

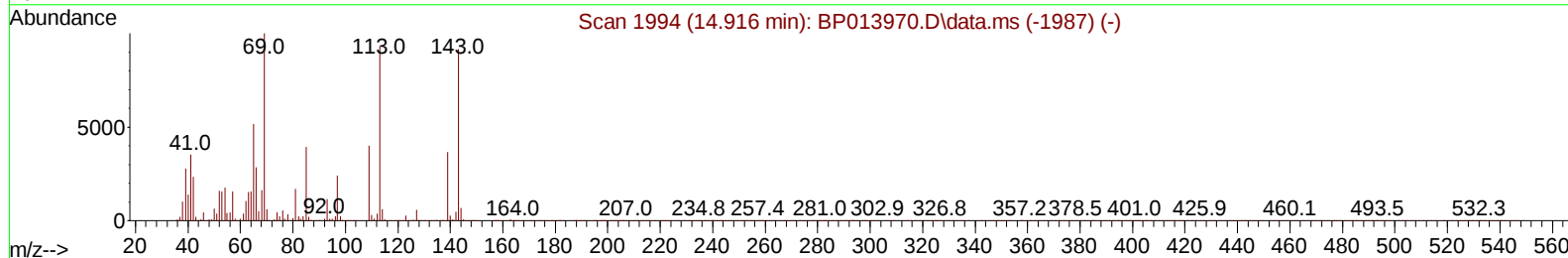
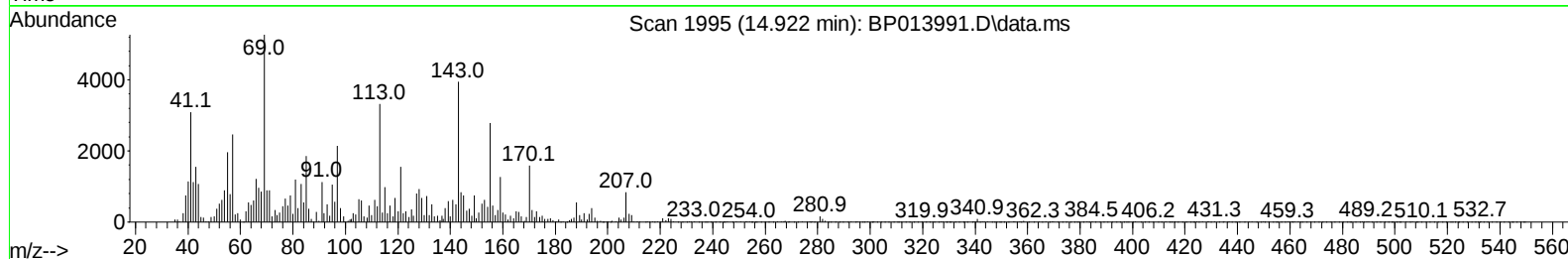
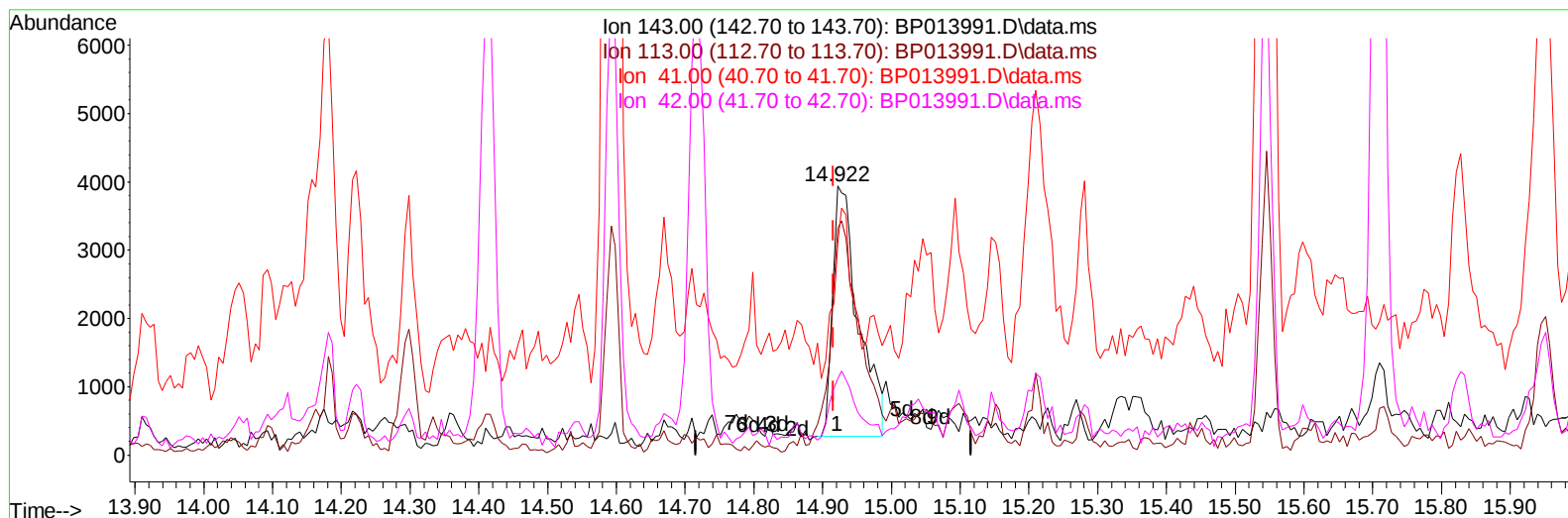
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030823\
Data File : BP013991.D
Acq On : 08 Mar 2023 22:53
Operator : CG/JU
Sample : 01729-06
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COBZ5

Quant Time: Mar 09 01:17:52 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

Manual IntegrationsAPPROVED

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



TIC: BP013991.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.922min (+ 0.006) 1.97 ng/u1

response 9571

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	84.41
41.00	40.20	78.57#
42.00	28.50	28.47

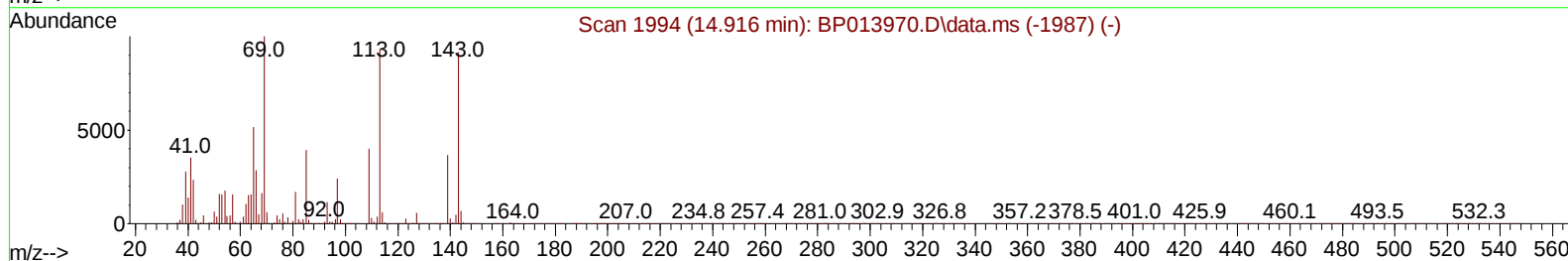
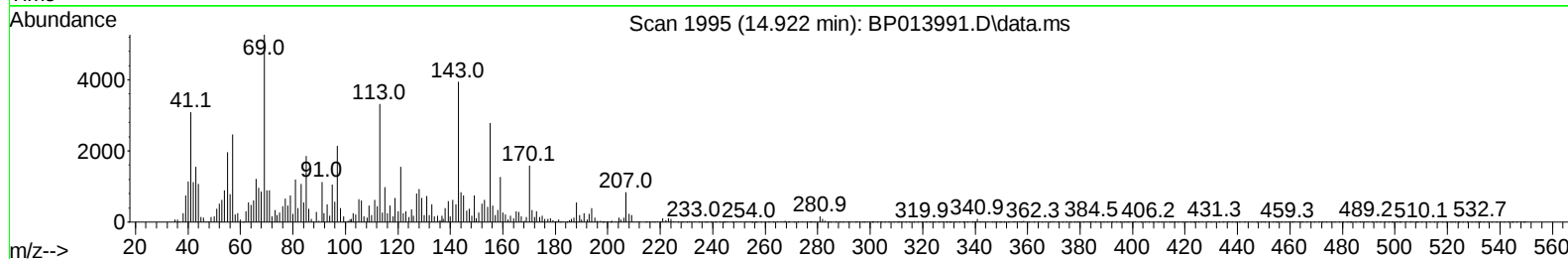
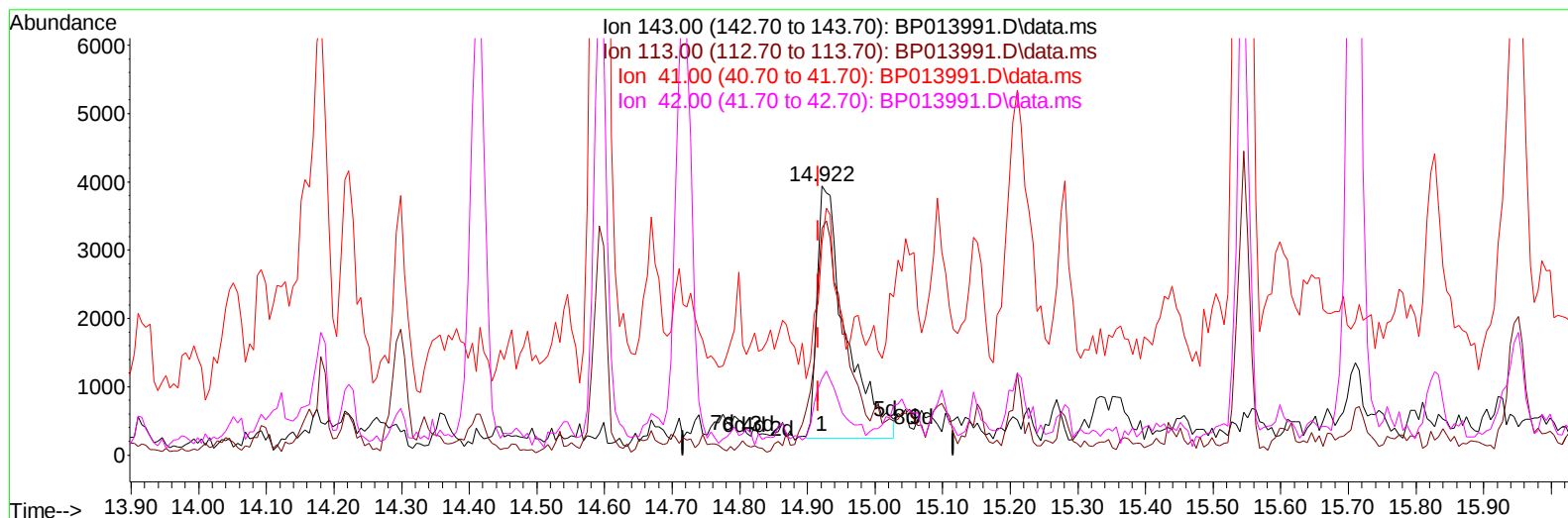
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(54) 4-Nitrophenol-d4 (S)

14.922min (+ 0.006) 2.23 ng/ul m

response 10863

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	84.41
41.00	40.20	78.57#
42.00	28.50	28.47

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Instrument :
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 C0BZ5

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.081	152	114773	20.000	ng/ul	0.00
20) Naphthalene-d8	10.904	136	454438	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.716	164	325200	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.475	188	816861	20.000	ng/ul	0.00
79) Chrysene-d12	21.551	240	987538	20.000	ng/ul	0.00
88) Perylene-d12	24.086	264	1131337	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.434	96	11103	3.655	ng/uL	0.00
4) Pyridine-d5	3.875	84	64359	7.652	ng/ul	0.00
7) Phenol-d5	7.234	99	32480	3.207	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.399	67	148443	25.130	ng/ul	0.00
11) 2-Chlorophenol-d4	7.610	132	96347	12.486	ng/ul	0.00
15) 4-Methylphenol-d8	8.787	113	72548	8.769	ng/ul	0.00
21) Nitrobenzene-d5	9.252	128	95185	26.251	ng/ul	0.00
24) 2-Nitrophenol-d4	9.981	143	67345	15.516	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.522	165	123012	15.062	ng/ul	0.00
31) 4-Chloroaniline-d4	11.040	131	184353	17.057	ng/ul	0.00
46) Dimethylphthalate-d6	14.122	166	763970	27.681	ng/ul	0.00
49) Acenaphthylene-d8	14.416	160	793421	27.138	ng/ul	0.00
54) 4-Nitrophenol-d4	14.922	143	10863m	2.234	ng/ul	0.00
60) Fluorene-d10	15.710	176	675096	28.862	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.828	200	53591	8.680	ng/ul	0.00
73) Anthracene-d10	17.575	188	1195395	30.145	ng/ul	0.00
81) Pyrene-d10	19.792	212	1642073	30.149	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.921	264	1836906	30.524	ng/ul	0.00

Target Compounds Qvalue

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

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