

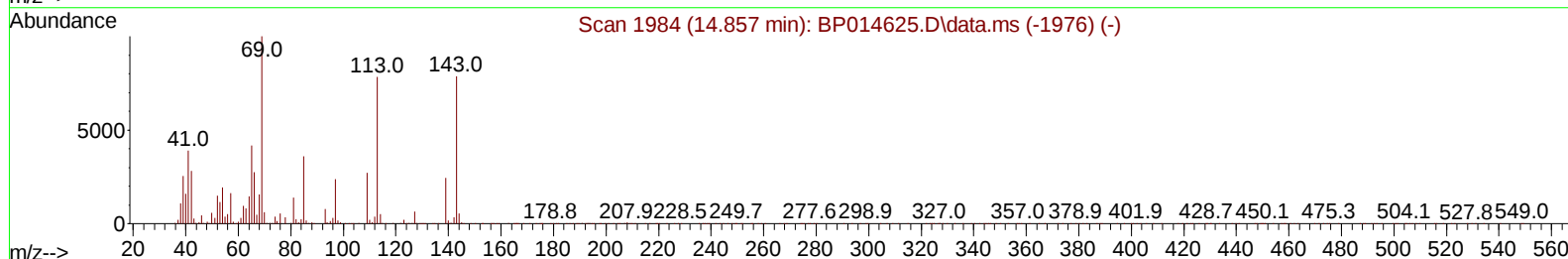
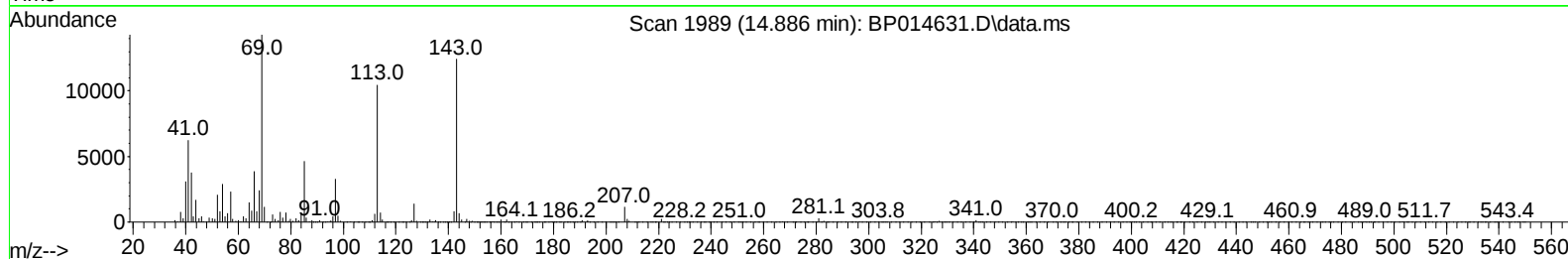
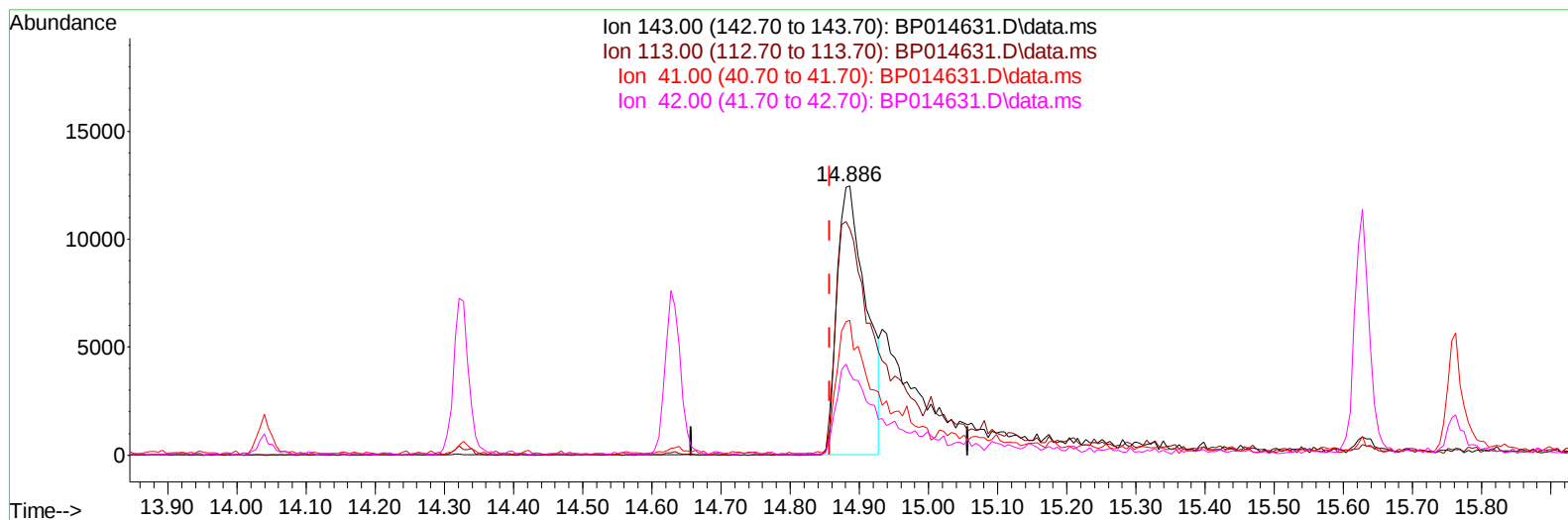
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041023\
Data File : BP014631.D
Acq On : 10 Apr 2023 17:40
Operator : CG/JU
Sample : PB152002BL
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK002

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 04/11/2023
Supervised By :Jagrut Upadhyay 04/11/2023

Quant Time: Apr 11 01:23:29 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Apr 11 01:21:43 2023
Response via : Initial Calibration



TIC: BP014631.D\data.ms

(54) 4-Nitrophenol-d4 (S)**14.886min (+ 0.029) 10.71 ng/u1****response 36292**

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	84.04
41.00	44.20	50.17
42.00	31.50	30.20

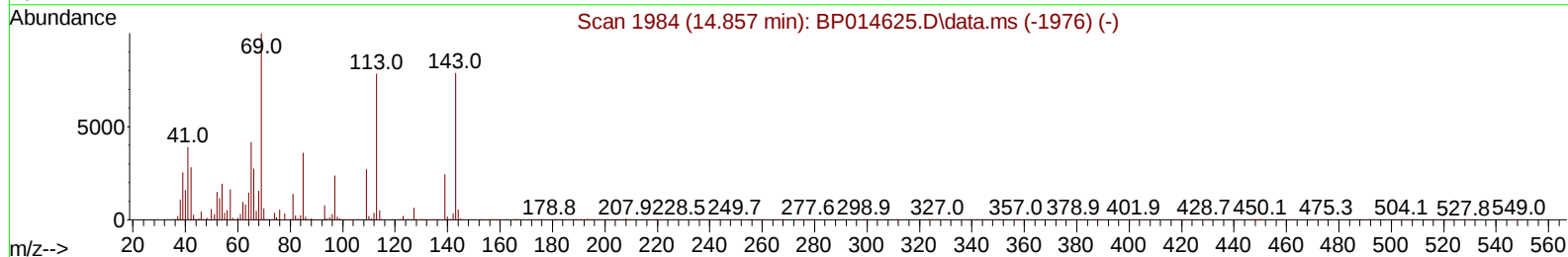
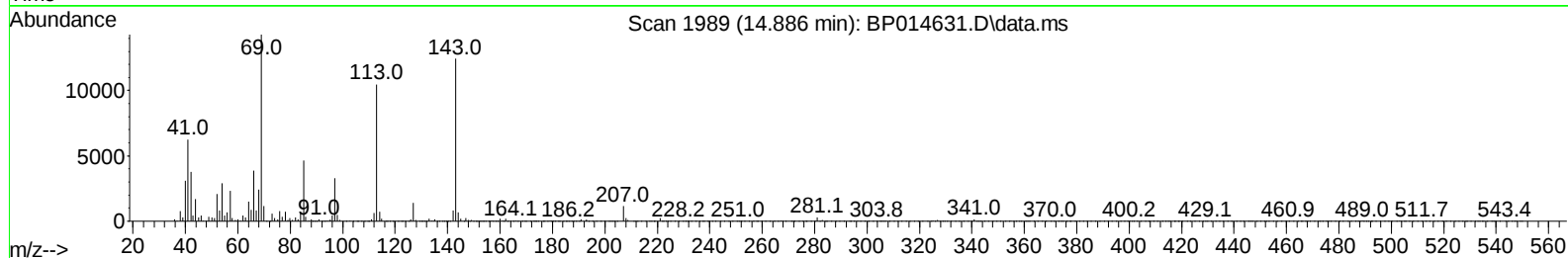
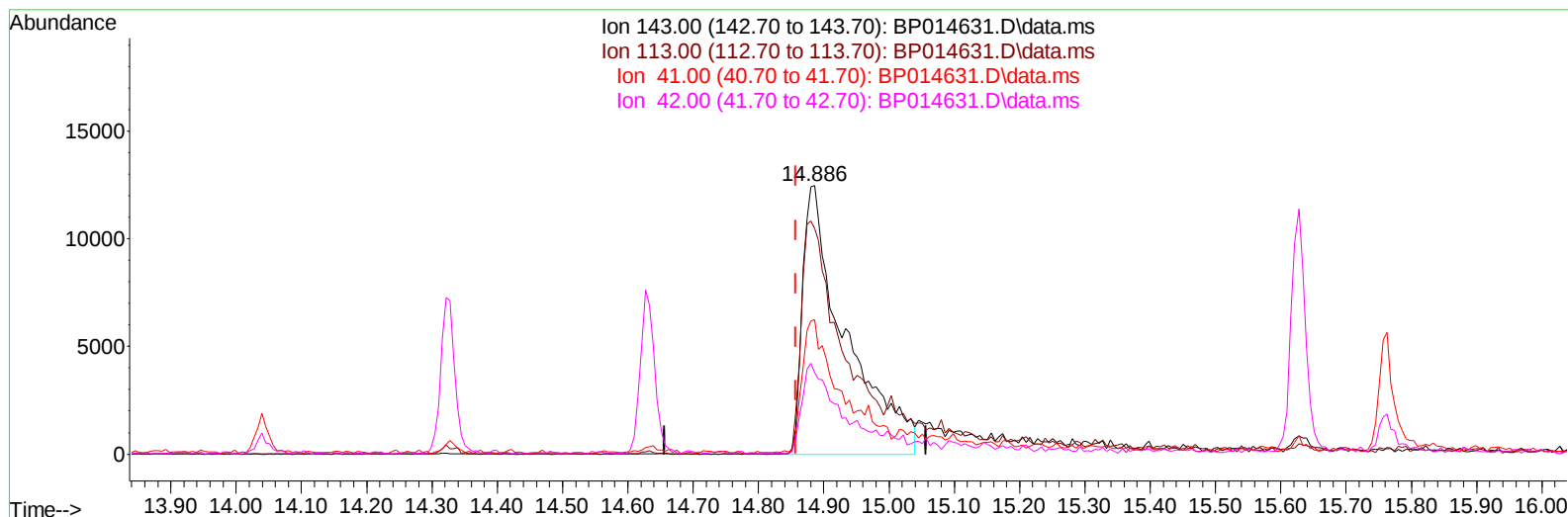
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TIC: BP014631.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.886min (+ 0.029) 16.75 ng/ul m

response 56758

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	84.04
41.00	44.20	50.17
42.00	31.50	30.20

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.975	152	107236	20.000	ng/ul	0.00
20) Naphthalene-d8	10.798	136	445456	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.633	164	276670	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.392	188	585631	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	576388	20.000	ng/ul	0.00
88) Perylene-d12	23.986	264	673679	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.363	96	16250	5.876	ng/uL	0.00
4) Pyridine-d5	3.793	84	189112	26.686	ng/ul	0.00
7) Phenol-d5	7.145	99	251239	25.404	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.304	67	184602	29.092	ng/ul	0.00
11) 2-Chlorophenol-d4	7.510	132	198763	27.807	ng/ul	0.00
15) 4-Methylphenol-d8	8.698	113	196004	24.570	ng/ul	0.00
21) Nitrobenzene-d5	9.157	128	93676	26.049	ng/ul	0.00
24) 2-Nitrophenol-d4	9.881	143	105320	25.577	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.428	165	171107	22.817	ng/ul	0.00
31) 4-Chloroaniline-d4	10.951	131	244934	24.796	ng/ul	0.00
46) Dimethylphthalate-d6	14.039	166	601875	26.830	ng/ul	0.00
49) Acenaphthylene-d8	14.327	160	675535	27.139	ng/ul	0.00
54) 4-Nitrophenol-d4	14.886	143	56758m	16.748	ng/ul	0.03
60) Fluorene-d10	15.627	176	501816	26.824	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.763	200	86016	19.961	ng/ul	0.00
73) Anthracene-d10	17.492	188	808865	28.539	ng/ul	0.00
81) Pyrene-d10	19.727	212	988699	25.537	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.821	264	1039174	29.034	ng/ul	0.00

Target Compounds Qvalue

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

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