

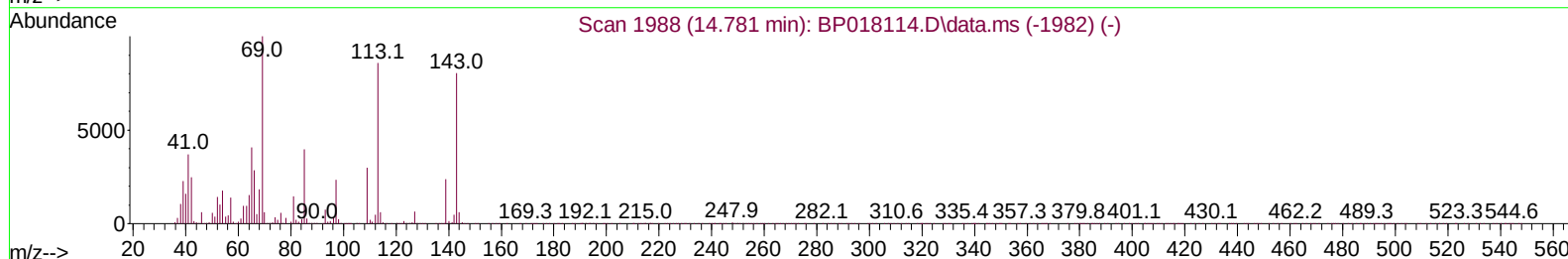
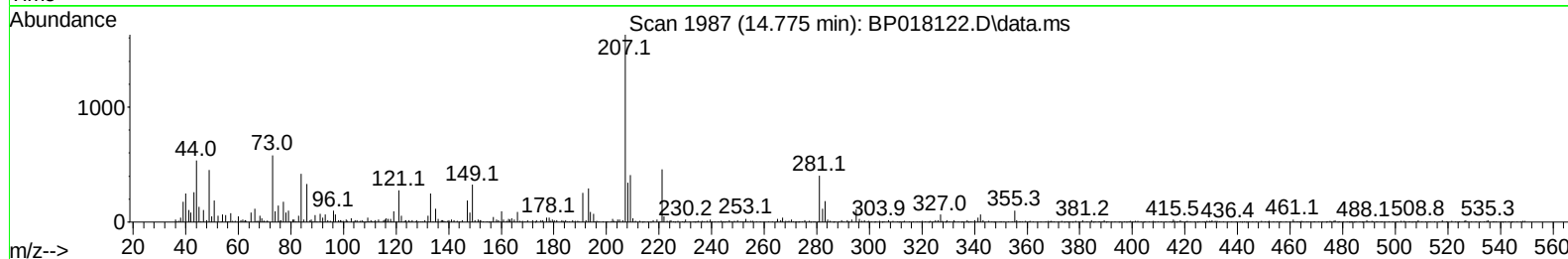
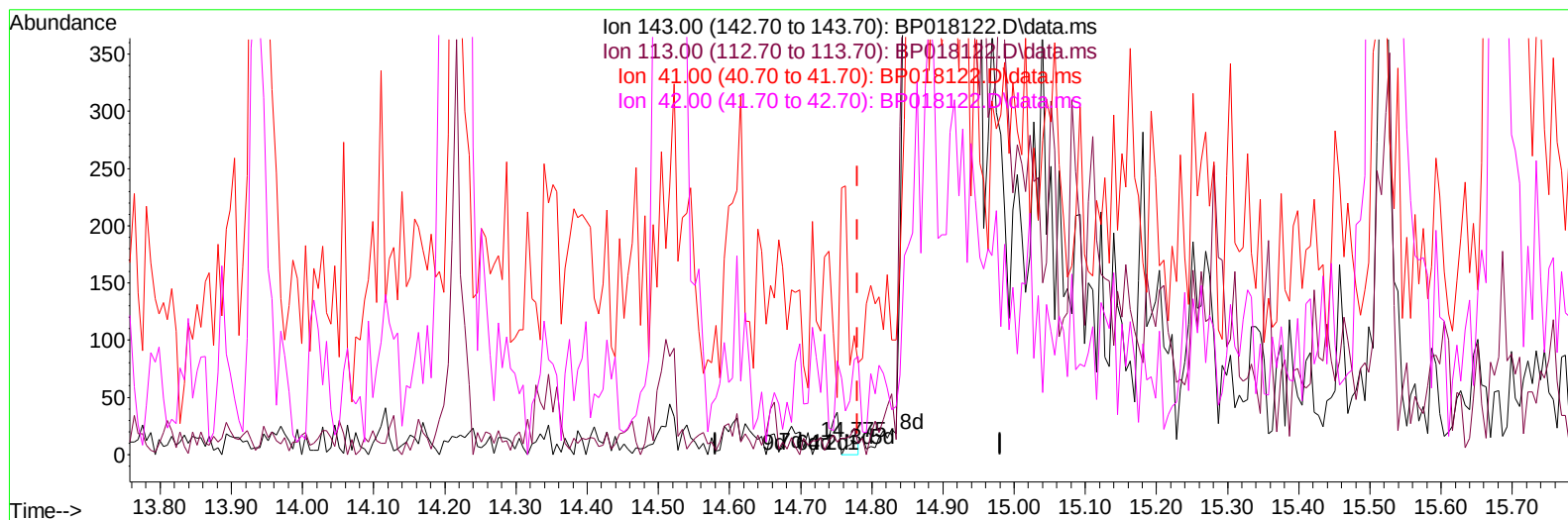
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
 Data File : BP018122.D
 Acq On : 13 Nov 2023 01:39
 Operator : MA/JU
 Sample : 05203-01 10X
 Misc :
 ALS Vial : 103 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AN8

Manual IntegrationsAPPROVED

Quant Time: Nov 13 03:37:19 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 10 21:55:18 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/14/2023
 Supervised By :mohammad ahmed 11/15/2023



TIC: BP018122.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.775min (-0.005) 0.01 ng/ul

response 13

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	191.67#
41.00	47.70	866.67#
42.00	30.10	691.67#

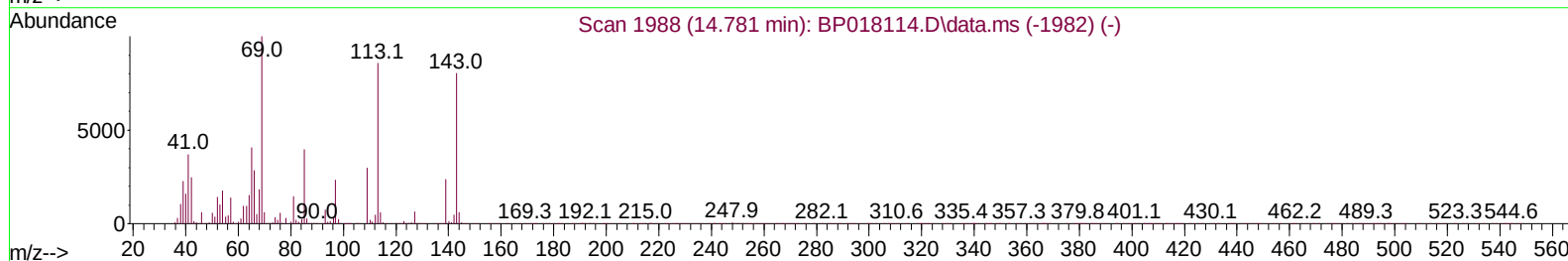
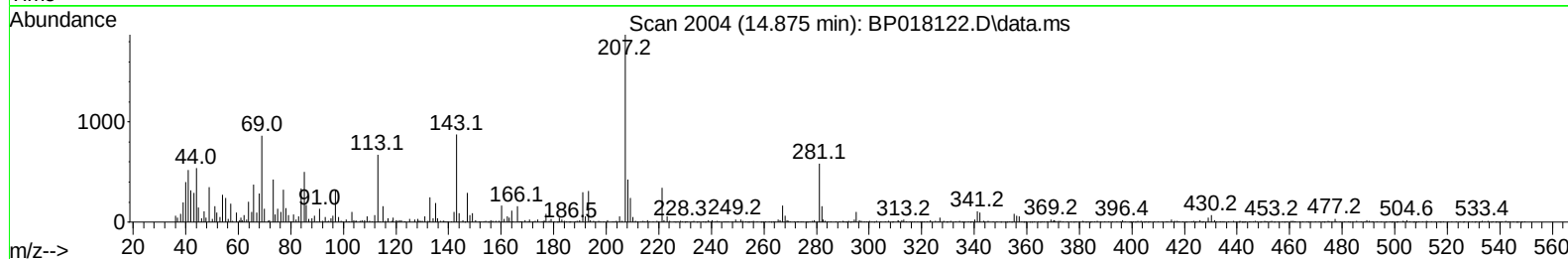
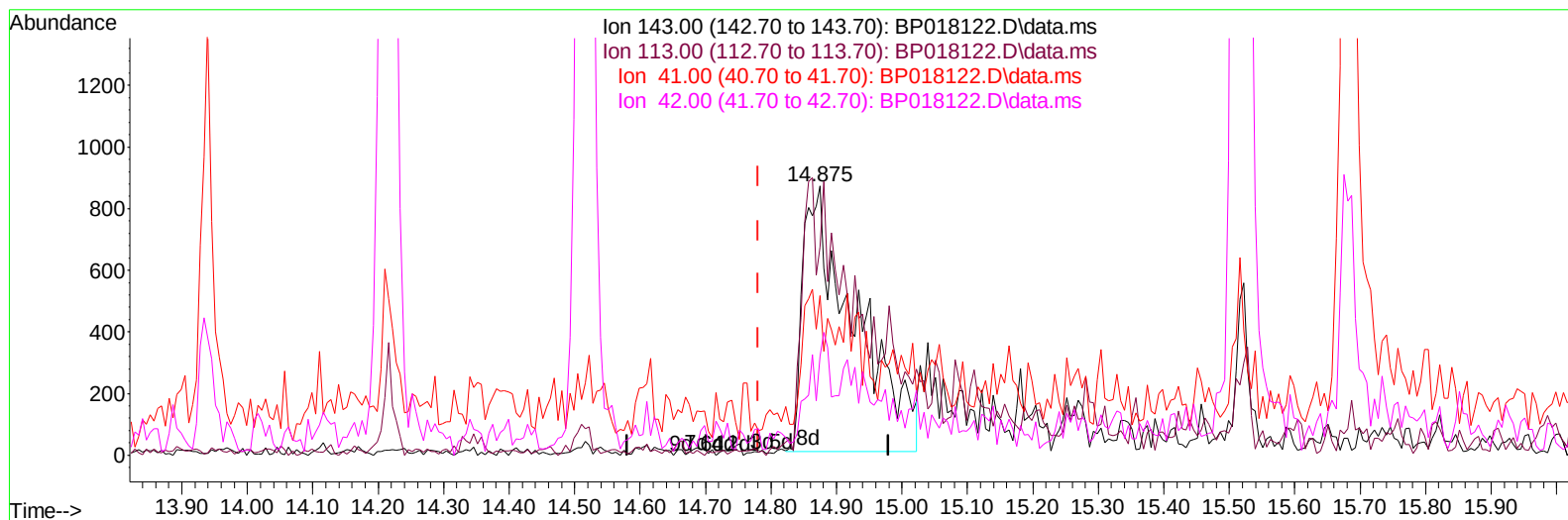
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TIC: BP018122.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.875min (+ 0.095) 2.40 ng/ul m

response 4795

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	77.43#
41.00	47.70	59.34#
42.00	30.10	36.66#

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Quant Time: Nov 13 03:38:35 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.840	152	58750	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.663	136	235886	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.516	164	145049	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.298	188	338829	20.000	ng/ul	-0.01
79) Chrysene-d12	21.428	240	337081	20.000	ng/ul	0.00
88) Perylene-d12	23.927	264	377465	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	2874	1.726	ng/uL	0.00
4) Pyridine-d5	3.676	84	30601	7.047	ng/ul	0.01
7) Phenol-d5	7.028	99	22118	3.900	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.193	67	88418	26.526	ng/ul	0.00
11) 2-Chlorophenol-d4	7.369	132	79013	17.591	ng/ul	0.00
15) 4-Methylphenol-d8	8.575	113	50025	10.763	ng/ul	0.00
21) Nitrobenzene-d5	9.052	128	54365	25.581	ng/ul	0.00
24) 2-Nitrophenol-d4	9.763	143	56844	23.610	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.293	165	84405	19.669	ng/ul	0.00
31) 4-Chloroaniline-d4	10.846	131	121318	20.616	ng/ul	0.00
46) Dimethylphthalate-d6	13.940	166	376782	28.153	ng/ul	-0.01
49) Acenaphthylene-d8	14.216	160	386961	26.546	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.875	143	4795m	2.398	ng/ul	0.09
60) Fluorene-d10	15.516	176	317776	28.759	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.681	200	47594	19.430	ng/ul	0.00
73) Anthracene-d10	17.398	188	551280	30.083	ng/ul	-0.01
81) Pyrene-d10	19.651	212	677586	30.094	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.763	264	670795	30.562	ng/ul	-0.01

Target Compounds Qvalue

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

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