

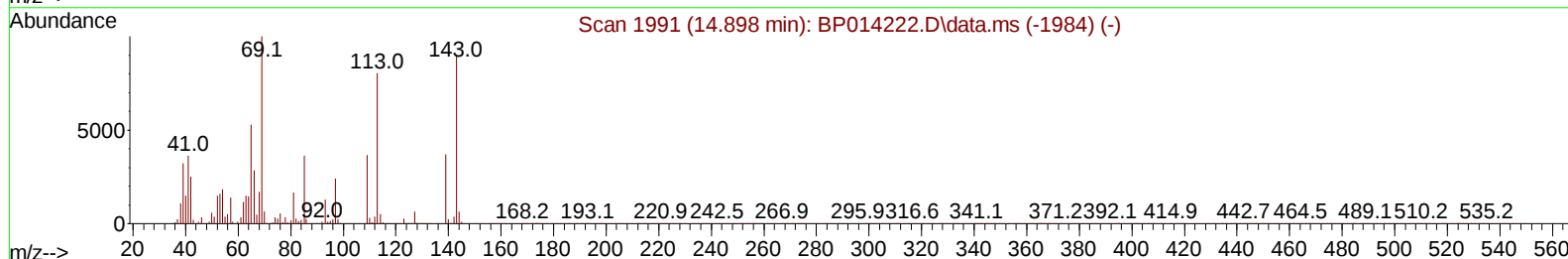
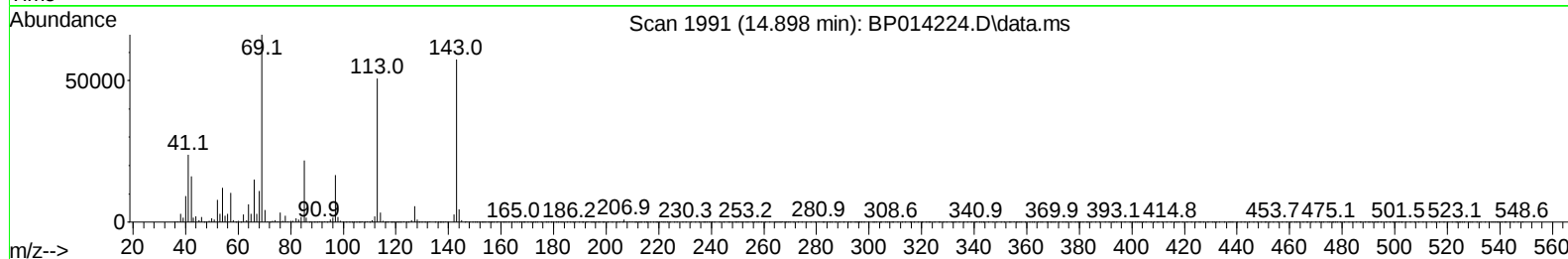
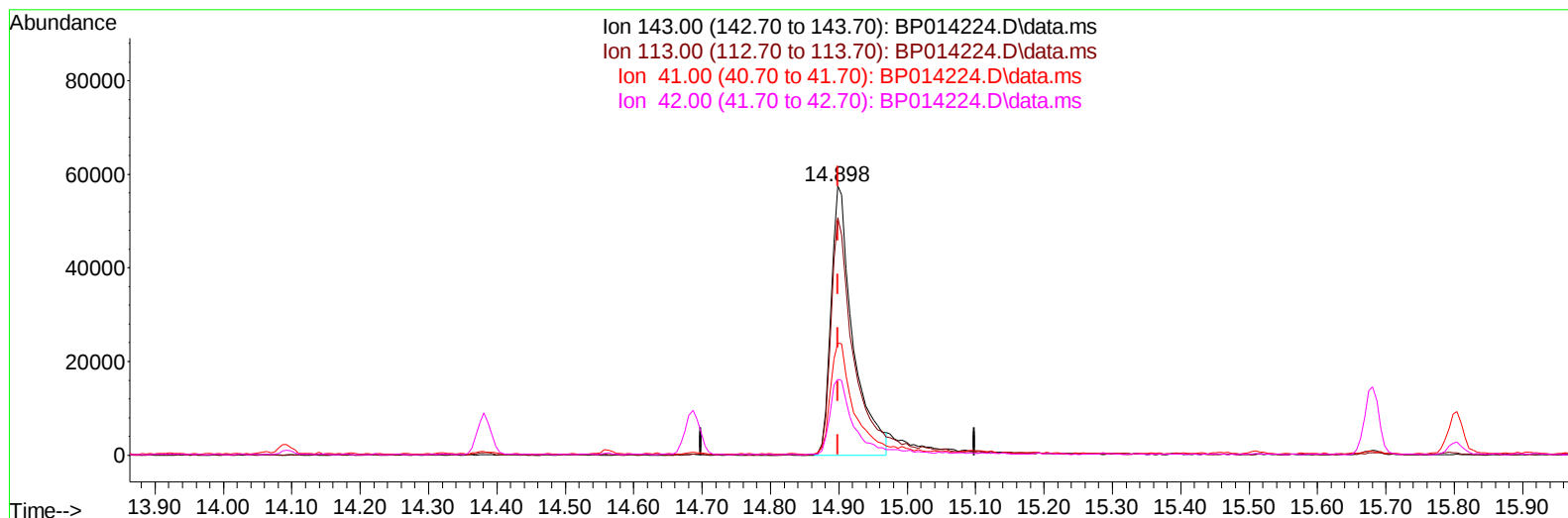
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032223\
Data File : BP014224.D
Acq On : 22 Mar 2023 19:20
Operator : CG/JU
Sample : 01977-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EYG87

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 03/23/2023
Supervised By :Jagrut Upadhyay 03/23/2023

Quant Time: Mar 23 02:35:26 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Mar 23 02:31:48 2023
Response via : Initial Calibration



TIC: BP014224.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.898min (-0.000) 18.07 ng/u1

response 127667

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	88.41
41.00	40.20	41.58
42.00	28.50	28.31

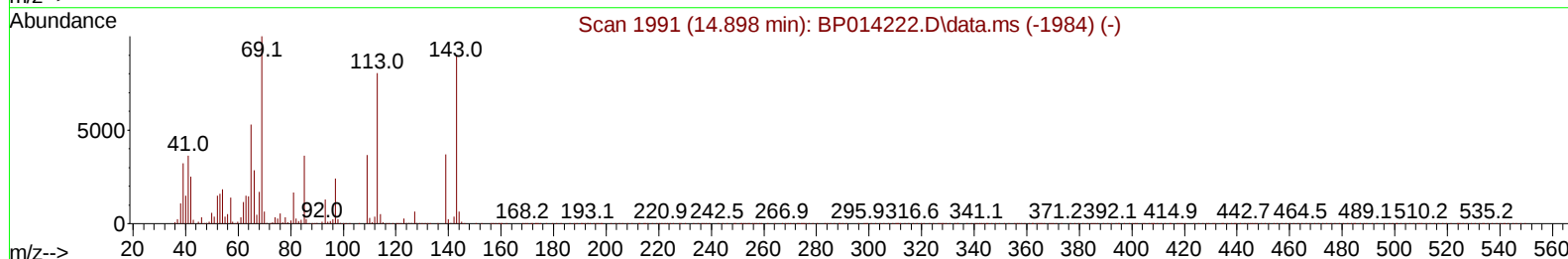
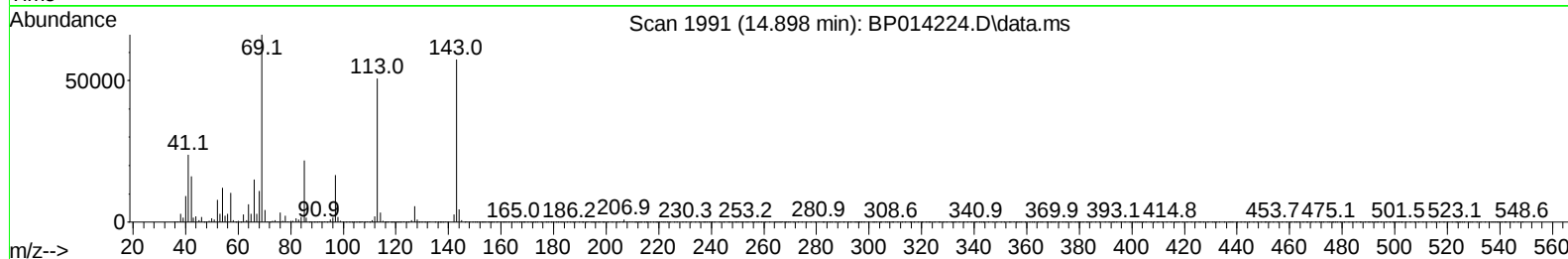
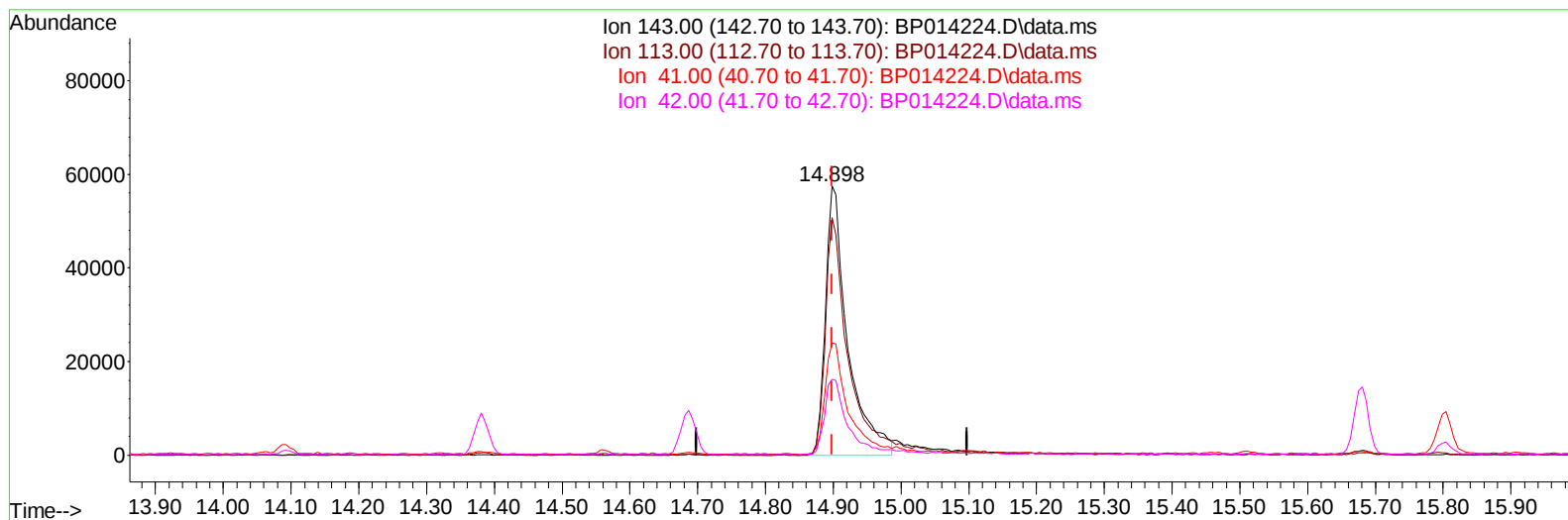
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TIC: BP014224.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.898min (-0.000) 18.63 ng/ul m

response 131687

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	88.41
41.00	40.20	41.58
42.00	28.50	28.31

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.040	152	142515	20.000	ng/ul	0.00
20) Naphthalene-d8	10.863	136	636813	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.687	164	472516	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.445	188	1145767	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.527	240	1169246	20.000	ng/ul	# 0.00
88) Perylene-d12	24.051	264	1212312	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.405	96	16551	4.388	ng/uL	0.00
4) Pyridine-d5	3.846	84	197388	18.900	ng/ul	0.00
7) Phenol-d5	7.199	99	300008	23.855	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.363	67	199257	27.166	ng/ul	0.00
11) 2-Chlorophenol-d4	7.569	132	227843	23.779	ng/ul	0.00
15) 4-Methylphenol-d8	8.757	113	248626	24.202	ng/ul	0.00
21) Nitrobenzene-d5	9.216	128	120723	23.759	ng/ul	0.00
24) 2-Nitrophenol-d4	9.946	143	144029	23.680	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.487	165	242700	21.207	ng/ul	0.00
31) 4-Chloroaniline-d4	11.010	131	325996	21.525	ng/ul	0.00
46) Dimethylphthalate-d6	14.092	166	981948	24.487	ng/ul	0.00
49) Acenaphthylene-d8	14.381	160	1029779	24.241	ng/ul	0.00
54) 4-Nitrophenol-d4	14.898	143	131687m	18.635	ng/ul	0.00
60) Fluorene-d10	15.681	176	821421	24.169	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.804	200	160175	18.496	ng/ul	0.00
73) Anthracene-d10	17.545	188	1361233	24.473	ng/ul	0.00
81) Pyrene-d10	19.769	212	1673569	25.952	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.886	264	1602356	24.848	ng/ul	0.00

Target Compounds Qvalue

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

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