

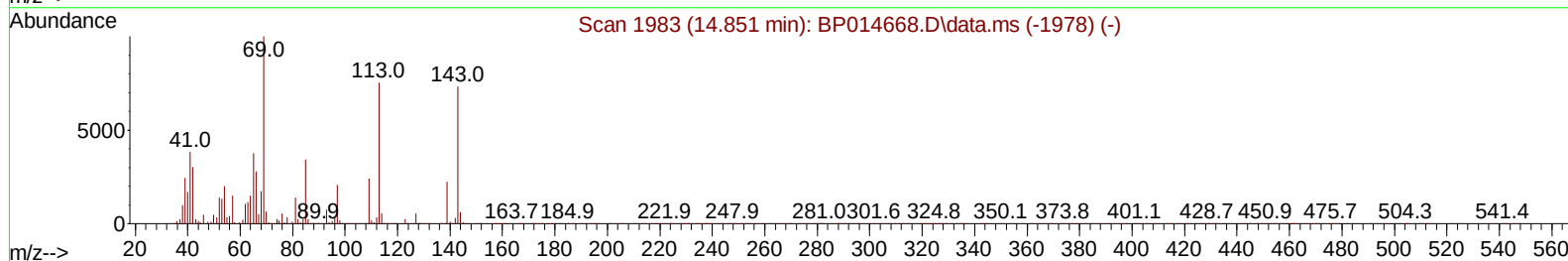
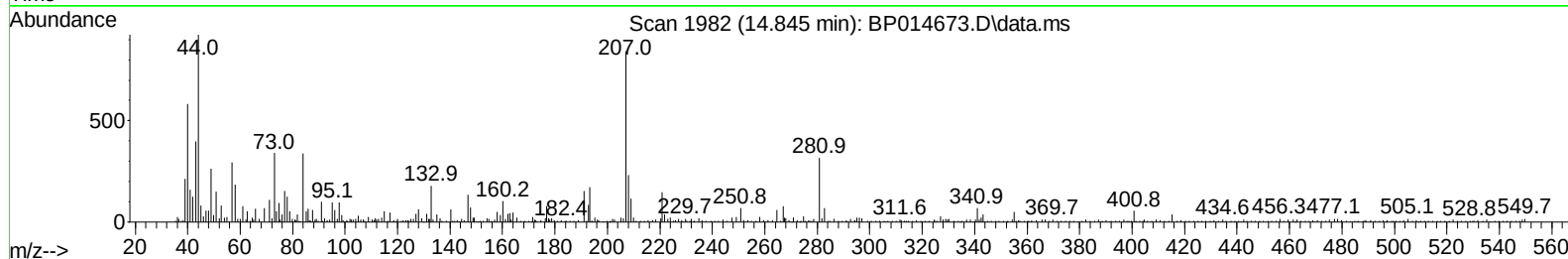
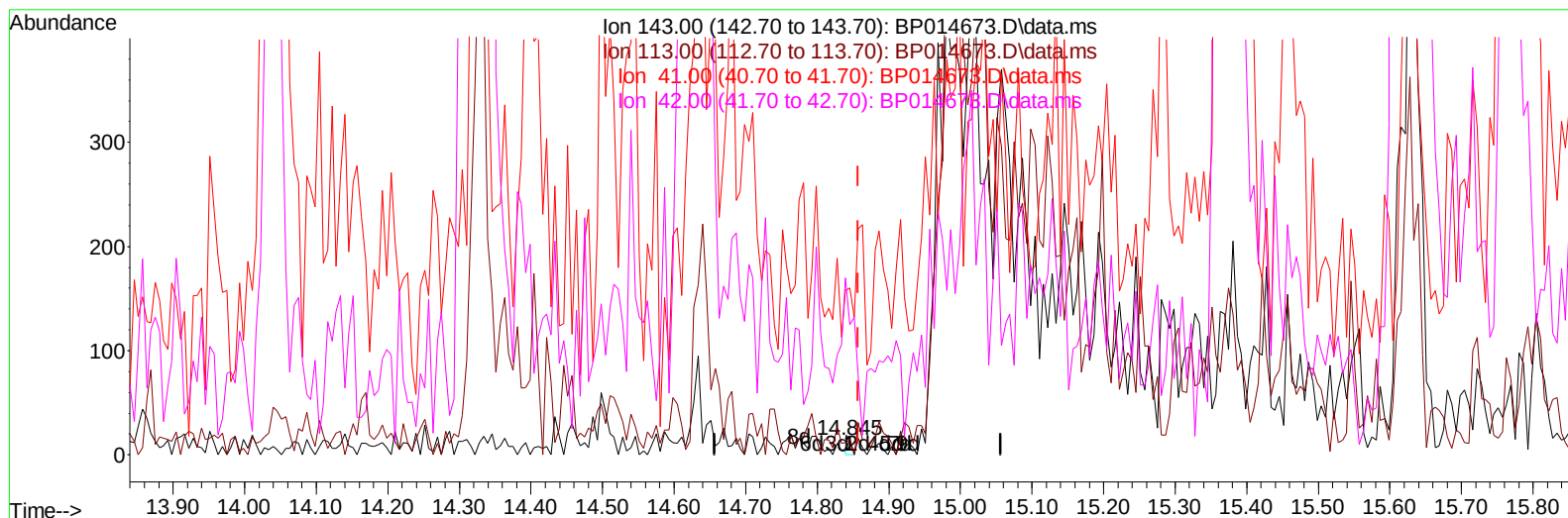
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041223\
Data File : BP014673.D
Acq On : 12 Apr 2023 14:56
Operator : CG/JU
Sample : 02251-15
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COMD3

Manual IntegrationsAPPROVED

Quant Time: Apr 13 00:49:30 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Apr 11 01:21:43 2023
Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/13/2023
Supervised By :Jagrut Upadhyay 04/13/2023



TIC: BP014673.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.845min (-0.012) 0.00 ng/u1

response 5

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	100.00
41.00	44.20	1066.67#
42.00	31.50	833.33#

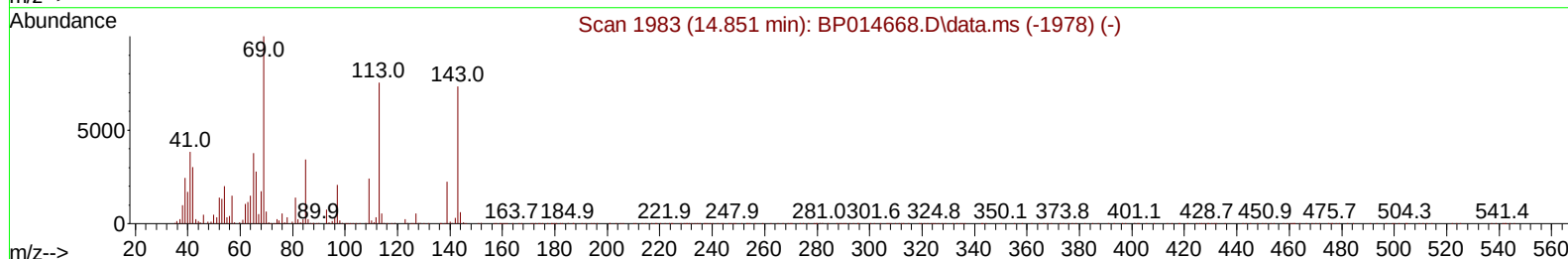
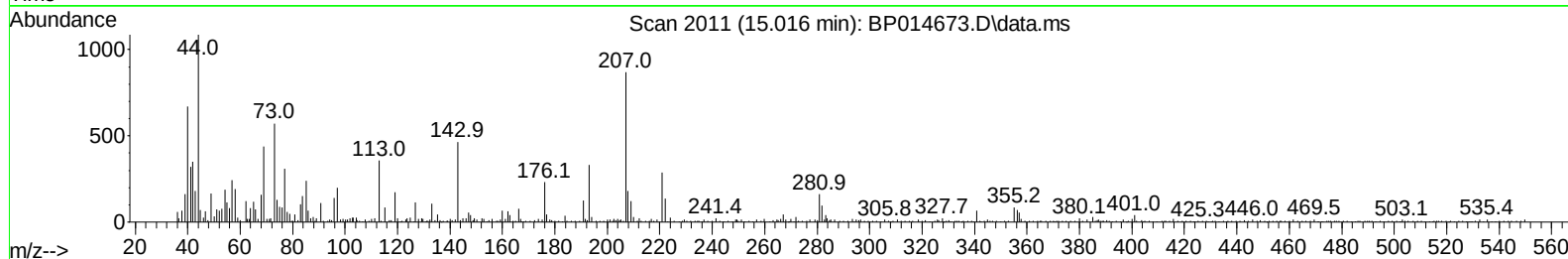
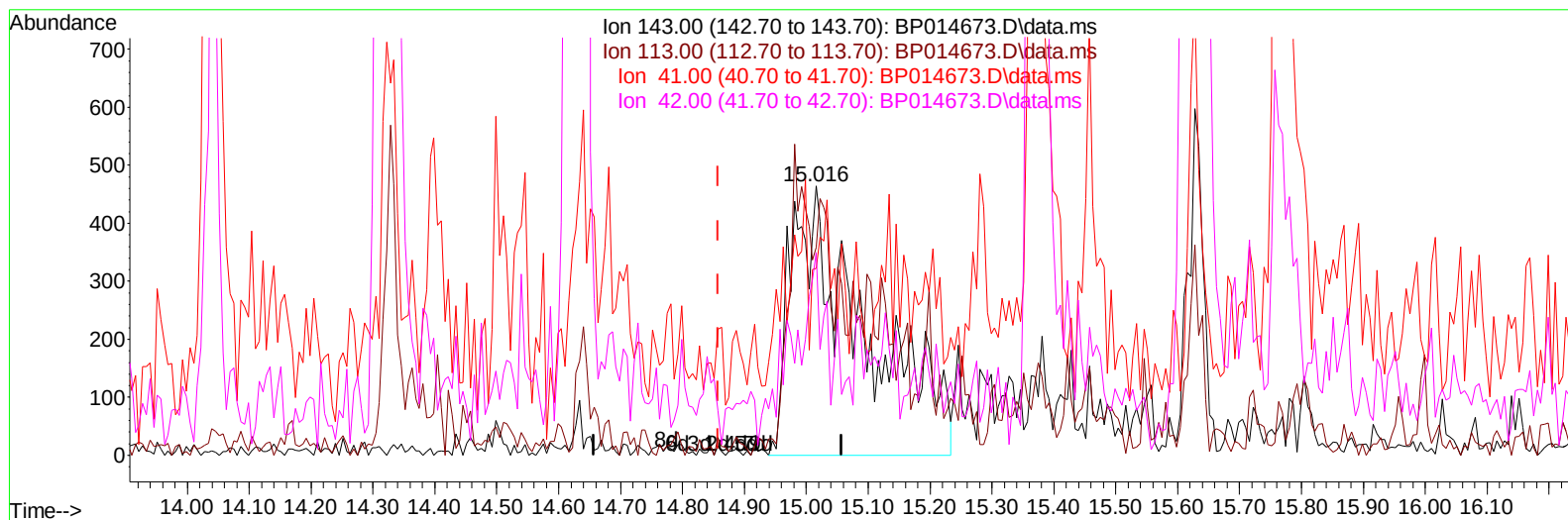
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(54) 4-Nitrophenol-d4 (S)

15.016min (+ 0.159) 1.23 ng/ul m

response 3821

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	77.16#
41.00	44.20	69.40#
42.00	31.50	75.43#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.981	152	110462	20.000	ng/ul	0.00
20) Naphthalene-d8	10.799	136	466659	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.628	164	254284	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.392	188	480432	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	322999	20.000	ng/ul	0.00
88) Perylene-d12	23.992	264	433992	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	12876	4.520	ng/uL	0.00
4) Pyridine-d5	3.805	84	74386	10.190	ng/ul	0.01
7) Phenol-d5	7.158	99	31140	3.057	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.305	67	190613	29.162	ng/ul	0.00
11) 2-Chlorophenol-d4	7.511	132	97555	13.249	ng/ul	0.00
15) 4-Methylphenol-d8	8.705	113	79097	9.626	ng/ul	0.00
21) Nitrobenzene-d5	9.158	128	93848	24.911	ng/ul	0.00
24) 2-Nitrophenol-d4	9.881	143	61226	14.193	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.428	165	100253	12.761	ng/ul	0.00
31) 4-Chloroaniline-d4	10.952	131	166410	16.081	ng/ul	0.00
46) Dimethylphthalate-d6	14.040	166	580909	28.175	ng/ul	0.00
49) Acenaphthylene-d8	14.322	160	630089	27.542	ng/ul	0.00
54) 4-Nitrophenol-d4	15.016	143	3821m	1.227	ng/ul	0.16
60) Fluorene-d10	15.628	176	472005	27.452	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.769	200	30453	8.614	ng/ul	0.01
73) Anthracene-d10	17.492	188	685105	29.466	ng/ul	0.00
81) Pyrene-d10	19.728	212	641917	29.587	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.821	264	701813	30.437	ng/ul	0.00

Target Compounds Qvalue

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

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