

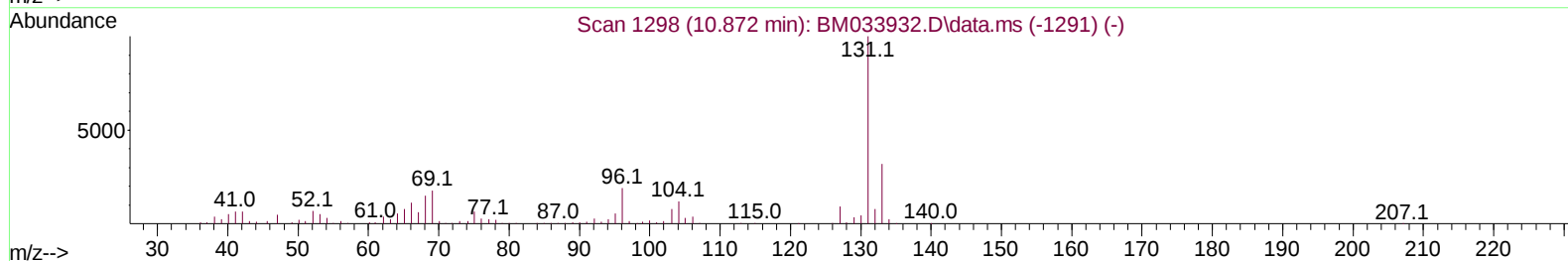
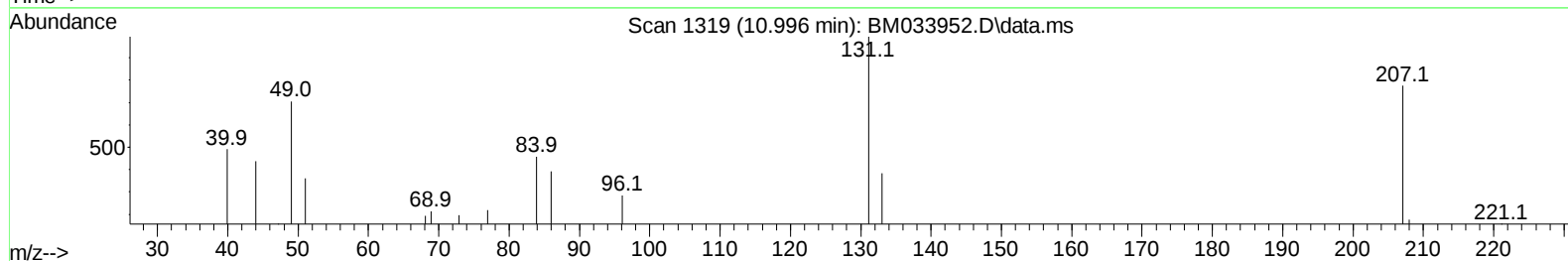
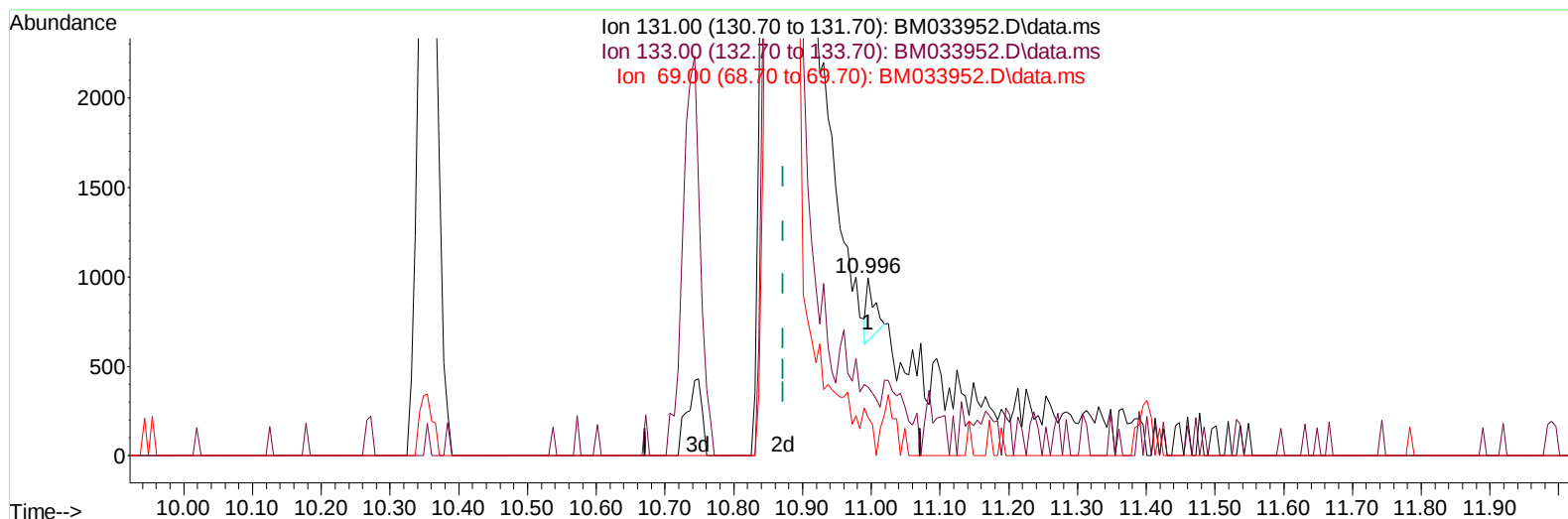
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033952.D
Acq On : 01 Feb 2022 23:16
Operator : CG/JU
Sample : N1307-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0Q33

Manual IntegrationsAPPROVED

Quant Time: Feb 02 01:52:05 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Feb 01 02:16:28 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 02/02/2022
Supervised By :Yogesh Patel 02/04/2022



TIC: BM033952.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.996min (+ 0.124) 0.03 ng/u1

response 277

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	33.10	38.63
69.00	22.50	21.53
0.00	0.00	0.00

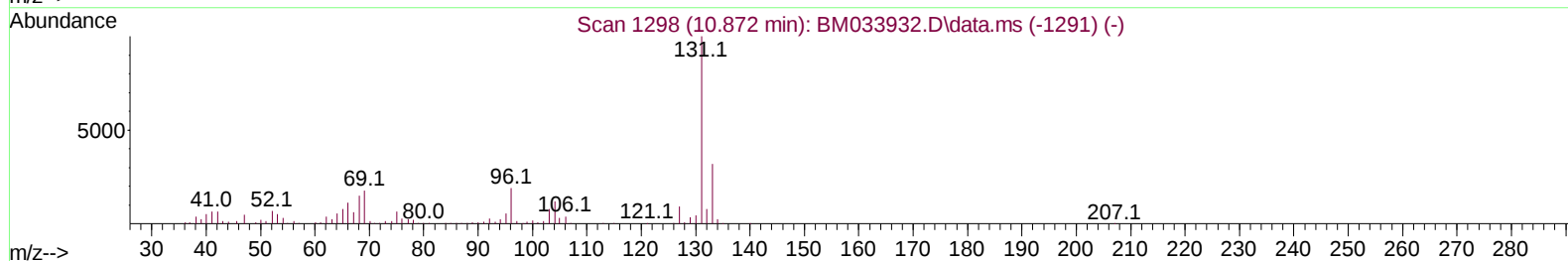
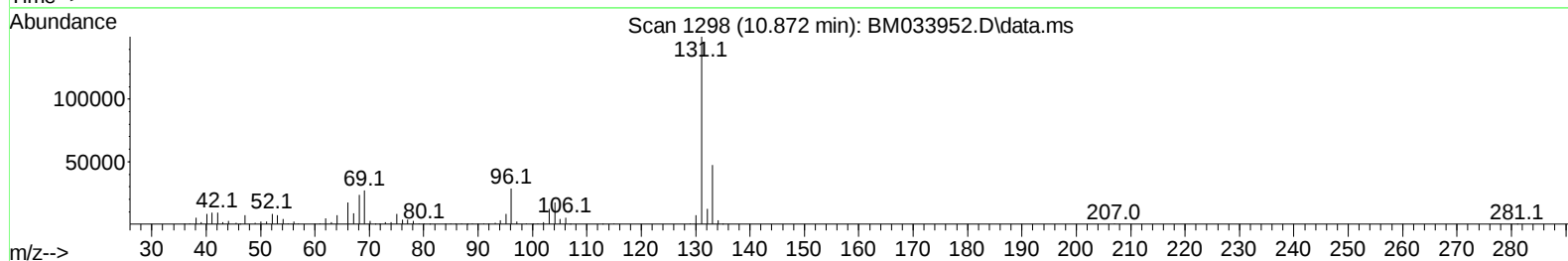
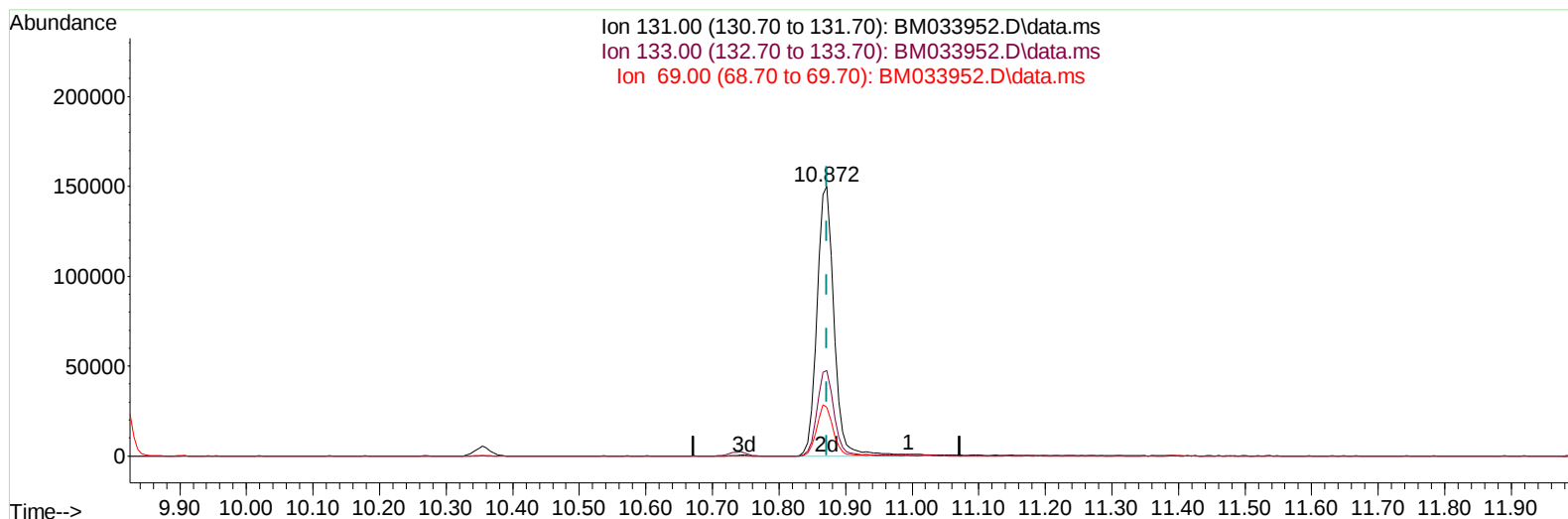
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(31) 4-Chloroaniline-d4 (S)

10.872min (+ 0.000) 25.88 ng/ul m

response 265678

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	33.10	31.69
69.00	22.50	17.95#
0.00	0.00	0.00

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.925	152		112528	20.000	ng/ul	0.00
20) Naphthalene-d8	10.737	136		474106	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.566	164		288829	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.301	188		538608	20.000	ng/ul	0.00
79) Chrysene-d12	21.460	240		458068	20.000	ng/ul	0.00
88) Perylene-d12	23.853	264		464652	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.290	96		13462	4.718	ng/uL	0.00
4) Pyridine-d5	3.720	84		70982	8.913	ng/ul	0.00
7) Phenol-d5	7.084	99		62638	6.644	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.249	67		184329	32.072	ng/ul	0.00
11) 2-Chlorophenol-d4	7.455	132		188031	24.895	ng/ul	0.00
15) 4-Methylphenol-d8	8.631	113		118972	15.926	ng/ul	-0.01
21) Nitrobenzene-d5	9.090	128		119726	31.472	ng/ul	0.00
24) 2-Nitrophenol-d4	9.813	143		122020	29.702	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.354	165		196719	26.461	ng/ul	0.00
31) 4-Chloroaniline-d4	10.872	131		265678m	25.879	ng/ul	0.00
46) Dimethylphthalate-d6	13.972	166		732549	33.941	ng/ul	0.00
49) Acenaphthylene-d8	14.260	160		850305	32.160	ng/ul	0.00
54) 4-Nitrophenol-d4	14.760	143		14483	3.928	ng/ul	0.00
60) Fluorene-d10	15.554	176		597752	33.168	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.666	200		86685	25.636	ng/ul	0.00
73) Anthracene-d10	17.395	188		897467	34.794	ng/ul	0.00
81) Pyrene-d10	19.683	212		957901	35.421	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.695	264		858561	35.179	ng/ul	-0.01
Target Compounds							
59) Diethylphthalate	15.378	149		50738	2.298	ng/ul	Qvalue 99

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

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