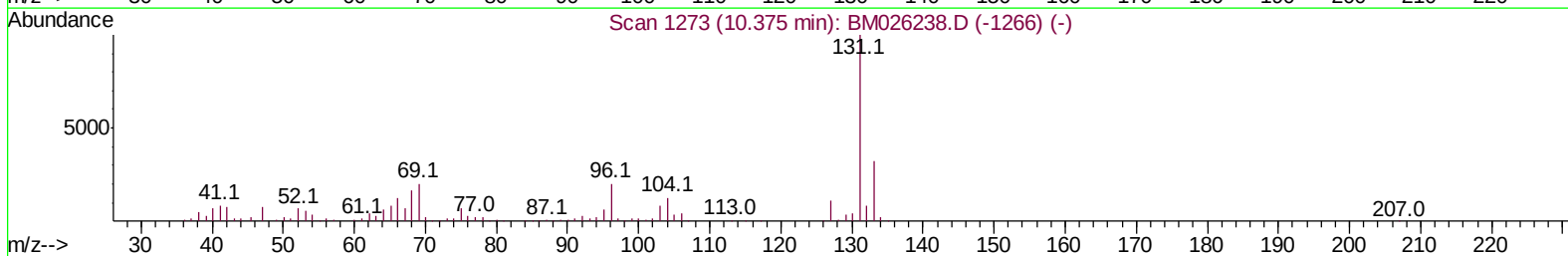
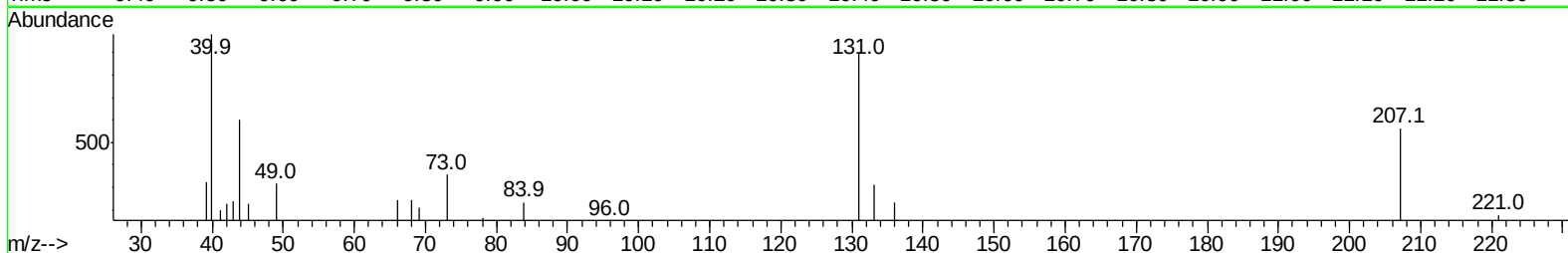
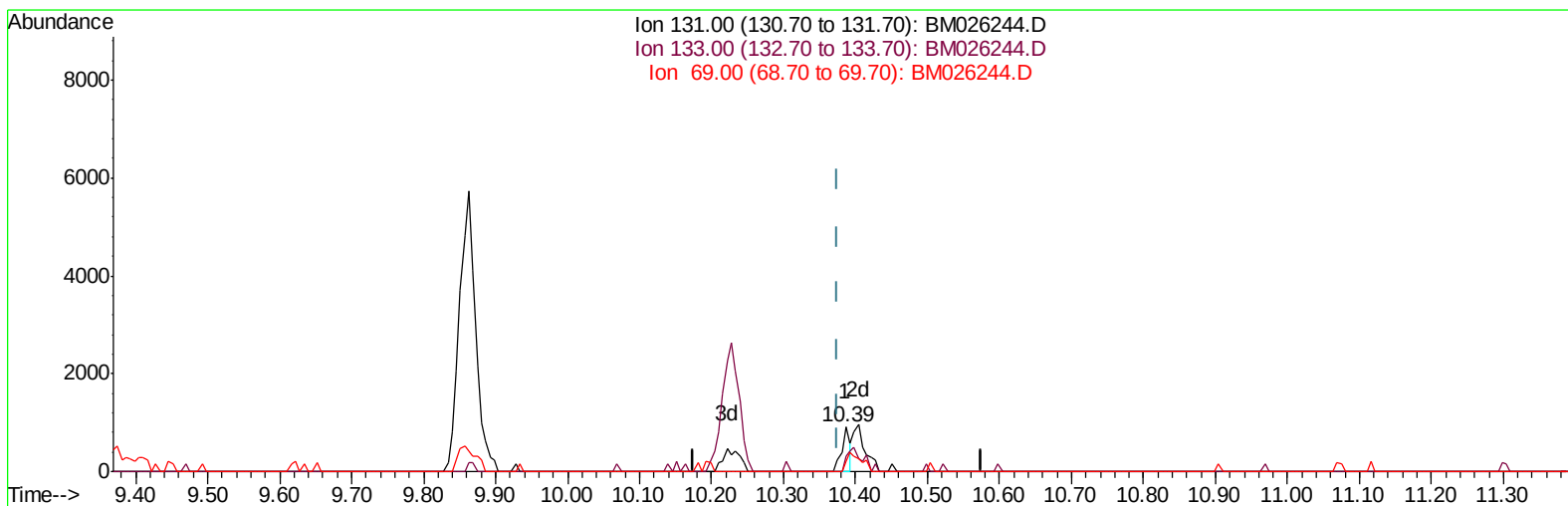


Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM060320\
Data File : BM026244.D
Acq On : 03 Jun 2020 13:35
Operator : JU/CG
Sample : L2836-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Quant Time: Jun 03 14:09:16 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jun 03 12:10:52 2020
Response via : Initial Calibration



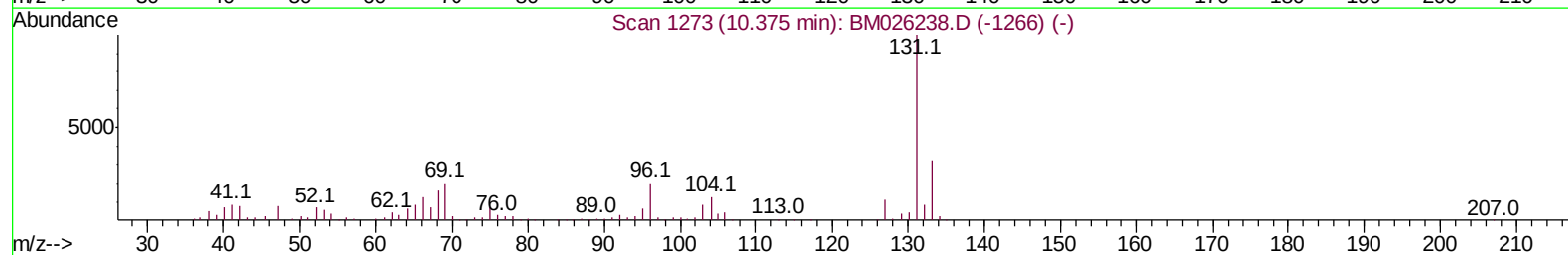
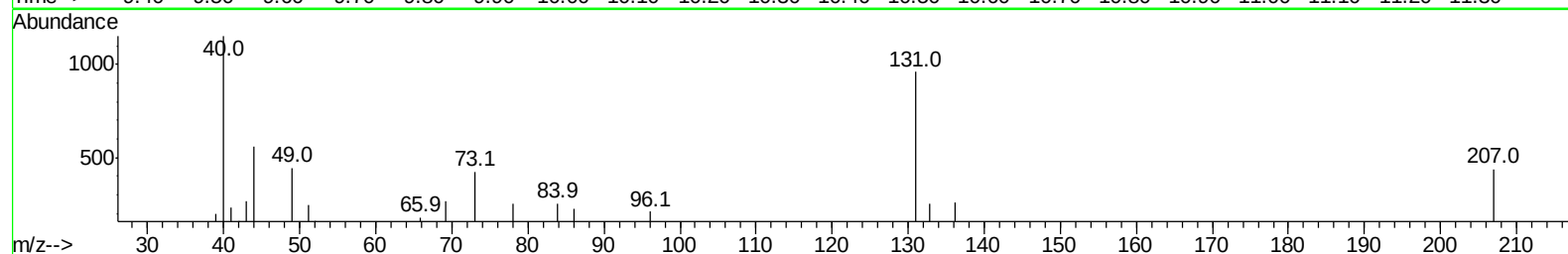
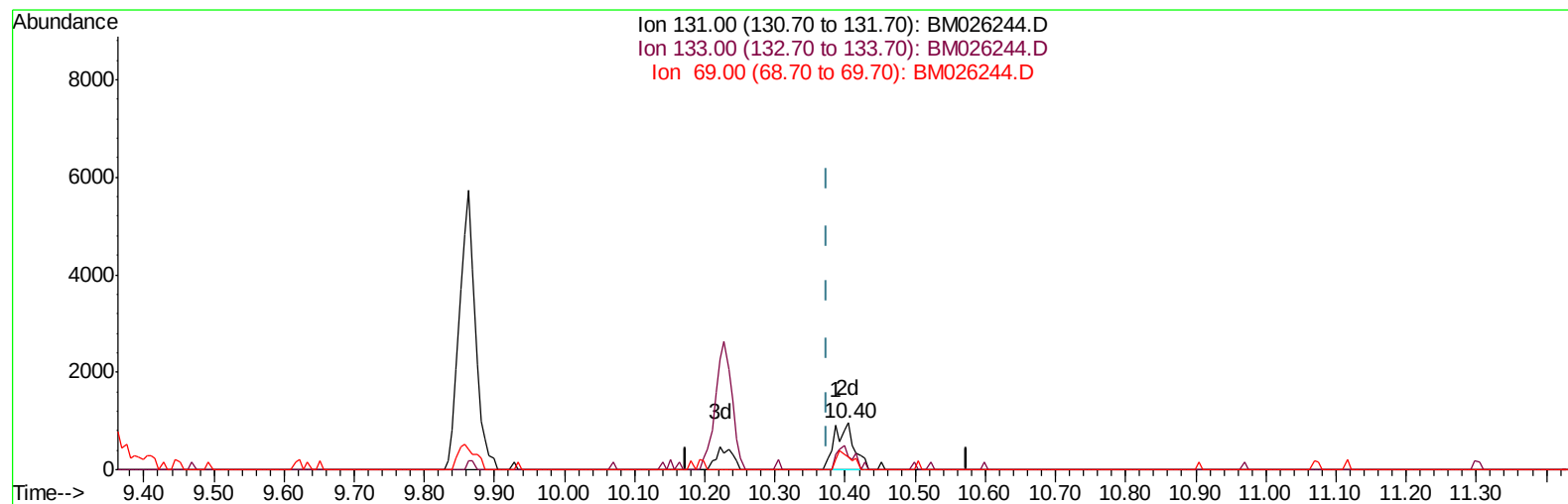
TIC: BM026244.D

(29) 4-Chloroaniline-d4 (S)
10.386min (+0.012) 0.08ng/ul
response 736

Ion	Exp%	Act%
131.00	100	100
133.00	34.00	34.30
69.00	21.90	23.20
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM060320\
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Operator : JU/CG
Sample : L2836-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Quant Time: Jun 03 14:09:16 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jun 03 12:10:52 2020
Response via : Initial Calibration



TIC: BM026244.D

(29) 4-Chloroaniline-d4 (S)

10.404min (+0.030) 0.20ng/ul m

response 1843

Ion	Exp%	Act%
131.00	100	100
133.00	34.00	26.38#
69.00	21.90	27.63#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM060320\
 Data File : BM026244.D
 Acq On : 03 Jun 2020 13:35
 Operator : JU/CG
 Sample : L2836-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Quant Time: Jun 03 14:12:03 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jun 03 12:10:52 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.46	152	139452	20.00	ng/ul	0.00
18) Naphthalene-d8	10.22	136	565972	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.11	164	330312	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.86	188	704204	20.00	ng/ul	0.00
78) Chrysene-d12	21.07	240	749828	20.00	ng/ul	0.00
86) Perylene-d12	23.18	264	798943	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.06	96	23157	5.01	ng/uL	0.00
5) Phenol-d5	6.66	99	87864	6.95	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.82	67	244251	27.91	ng/ul	0.00
9) 2-Chlorophenol-d4	7.00	132	227670	22.99	ng/ul	0.00
13) 4-Methylphenol-d8	8.18	113	148976	15.63	ng/ul	0.00
19) Nitrobenzene-d5	8.61	128	133123	28.85	ng/ul	0.00
22) 2-Nitrophenol-d4	9.33	143	140770	30.84	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.86	165	231718	25.29	ng/ul	0.00
29) 4-Chloroaniline-d4	10.40	131	1843m	0.20	ng/ul	0.03
44) Dimethylphthalate-d6	13.54	166	778422	29.39	ng/ul	0.00
47) Acenaphthylene-d8	13.80	160	890716	28.38	ng/ul	0.00
52) 4-Nitrophenol-d4	14.35	143	25718	5.48	ng/ul	0.00
58) Fluorene-d10	15.12	176	680245	29.53	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.24	200	108033	26.31	ng/ul	0.00
71) Anthracene-d10	16.96	188	1070366	30.84	ng/ul	0.00
79) Pyrene-d10	19.26	212	1224118	31.78	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.05	264	1208970	28.43	ng/ul	0.00

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

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