

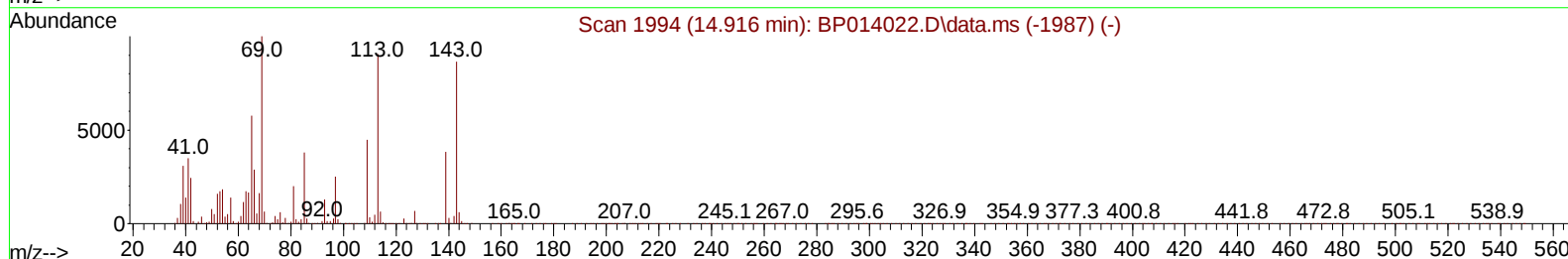
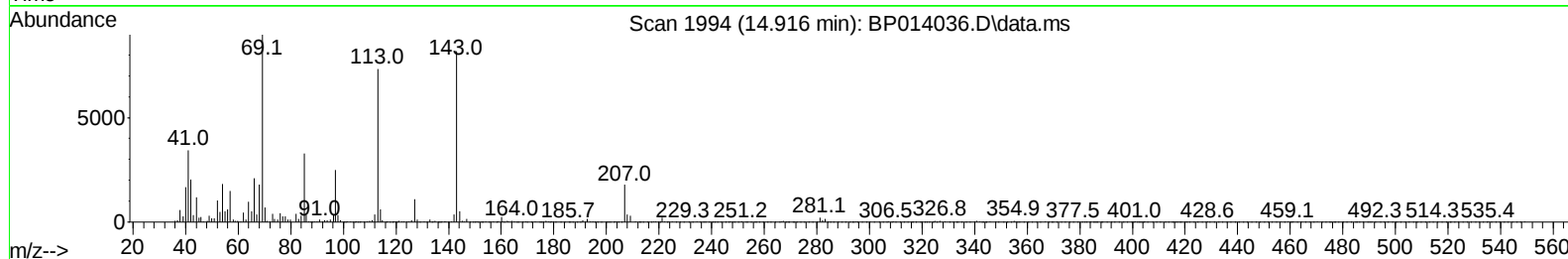
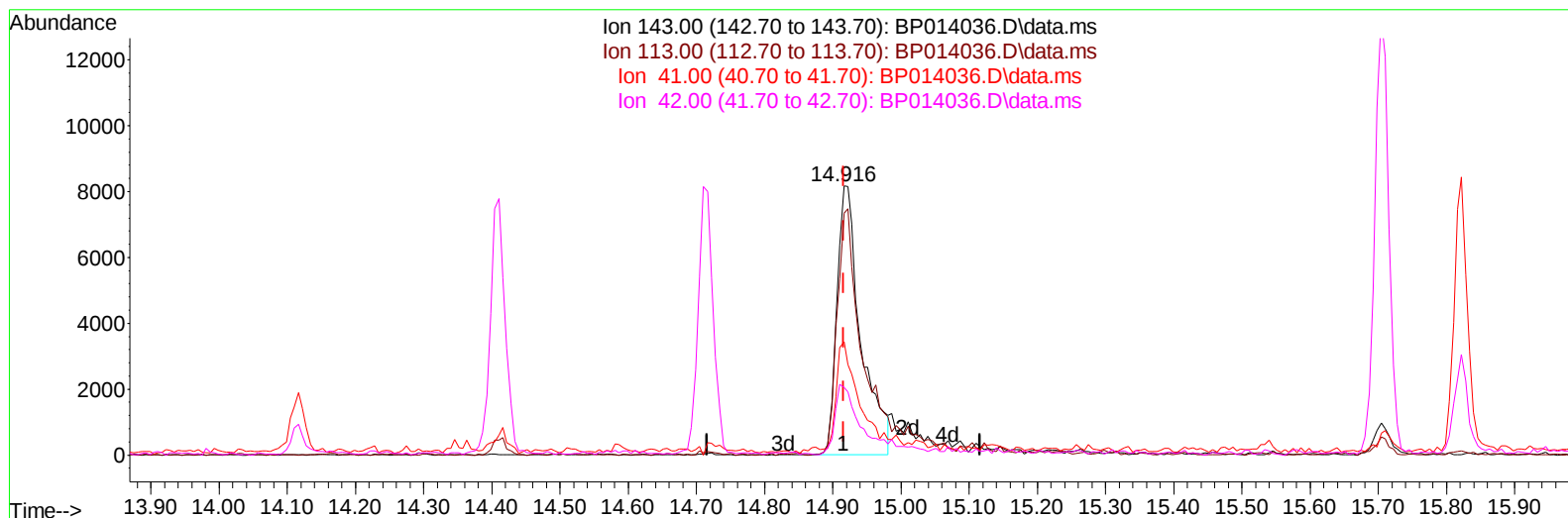
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031023\
 Data File : BP014036.D
 Acq On : 10 Mar 2023 18:22
 Operator : CG/JU
 Sample : 01804-04
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AA1

Manual IntegrationsAPPROVED

Quant Time: Mar 10 21:31:12 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 03/13/2023
 Supervised By :Jagrut Upadhyay 03/13/2023



TIC: BP014036.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.916min (-0.000) 3.56 ng/u1

response 20131

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	89.84
41.00	40.20	42.21
42.00	28.50	25.12

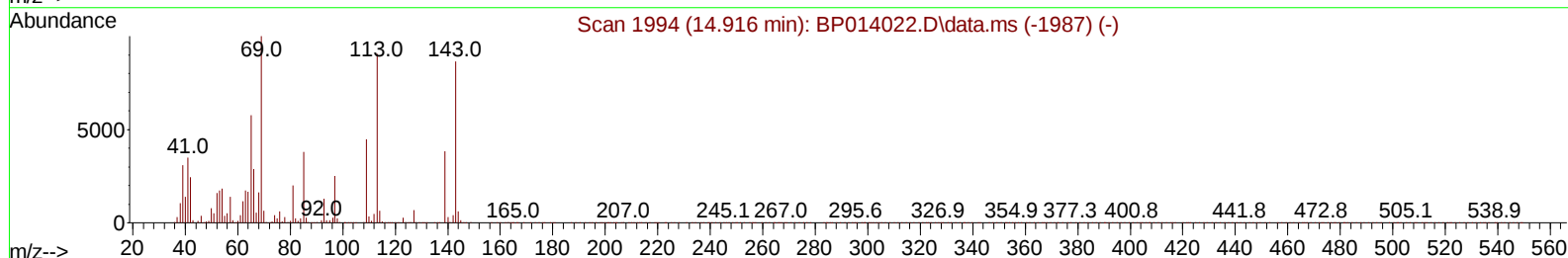
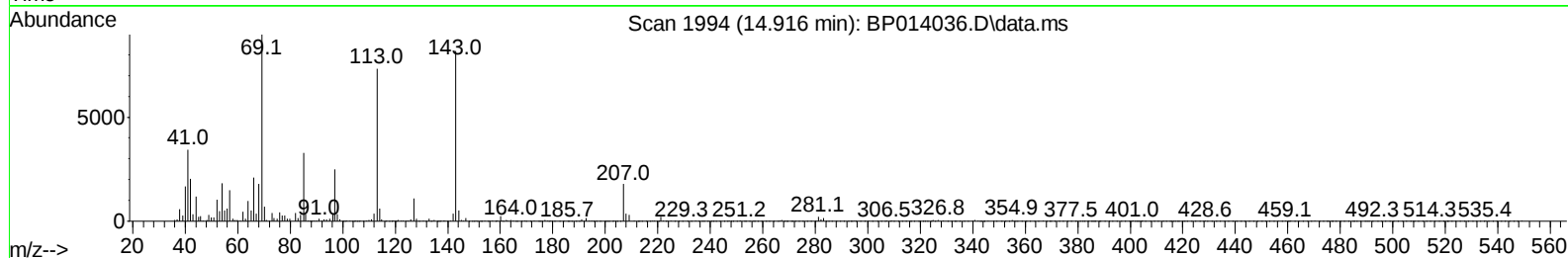
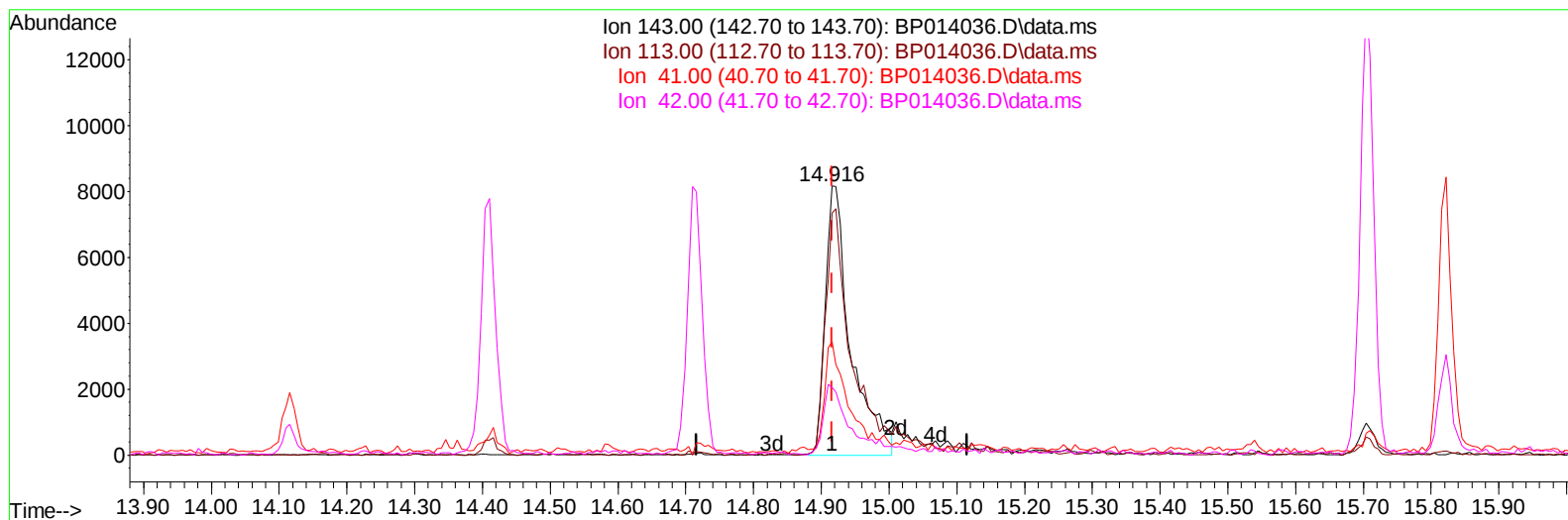
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TIC: BP014036.D\data.ms

(54) 4-Nitrophenol-d4 (S)**14.916min (-0.000) 3.78 ng/ul m****response 21374**

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	89.84
41.00	40.20	42.21
42.00	28.50	25.12

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.075	152	128209	20.000	ng/ul	0.00
20) Naphthalene-d8	10.899	136	551186	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.716	164	378520	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.469	188	909019	20.000	ng/ul	0.00
79) Chrysene-d12	21.545	240	973758	20.000	ng/ul	-0.01
88) Perylene-d12	24.086	264	1008751	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.434	96	12702	3.743	ng/uL	0.00
4) Pyridine-d5	3.875	84	64551	6.871	ng/ul	0.00
7) Phenol-d5	7.228	99	55505	4.906	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.399	67	179037	27.133	ng/ul	0.00
11) 2-Chlorophenol-d4	7.605	132	165774	19.231	ng/ul	0.00
15) 4-Methylphenol-d8	8.781	113	111636	12.079	ng/ul	-0.01
21) Nitrobenzene-d5	9.252	128	109618	24.925	ng/ul	0.00
24) 2-Nitrophenol-d4	9.975	143	122847	23.335	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.516	165	207676	20.965	ng/ul	0.00
31) 4-Chloroaniline-d4	11.040	131	252917	19.294	ng/ul	0.00
46) Dimethylphthalate-d6	14.116	166	823694	25.641	ng/ul	0.00
49) Acenaphthylene-d8	14.410	160	866029	25.449	ng/ul	0.00
54) 4-Nitrophenol-d4	14.916	143	21374m	3.776	ng/ul	0.00
60) Fluorene-d10	15.704	176	727904	26.736	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.822	200	157414	22.912	ng/ul	-0.01
73) Anthracene-d10	17.569	188	1330799	30.157	ng/ul	0.00
81) Pyrene-d10	19.792	212	1710788	31.855	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.916	264	1680136	31.312	ng/ul	-0.01

Target Compounds Qvalue

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

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