

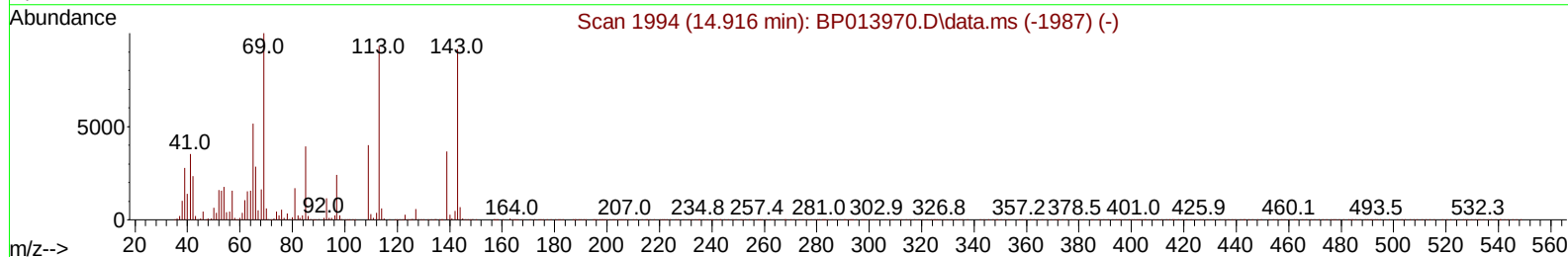
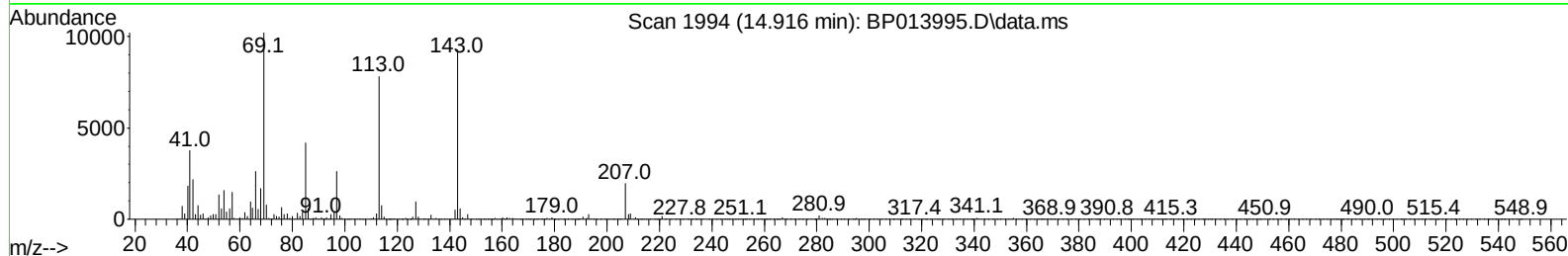
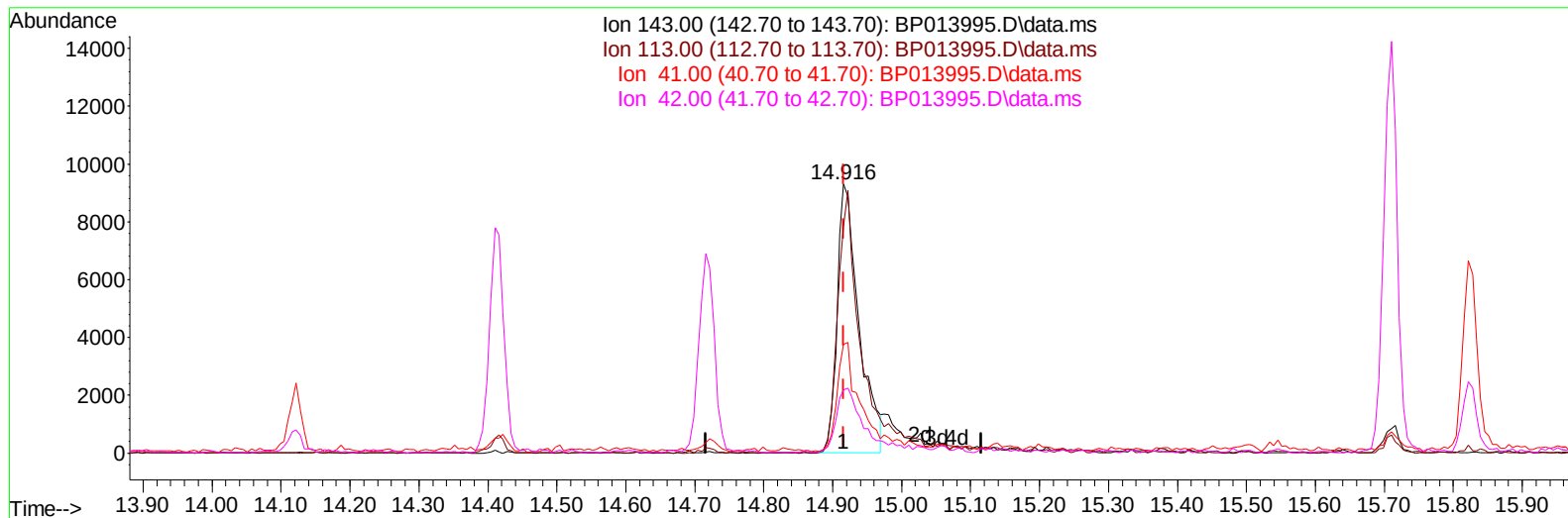
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030823\
Data File : BP013995.D
Acq On : 09 Mar 2023 01:09
Operator : CG/JU
Sample : 01729-12
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0BZ6

Manual IntegrationsAPPROVED

Quant Time: Mar 09 02:27:24 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

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



TIC: BP013995.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.916min (-0.000) 4.14 ng/u1

response 20426

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	84.24
41.00	40.20	40.46
42.00	28.50	23.45

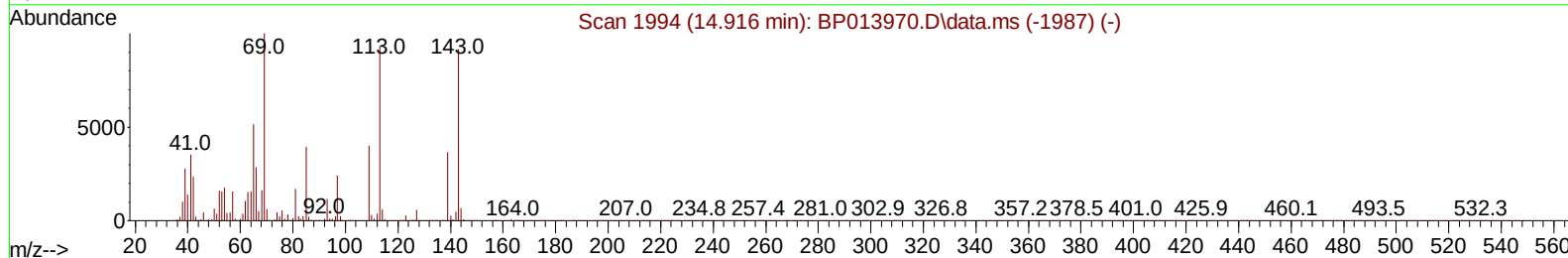
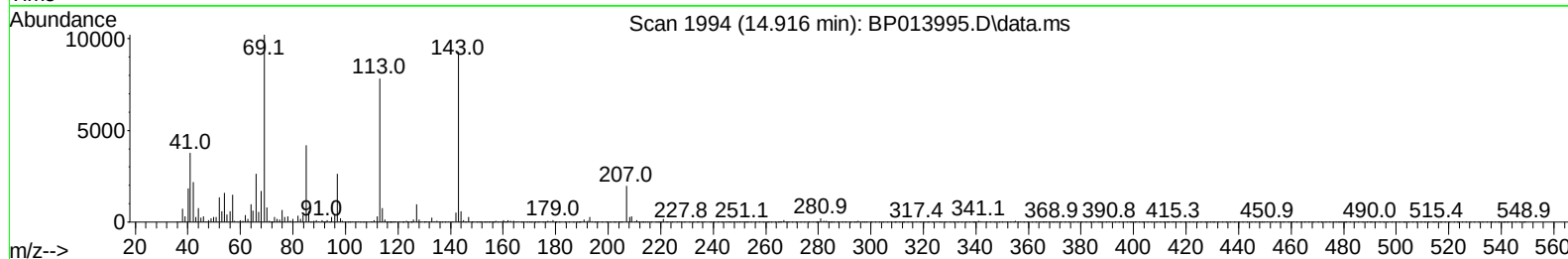
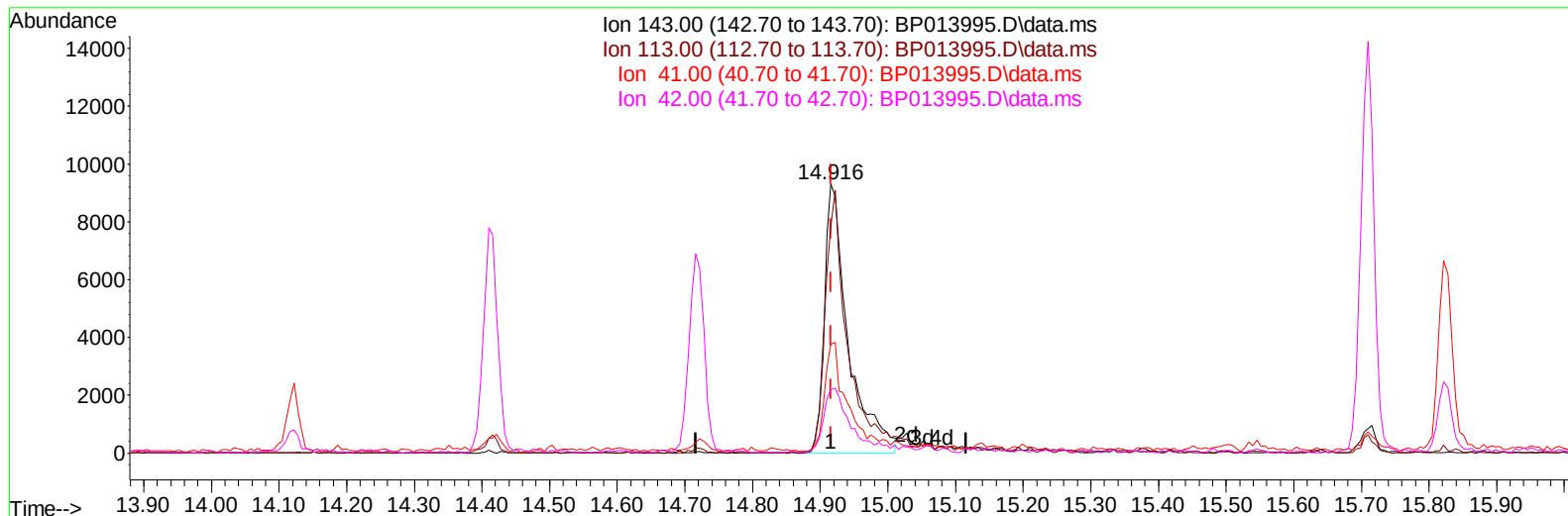
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TIC: BP013995.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.916min (-0.000) 4.59 ng/ul m

response 22692

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	84.24
41.00	40.20	40.46
42.00	28.50	23.45

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.081	152	116085	20.000	ng/ul	0.00
20) Naphthalene-d8	10.904	136	473062	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.716	164	330248	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.475	188	852978	20.000	ng/ul	0.00
79) Chrysene-d12	21.545	240	1015025	20.000	ng/ul	-0.01
88) Perylene-d12	24.086	264	1166253	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.440	96	14920	4.856	ng/uL	0.00
4) Pyridine-d5	3.875	84	84414	9.923	ng/ul	0.00
7) Phenol-d5	7.234	99	57631	5.626	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.399	67	173334	29.013	ng/ul	0.00
11) 2-Chlorophenol-d4	7.605	132	168931	21.644	ng/ul	0.00
15) 4-Methylphenol-d8	8.787	113	113047	13.510	ng/ul	0.00
21) Nitrobenzene-d5	9.252	128	111922	29.652	ng/ul	0.00
24) 2-Nitrophenol-d4	9.981	143	109090	24.144	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.522	165	210716	24.785	ng/ul	0.00
31) 4-Chloroaniline-d4	11.040	131	285037	25.335	ng/ul	0.00
46) Dimethylphthalate-d6	14.122	166	918155	32.759	ng/ul	0.00
49) Acenaphthylene-d8	14.416	160	904001	30.448	ng/ul	0.00
54) 4-Nitrophenol-d4	14.916	143	22692m	4.594	ng/ul	0.00
60) Fluorene-d10	15.710	176	784315	33.019	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.828	200	138335	21.458	ng/ul	0.00
73) Anthracene-d10	17.569	188	1418792	34.264	ng/ul	0.00
81) Pyrene-d10	19.792	212	1887470	33.716	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.921	264	2204697	35.539	ng/ul	0.00

Target Compounds Qvalue

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

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