

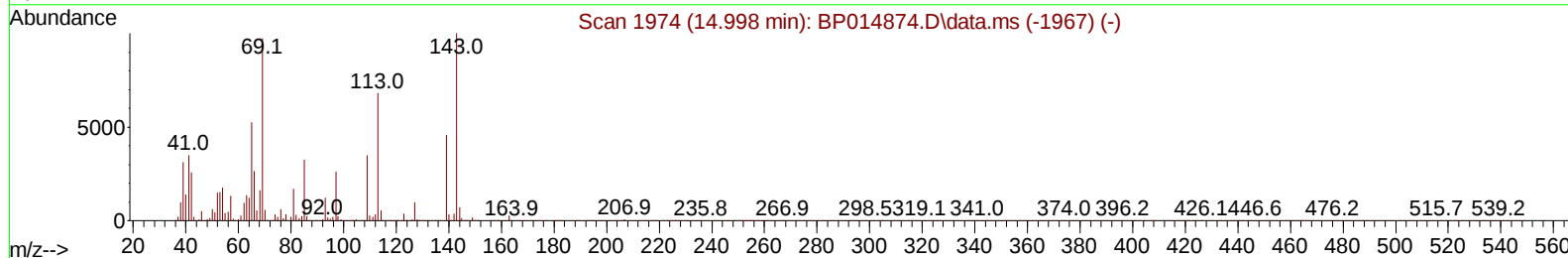
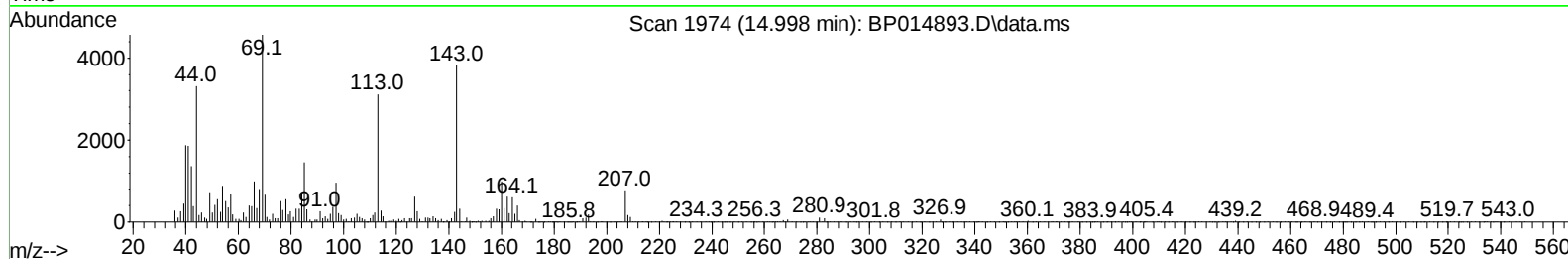
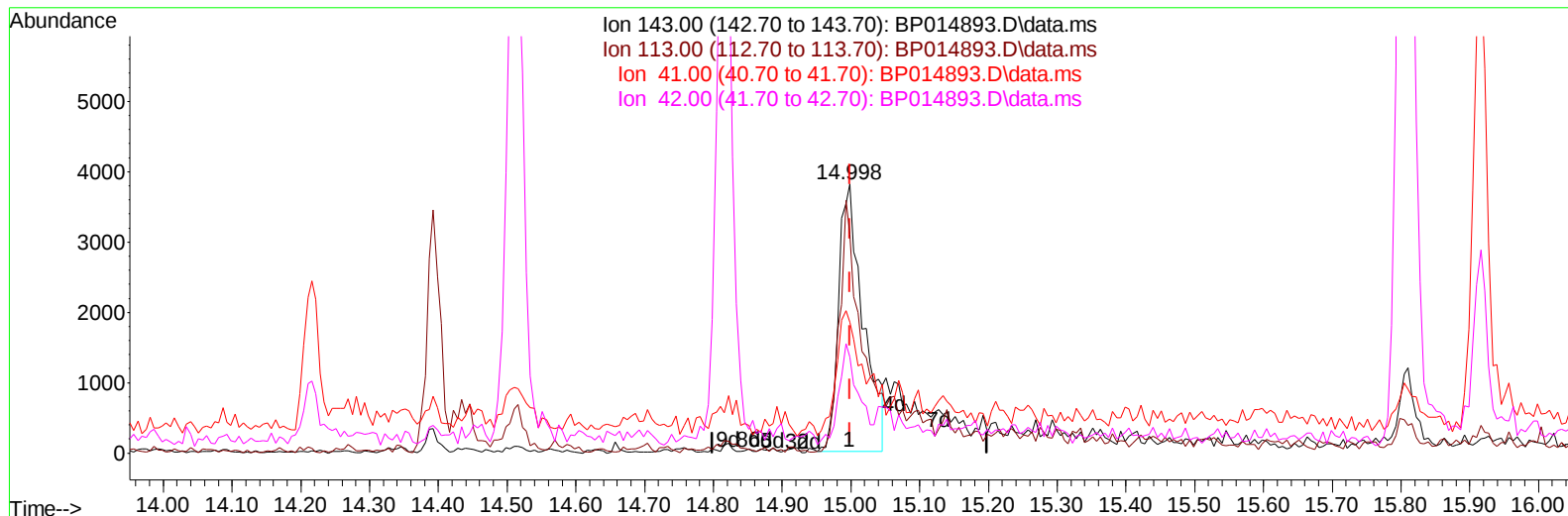
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042823\
Data File : BP014893.D
Acq On : 29 Apr 2023 02:03
Operator : CG/JU
Sample : 02528-03
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
HOA35

Manual IntegrationsAPPROVED

Quant Time: Apr 29 03:53:17 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042823.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Apr 29 00:40:33 2023
Response via : Initial Calibration

Reviewed By :Christian Giraldo 05/01/2023
Supervised By :Jagrut Upadhyay 05/01/2023



TIC: BP014893.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.998min (-0.000) 1.93 ng/u1

response 9470

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	68.20	81.56
41.00	36.10	48.73#
42.00	26.80	35.68#

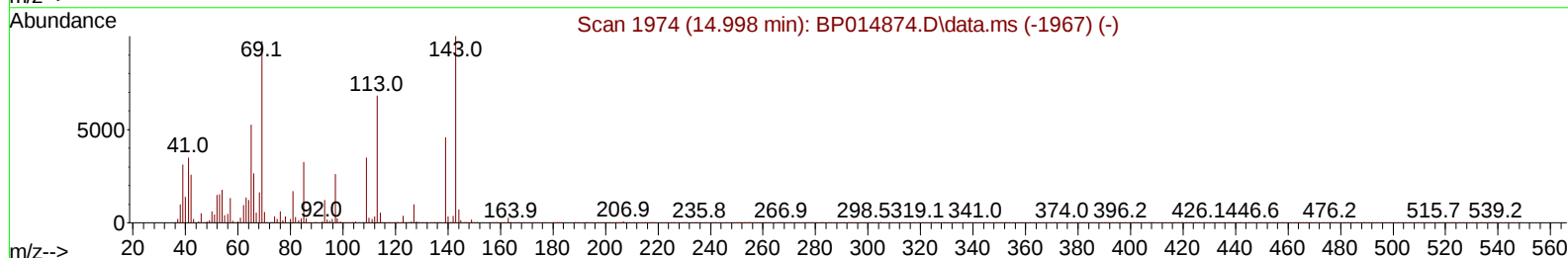
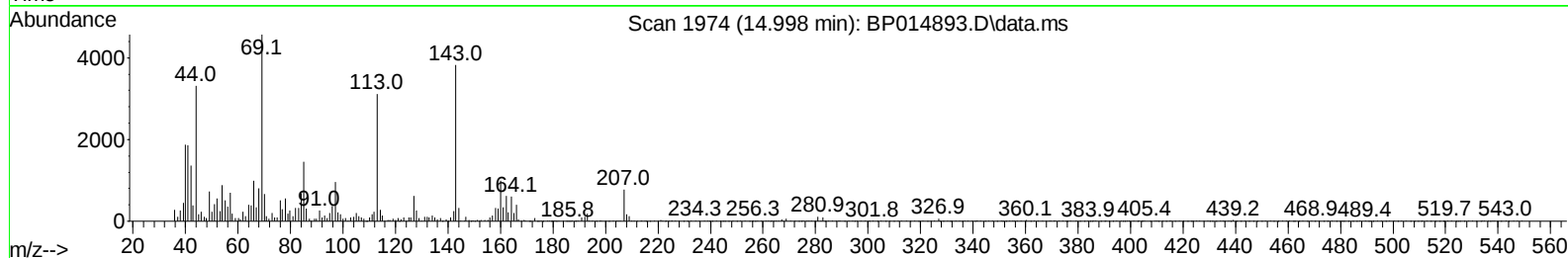
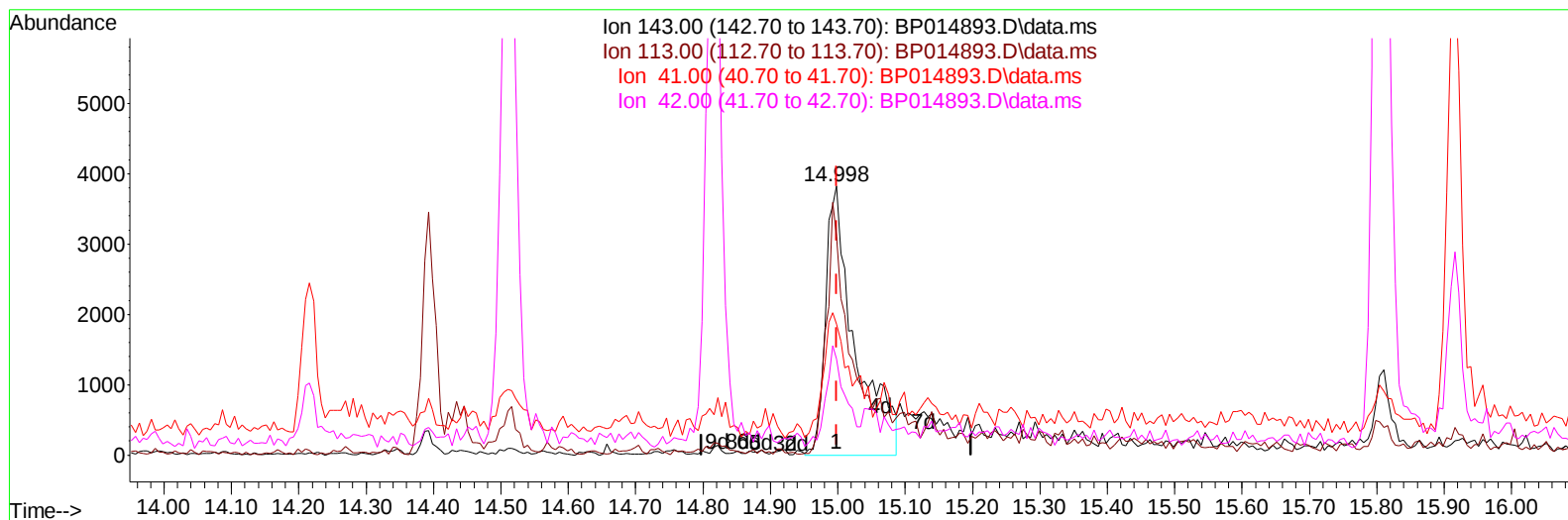
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TIC: BP014893.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.998min (-0.000) 2.38 ng/ul m

response 11656

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	68.20	81.56
41.00	36.10	48.73#
42.00	26.80	35.68#

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 Sample : 02528-03
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 ALS Vial : 24 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 05/01/2023
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Quant Time: Apr 29 04:11:40 2023
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 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Apr 29 00:40:33 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.175	152	202096	20.000	ng/ul	0.00
20) Naphthalene-d8	11.010	136	798061	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.816	164	459525	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.575	188	904340	20.000	ng/ul	0.00
79) Chrysene-d12	21.651	240	785662	20.000	ng/ul	-0.01
88) Perylene-d12	24.257	264	909678	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.464	96	24725	4.409	ng/uL	0.00
4) Pyridine-d5	3.911	84	110296	8.182	ng/ul	0.00
7) Phenol-d5	7.311	99	81934	5.180	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.487	67	295334	29.770	ng/ul	0.00
11) 2-Chlorophenol-d4	7.693	132	279107	23.305	ng/ul	0.00
15) 4-Methylphenol-d8	8.875	113	159199	13.209	ng/ul	0.00
21) Nitrobenzene-d5	9.346	128	168456	29.398	ng/ul	0.00
24) 2-Nitrophenol-d4	10.081	143	161277	30.660	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.616	165	287379	25.438	ng/ul	0.00
31) 4-Chloroaniline-d4	11.140	131	334819	20.499	ng/ul	0.00
46) Dimethylphthalate-d6	14.216	166	1148015	34.860	ng/ul	0.00
49) Acenaphthylene-d8	14.510	160	1275706	32.429	ng/ul	0.00
54) 4-Nitrophenol-d4	14.998	143	11656m	2.379	ng/ul	0.00
60) Fluorene-d10	15.810	176	991125	34.526	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.916	200	108892	28.467	ng/ul	0.00
73) Anthracene-d10	17.669	188	1593644	38.586	ng/ul	0.00
81) Pyrene-d10	19.886	212	1810643	41.494	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.080	264	1738131	38.685	ng/ul	-0.02
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.505	88	364395	61.028	ng/uL	98

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

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