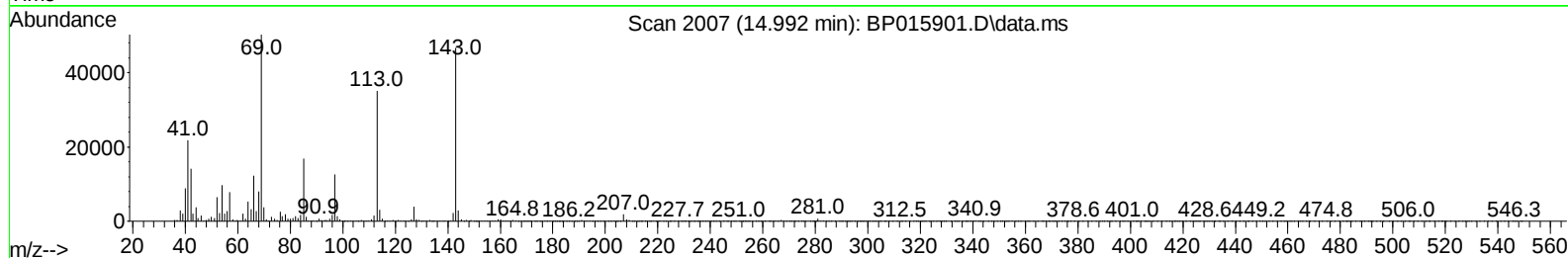
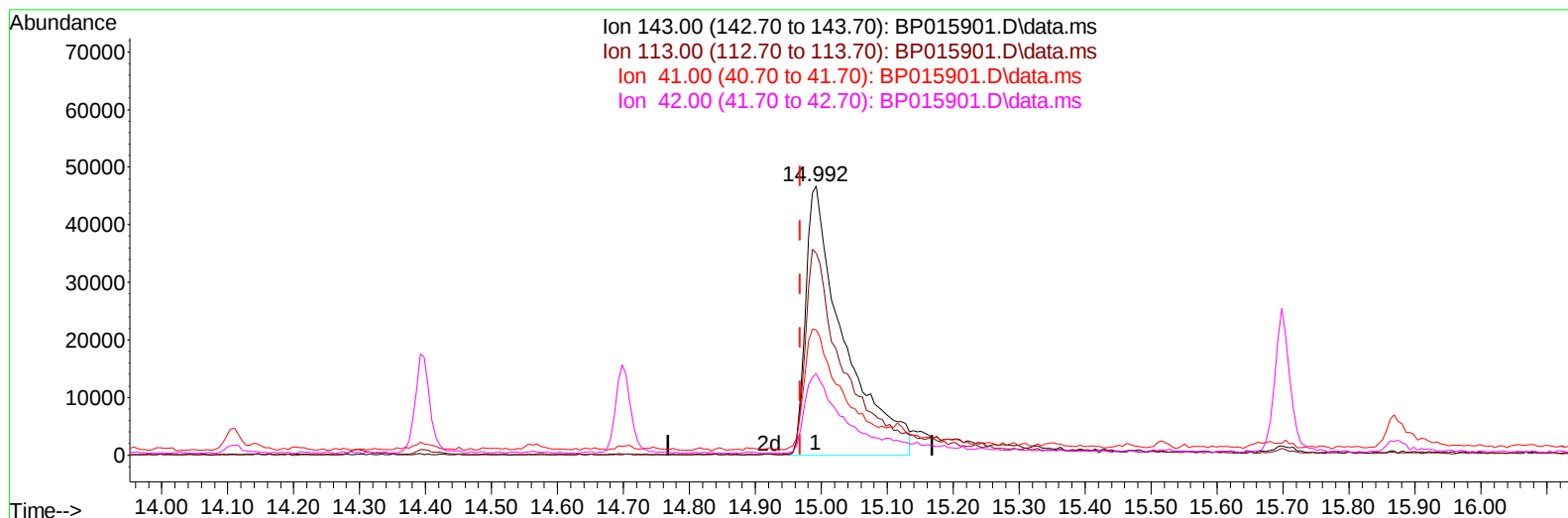
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TIC: BP015901.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.992min (+ 0.023) 28.12 ng/ul m

response 186717

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.00	75.16#
41.00	45.70	46.54
42.00	31.20	30.29

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Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/02/2023
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Quant Time: Jul 02 04:43:29 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
 Quant Title : SVOA CALIBRATION
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.051	152	252922	20.000	ng/ul	0.00
20) Naphthalene-d8	10.881	136	1117423	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.698	164	650971	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.463	188	1265790	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.557	240	756708	20.000	ng/ul	# 0.00
88) Perylene-d12	24.098	264	866150	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.375	96	34282	4.877	ng/uL	0.00
4) Pyridine-d5	3.828	84	221933	12.524	ng/ul	0.01
7) Phenol-d5	7.234	99	686856	31.740	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.375	67	443964	33.160	ng/ul	0.00
11) 2-Chlorophenol-d4	7.581	132	536046	32.814	ng/ul	0.00
15) 4-Methylphenol-d8	8.787	113	532703	31.160	ng/ul	0.00
21) Nitrobenzene-d5	9.240	128	268190	31.405	ng/ul	0.00
24) 2-Nitrophenol-d4	9.969	143	298581	29.957	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.534	165	518388	29.187	ng/ul	0.01
31) 4-Chloroaniline-d4	11.051	131	499288	21.883	ng/ul	0.01
46) Dimethylphthalate-d6	14.110	166	1662287	33.038	ng/ul	0.00
49) Acenaphthylene-d8	14.398	160	1880392	33.245	ng/ul	0.00
54) 4-Nitrophenol-d4	14.992	143	186717m	28.121	ng/ul	0.02
60) Fluorene-d10	15.698	176	1366090	32.974	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.869	200	128790	16.544	ng/ul	0.00
73) Anthracene-d10	17.563	188	1884003	32.585	ng/ul	0.00
81) Pyrene-d10	19.792	212	1876247	37.972	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.927	264	1538774	35.218	ng/ul	-0.01

Target Compounds Qvalue

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

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