

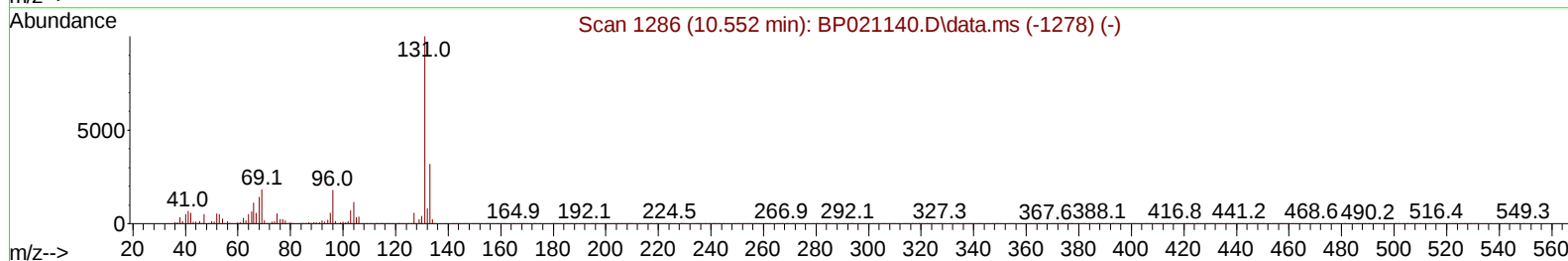
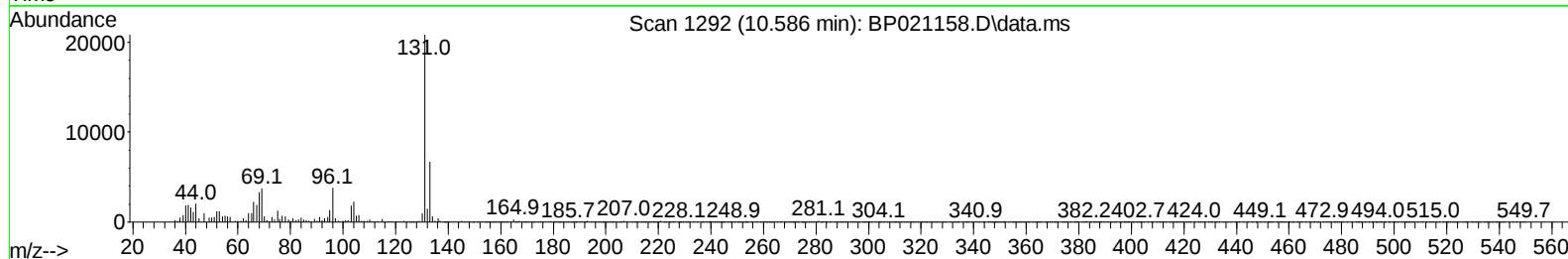
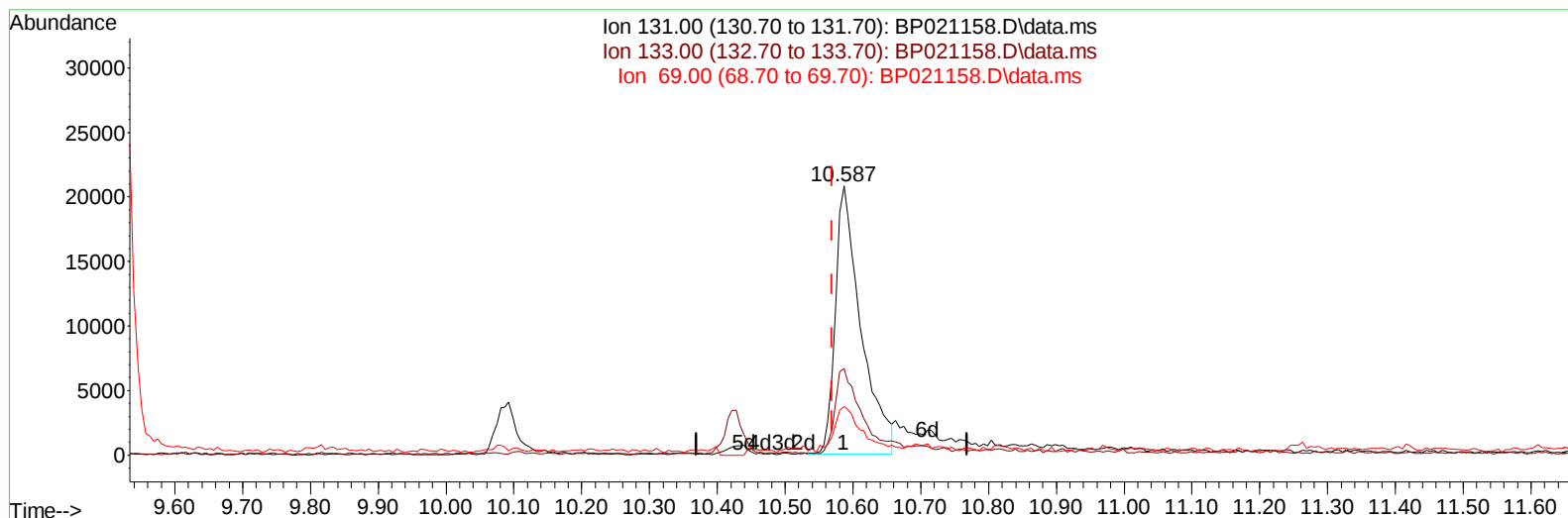
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072524\
Data File : BP021158.D
Acq On : 25 Jul 2024 22:26
Operator : MA/JU
Sample : P3263-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CG7

Manual IntegrationsAPPROVED

Quant Time: Jul 25 23:52:37 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jul 25 00:15:24 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 07/26/2024
Supervised By :mohammad ahmed 07/27/2024



TIC: BP021158.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.586min (+ 0.017) 3.23 ng/u1

response 54482

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.50	32.31
69.00	17.40	17.98
0.00	0.00	0.00

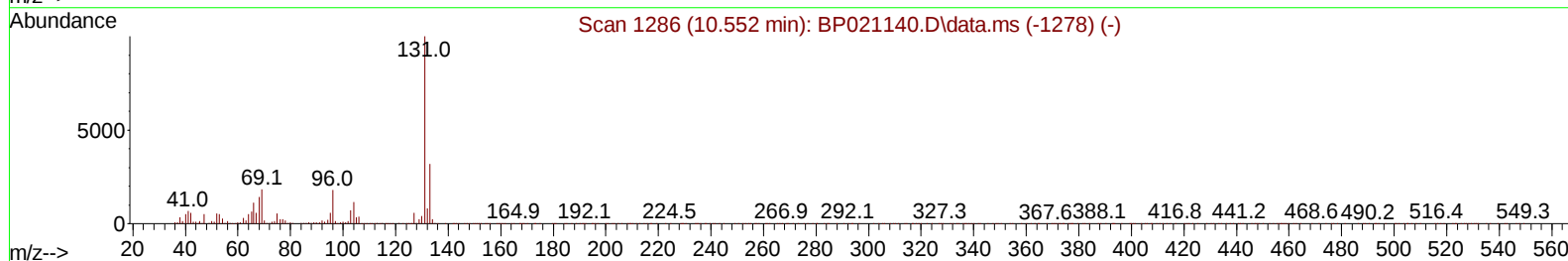
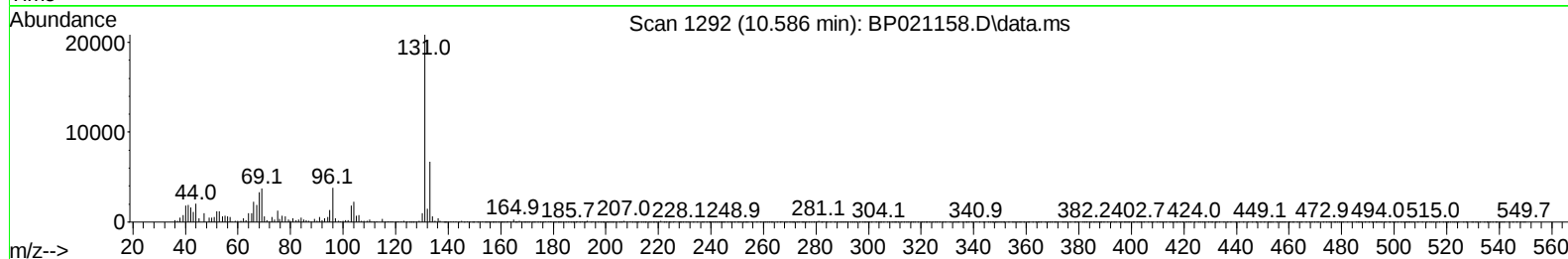
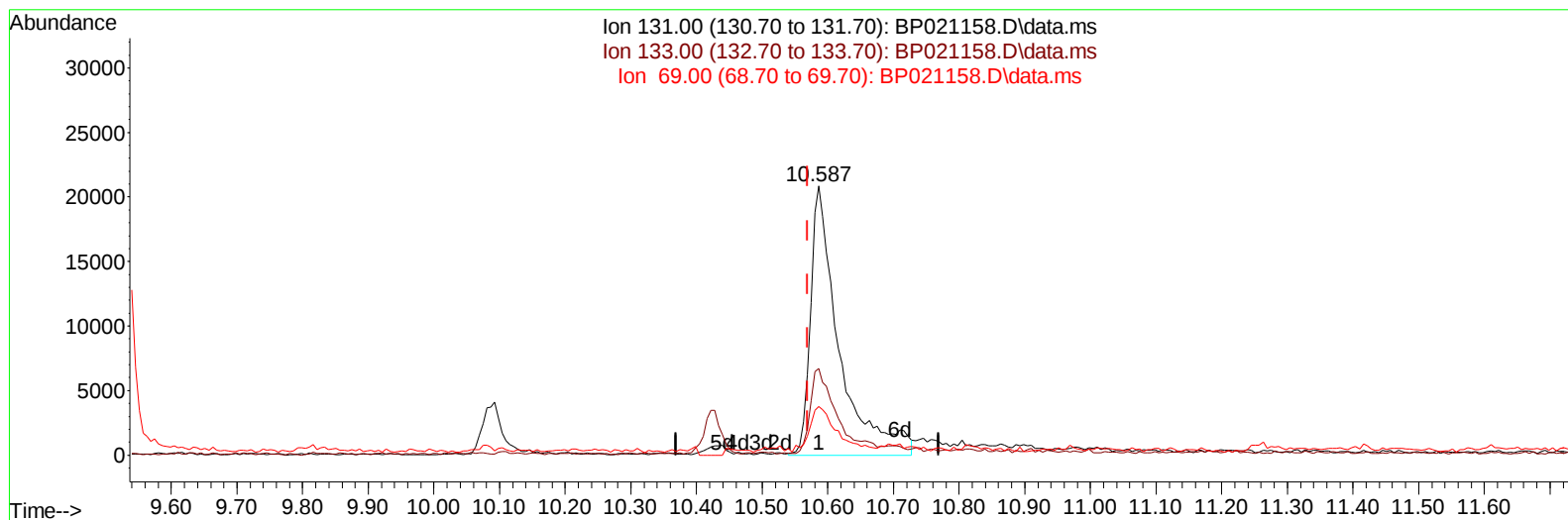
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TIC: BP021158.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.586min (+ 0.017) 3.71 ng/ul m

response 62664

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.50	32.31
69.00	17.40	17.98
0.00	0.00	0.00

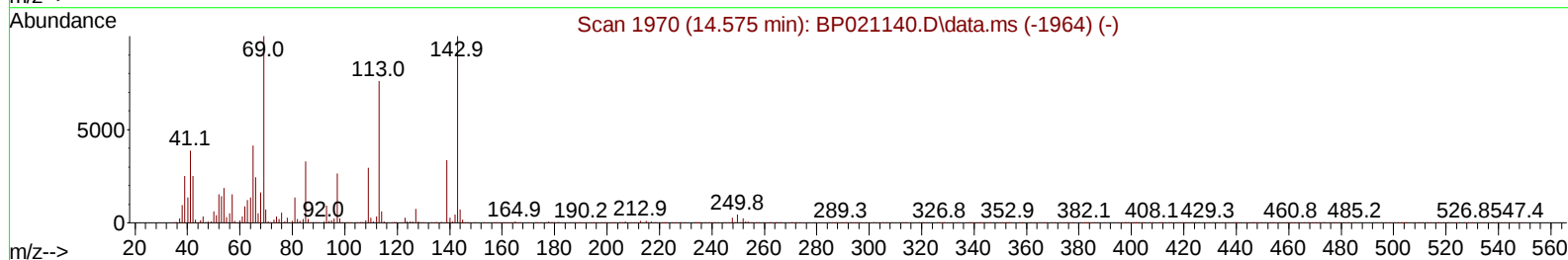
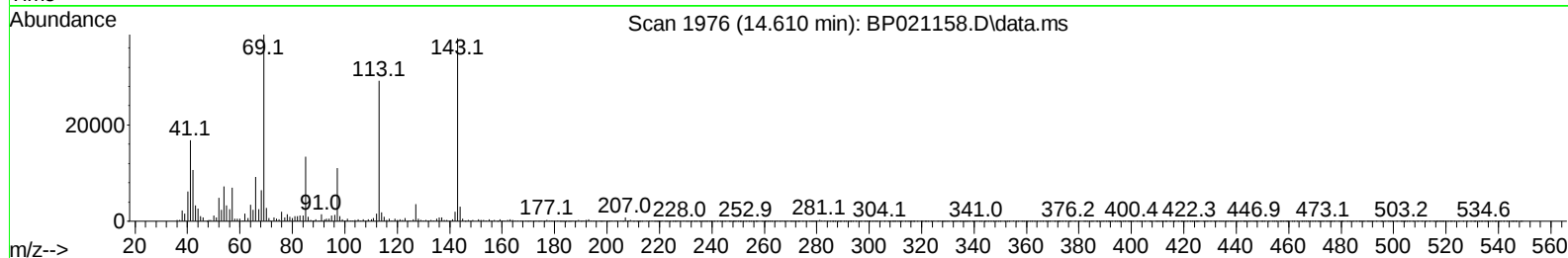
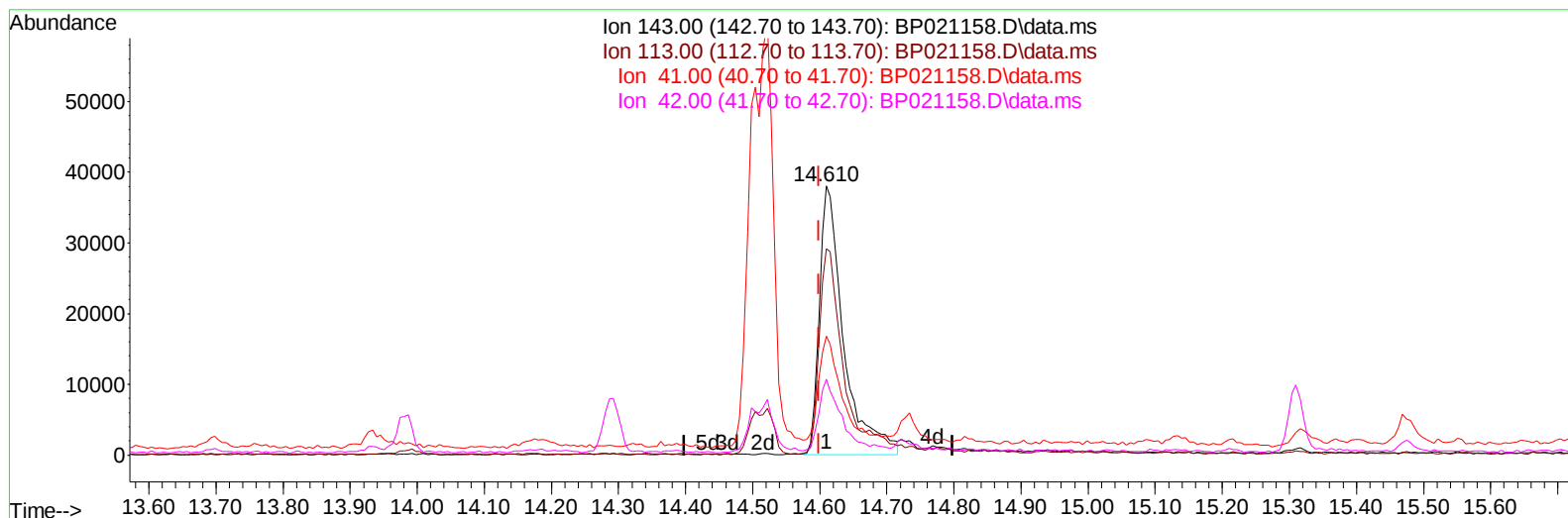
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TIC: BP021158.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.610min (+ 0.011) 19.55 ng/ul m

response 95969

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	76.76
41.00	37.90	44.26
42.00	27.10	28.08

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 Misc :
 ALS Vial : 18 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.651	152	208731	20.000	ng/ul	0.00
20) Naphthalene-d8	10.422	136	792264	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.292	164	474724	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.110	188	1011141	20.000	ng/ul	0.00
79) Chrysene-d12	21.551	240	1050521	20.000	ng/ul	0.00
88) Perylene-d12	24.880	264	1320753	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	15429	3.364	ng/uL	0.00
4) Pyridine-d5	3.622	84	16645	1.253	ng/ul	0.02
7) Phenol-d5	6.875	99	300658	17.668	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.998	67	173031	16.778	ng/ul	0.00
11) 2-Chlorophenol-d4	7.204	132	256630	18.304	ng/ul	0.00
15) 4-Methylphenol-d8	8.387	113	243177	17.206	ng/ul	0.00
21) Nitrobenzene-d5	8.816	128	118688	19.288	ng/ul	0.01
24) 2-Nitrophenol-d4	9.522	143	138678	19.690	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.087	165	237459	18.645	ng/ul	0.01
31) 4-Chloroaniline-d4	10.586	131	62664m	3.714	ng/ul	0.02
46) Dimethylphthalate-d6	13.698	166	670571	19.431	ng/ul	0.00
49) Acenaphthylene-d8	13.986	160	795454	20.528	ng/ul	0.00
54) 4-Nitrophenol-d4	14.610	143	95969m	19.545	ng/ul	0.01
60) Fluorene-d10	15.310	176	591148	20.880	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.474	200	76122	12.442	ng/ul	0.00
73) Anthracene-d10	17.210	188	896079	19.953	ng/ul	0.00
81) Pyrene-d10	19.592	212	1158405	17.836	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.645	264	1324276	20.330	ng/ul	0.00

Target Compounds Qvalue

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

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