

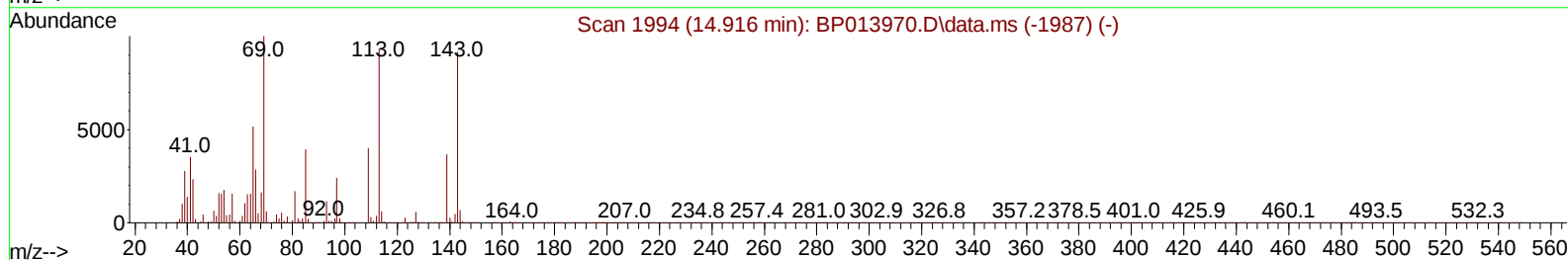
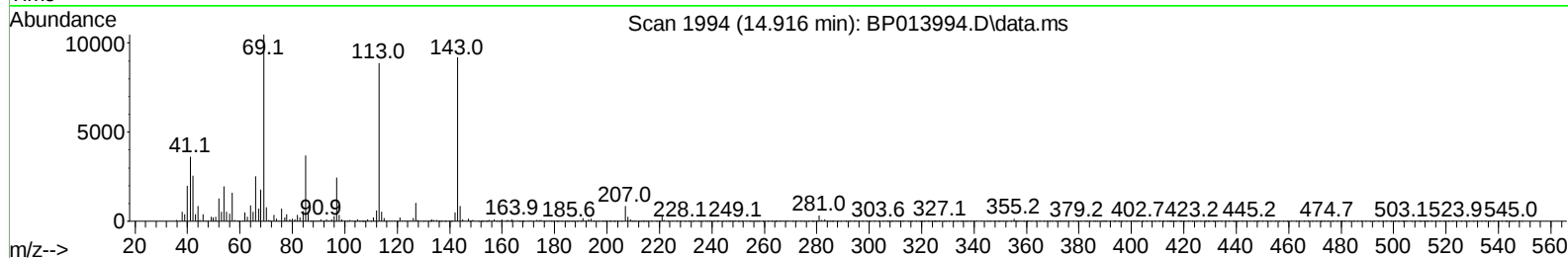
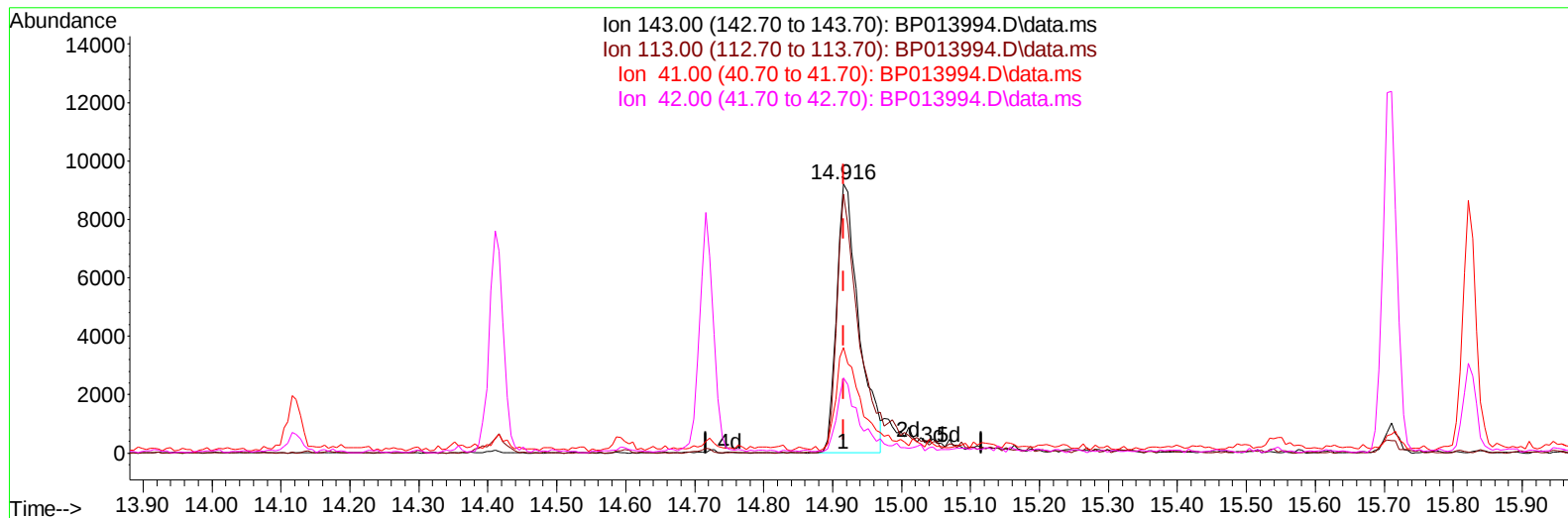
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030823\
Data File : BP013994.D
Acq On : 09 Mar 2023 00:35
Operator : CG/JU
Sample : 01729-11
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0BZ1

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 03/09/2023
Supervised By :Jagrut Upadhyay 03/09/2023

Quant Time: Mar 09 02:27:10 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Mar 09 01:10:12 2023
Response via : Initial Calibration



TIC: BP013994.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.916min (-0.000) 4.15 ng/u1

response 20796

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	96.43
41.00	40.20	39.33
42.00	28.50	27.84

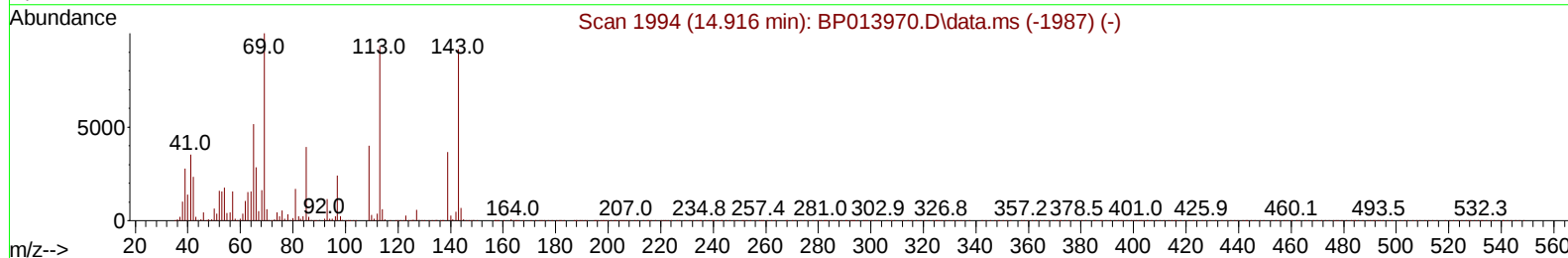
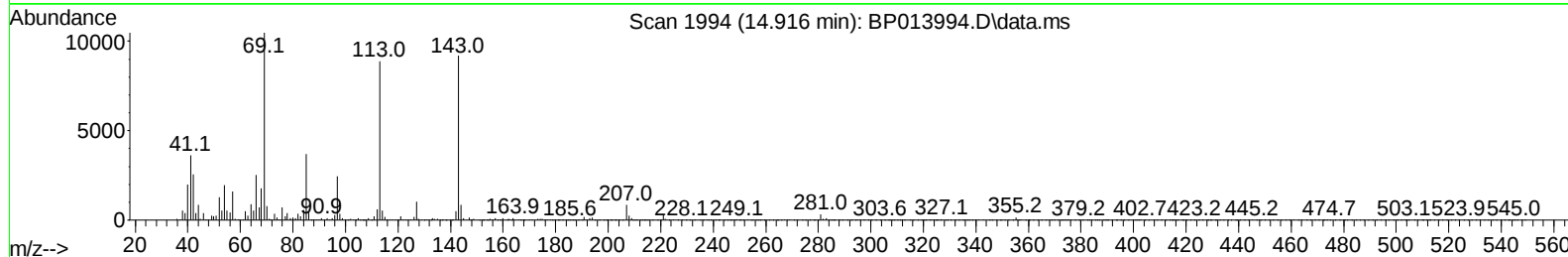
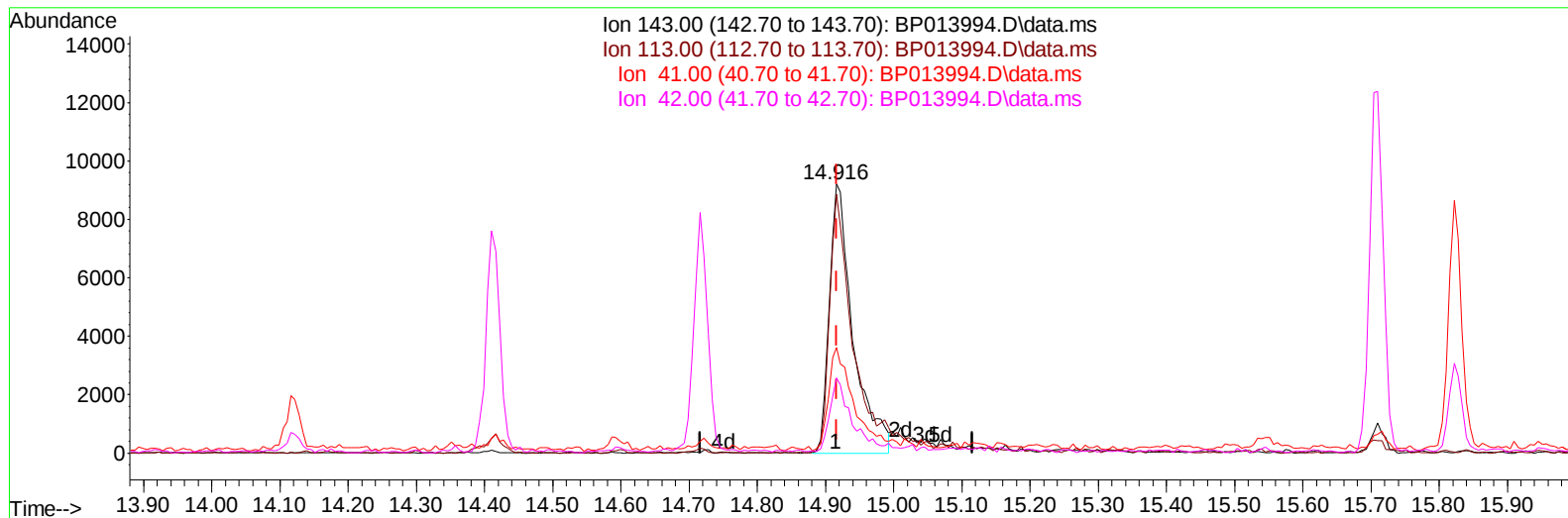
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TIC: BP013994.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.916min (-0.000) 4.43 ng/u1 m

response 22212

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	96.43
41.00	40.20	39.33
42.00	28.50	27.84

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.081	152	113266	20.000	ng/ul	0.00
20) Naphthalene-d8	10.905	136	461980	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.716	164	335104	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.469	188	824769	20.000	ng/ul	0.00
79) Chrysene-d12	21.545	240	904507	20.000	ng/ul	-0.01
88) Perylene-d12	24.086	264	1055658	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.434	96	12043	4.017	ng/uL	0.00
4) Pyridine-d5	3.875	84	69413	8.363	ng/ul	0.00
7) Phenol-d5	7.234	99	50106	5.013	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.399	67	155764	26.721	ng/ul	0.00
11) 2-Chlorophenol-d4	7.605	132	151316	19.870	ng/ul	0.00
15) 4-Methylphenol-d8	8.787	113	100137	12.265	ng/ul	0.00
21) Nitrobenzene-d5	9.252	128	101067	27.418	ng/ul	0.00
24) 2-Nitrophenol-d4	9.981	143	111968	25.375	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.522	165	189585	22.835	ng/ul	0.00
31) 4-Chloroaniline-d4	11.040	131	234658	21.357	ng/ul	0.00
46) Dimethylphthalate-d6	14.122	166	810143	28.487	ng/ul	0.00
49) Acenaphthylene-d8	14.410	160	820716	27.242	ng/ul	0.00
54) 4-Nitrophenol-d4	14.916	143	22212m	4.432	ng/ul	0.00
60) Fluorene-d10	15.710	176	696733	28.907	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.828	200	163109	26.166	ng/ul	0.00
73) Anthracene-d10	17.569	188	1278343	31.928	ng/ul	0.00
81) Pyrene-d10	19.792	212	1707022	34.218	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.921	264	1917549	34.149	ng/ul	0.00

Target Compounds Qvalue

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

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