

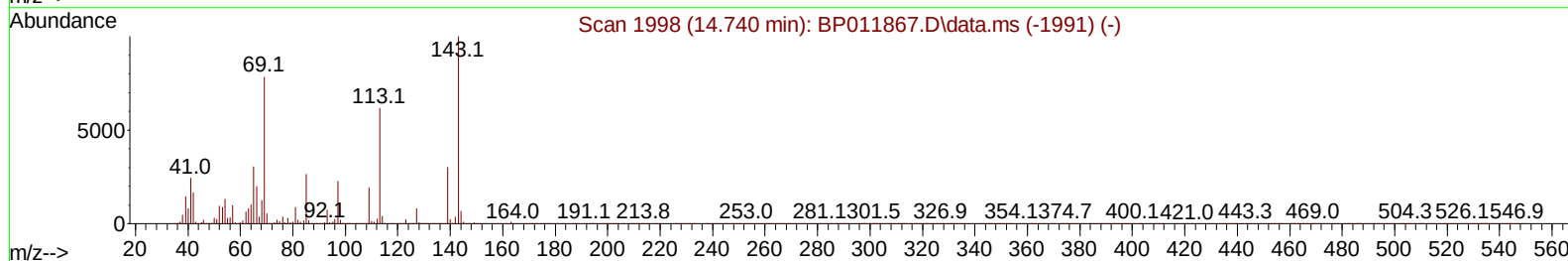
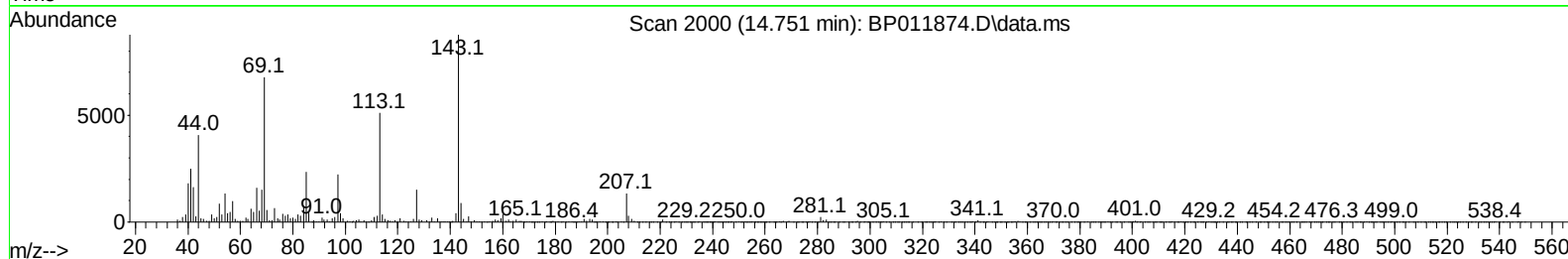
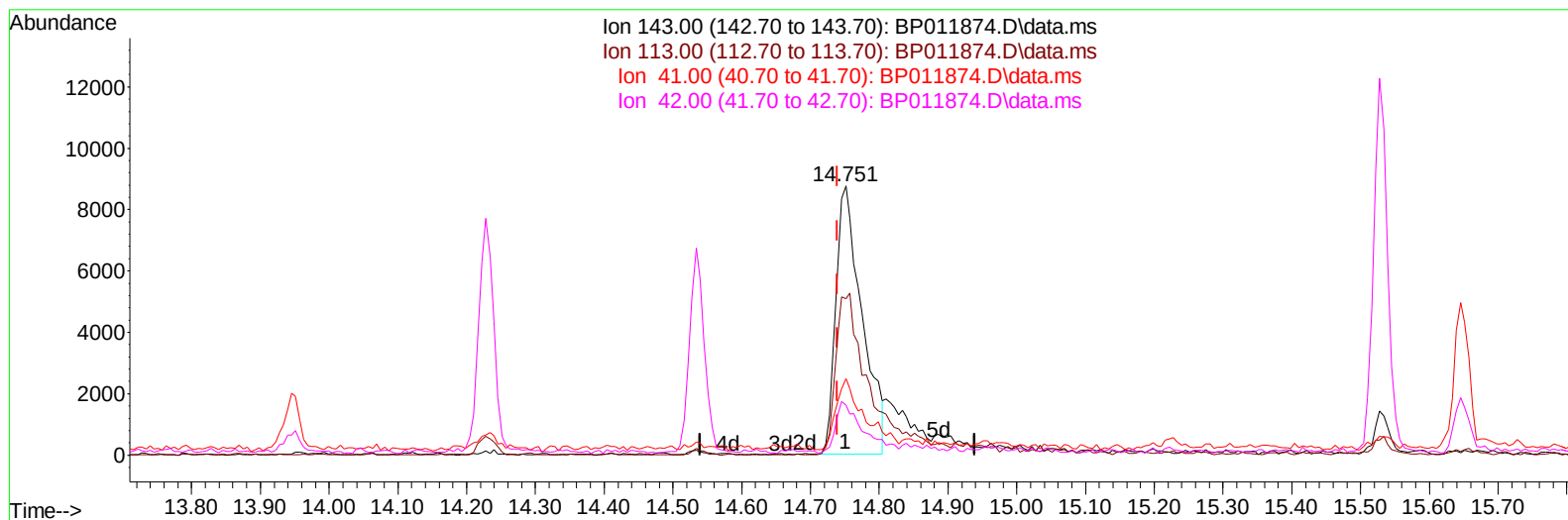
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100322\
 Data File : BP011874.D
 Acq On : 03 Oct 2022 16:15
 Operator : CG/JU
 Sample : N4859-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E10019

Manual IntegrationsAPPROVED

Quant Time: Oct 03 23:32:01 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 01 00:53:59 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/04/2022
 Supervised By :mohammad ahmed 10/06/2022



TIC: BP011874.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.751min (+ 0.012) 2.65 ng/ul

response 22978

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.80	58.15
41.00	24.20	28.32
42.00	16.40	18.49

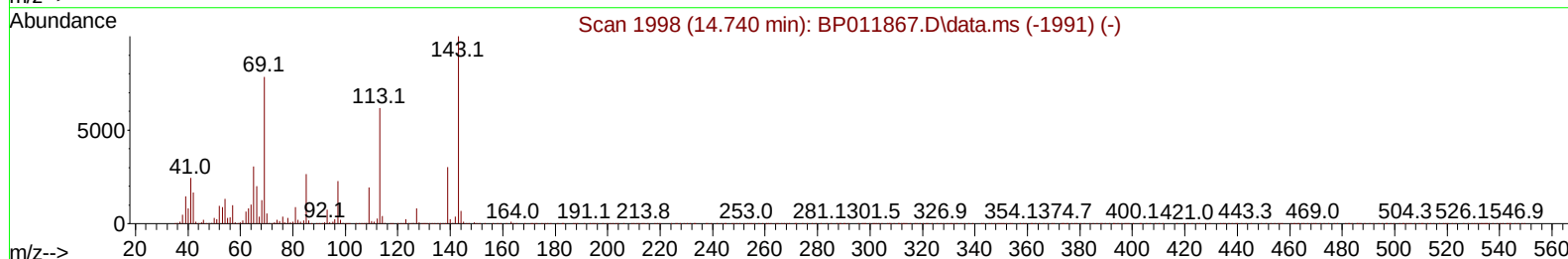
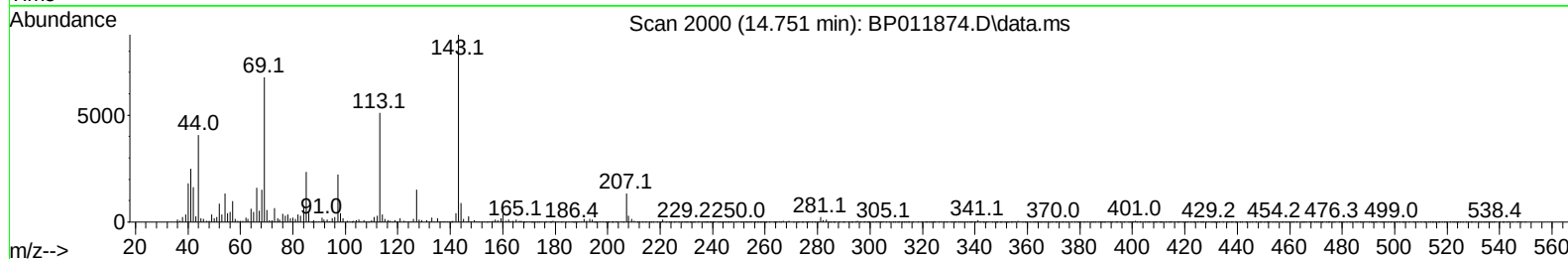
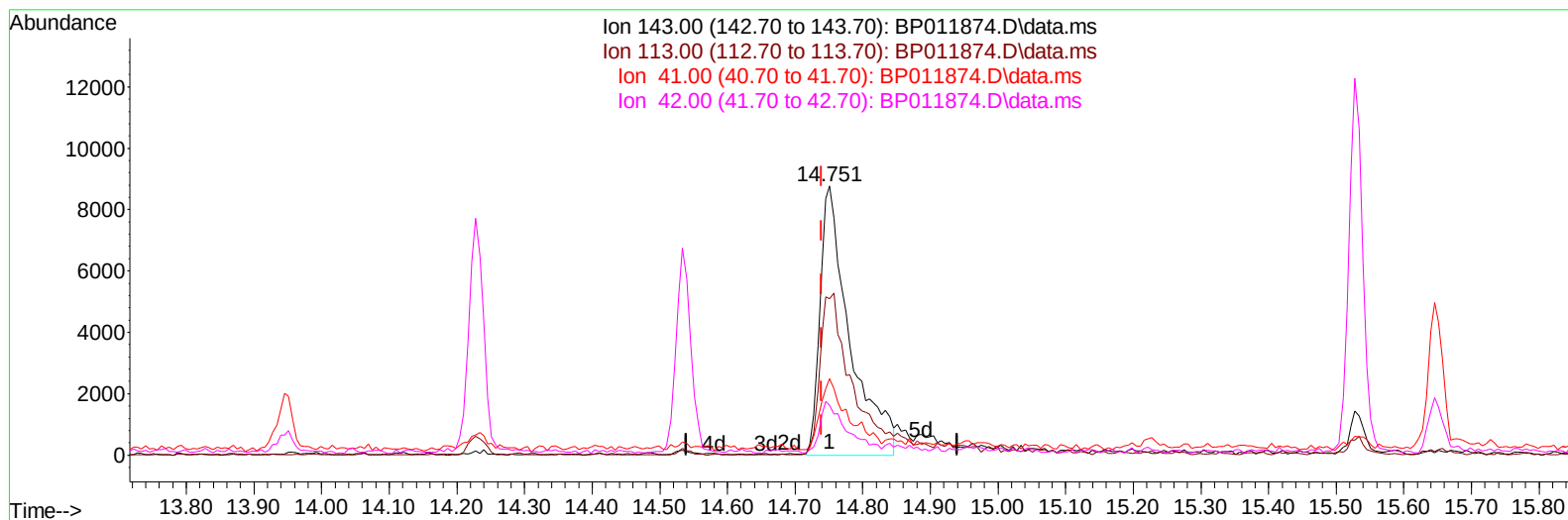
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TIC: BP011874.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.751min (+ 0.012) 3.08 ng/ul m

response 26698

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.80	58.15
41.00	24.20	28.32
42.00	16.40	18.49

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.893	152	256854	20.000	ng/ul	0.00
20) Naphthalene-d8	10.699	136	1032411	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.534	164	585415	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.292	188	1179815	20.000	ng/ul	0.00
79) Chrysene-d12	21.374	240	1294836	20.000	ng/ul	-0.01
88) Perylene-d12	23.810	264	1366497	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	32024	5.499	ng/uL	0.00
4) Pyridine-d5	3.734	84	74585	4.372	ng/ul	0.00
7) Phenol-d5	7.052	99	134022	6.044	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	381744	31.059	ng/ul	0.00
11) 2-Chlorophenol-d4	7.422	132	431825	24.669	ng/ul	0.00
15) 4-Methylphenol-d8	8.599	113	247984	13.529	ng/ul	0.00
21) Nitrobenzene-d5	9.057	128	270510	34.400	ng/ul	0.00
24) 2-Nitrophenol-d4	9.781	143	241237	28.673	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.322	165	433269	27.842	ng/ul	0.00
31) 4-Chloroaniline-d4	10.840	131	501173	21.125	ng/ul	0.00
46) Dimethylphthalate-d6	13.945	166	1403360	33.522	ng/ul	0.00
49) Acenaphthylene-d8	14.228	160	1645242	34.474	ng/ul	0.00
54) 4-Nitrophenol-d4	14.751	143	26698m	3.080	ng/ul	0.01
60) Fluorene-d10	15.528	176	1250700	34.402	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200	128886	18.049	ng/ul	0.00
73) Anthracene-d10	17.392	188	2058433	38.301	ng/ul	0.00
81) Pyrene-d10	19.628	212	2516751	38.379	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.651	264	2744181	40.116	ng/ul	-0.01

Target Compounds Qvalue

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

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