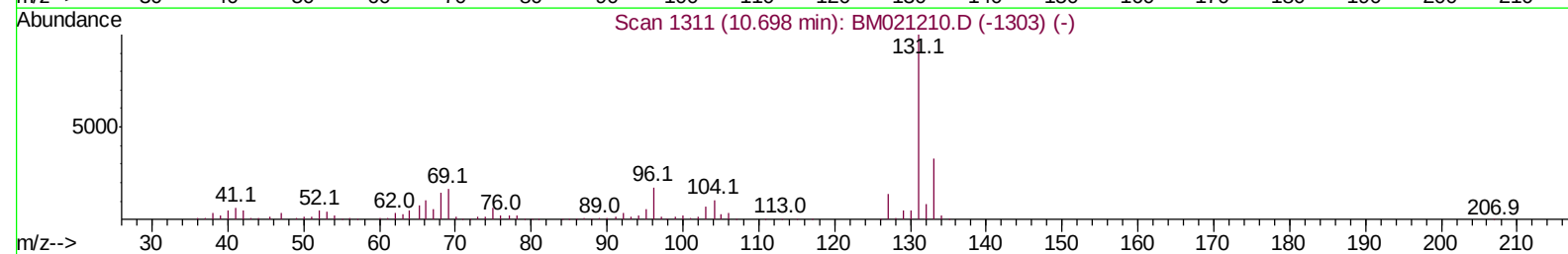
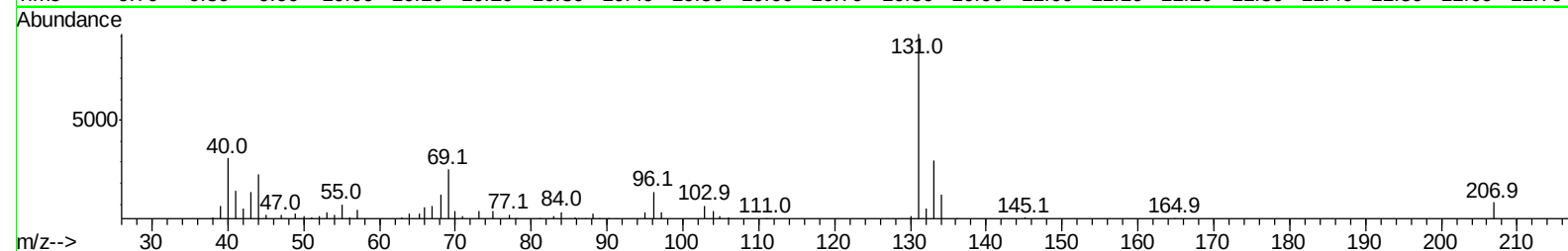
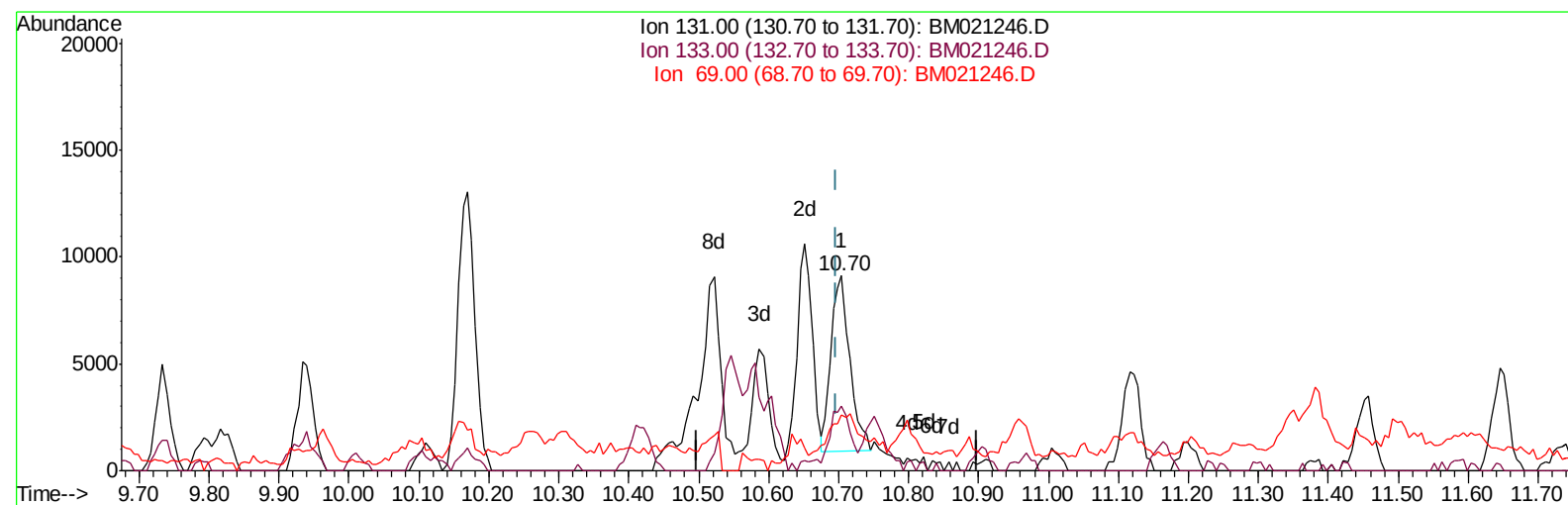


Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062819\
Data File : BM021246.D
Acq On : 29 Jun 2019 00:08
Operator : HP/JU
Sample : K3462-10
Misc :
ALS Vial : 24 Sample Multiplier: 1

Quant Time: Jun 29 00:57:33 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 27 05:07:12 2019
Response via : Initial Calibration



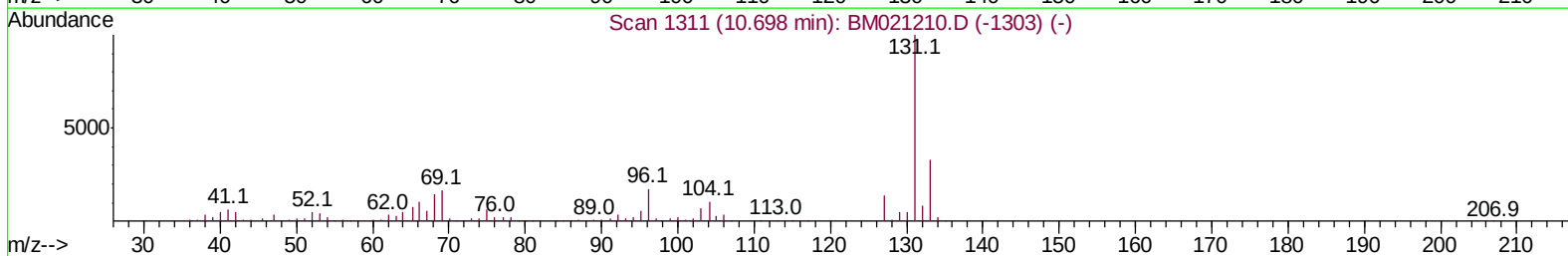
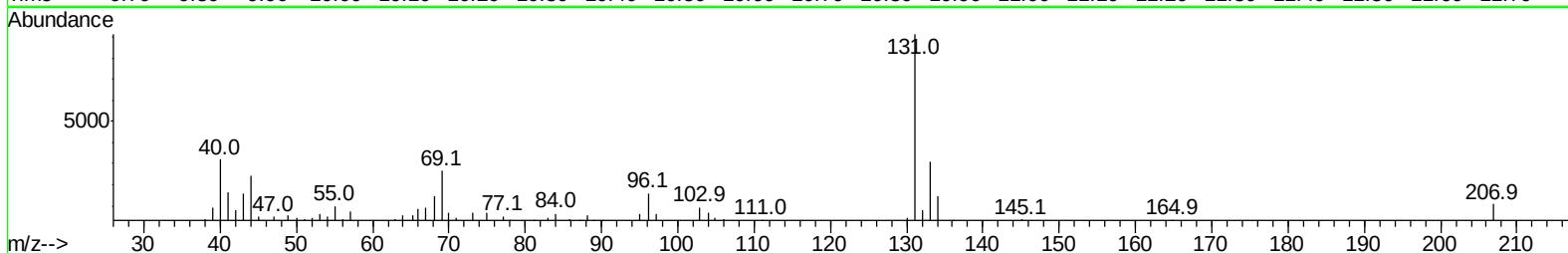
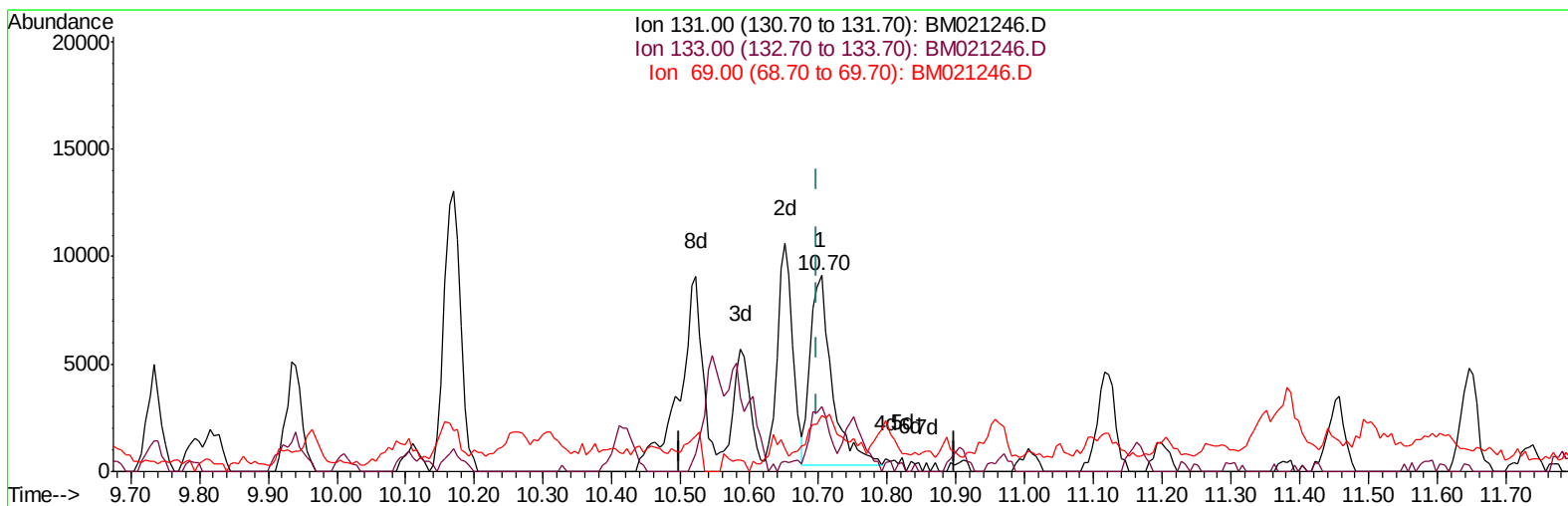
TIC: BM021246.D

(29) 4-Chloroaniline-d4 (S)
10.704min (+0.006) 0.69ng/ul
response 15435

Ion	Exp%	Act%
131.00	100	100
133.00	33.20	33.51
69.00	18.00	28.92#
0.00	0.00	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062819\
Data File : BM021246.D
Acq On : 29 Jun 2019 00:08
Operator : HP/JU
Sample : K3462-10
Misc :
ALS Vial : 24 Sample Multiplier: 1

Quant Time: Jun 29 00:57:33 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 27 05:07:12 2019
Response via : Initial Calibration



TIC: BM021246.D

(29) 4-Chloroaniline-d4 (S)

10.704min (+0.006) 0.86ng/ul m

response 19241

Ion	Exp%	Act%
131.00	100	100
133.00	33.20	33.51
69.00	18.00	28.92#
0.00	0.00	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062819\
 Data File : BM021246.D
 Acq On : 29 Jun 2019 00:08
 Operator : HP/JU
 Sample : K3462-10
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Quant Time: Jun 29 00:59:27 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jun 27 05:07:12 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.76	152	265487	20.00	ng/ul	-0.01
18) Naphthalene-d8	10.55	136	1165926	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.40	164	746270	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.15	188	1671258	20.00	ng/ul	0.00
77) Chrysene-d12	21.33	240	1383603	20.00	ng/ul	0.00
85) Perylene-d12	23.60	264	1498494	20.00	ng/ul	-0.02

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.28	96	38940	6.96	ng/uL	0.00
5) Phenol-d5	6.93	99	162541	7.03	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.10	67	374628	27.21	ng/ul	0.00
9) 2-Chlorophenol-d4	7.29	132	445299	23.71	ng/ul	0.00
13) 4-Methylphenol-d8	8.46	113	315091	16.33	ng/ul	0.00
19) Nitrobenzene-d5	8.92	128	282547	29.84	ng/ul	0.00
22) 2-Nitrophenol-d4	9.64	143	320481	30.49	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.17	165	629798	29.47	ng/ul	0.00
29) 4-Chloroaniline-d4	10.70	131	19241m	0.86	ng/ul	0.00
43) Dimethylphthalate-d6	13.82	166	2216638	32.87	ng/ul	0.00
46) Acenaphthylene-d8	14.09	160	2549602	30.89	ng/ul	0.00
51) 4-Nitrophenol-d4	14.62	143	57691	5.26	ng/ul	0.00
57) Fluorene-d10	15.40	176	1923402	32.79	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.52	200	273210	26.59	ng/ul	-0.01
70) Anthracene-d10	17.25	188	2894955	33.33	ng/ul	-0.01
78) Pyrene-d10	19.54	212	2998390	36.11	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.46	264	3001433	33.95	ng/ul	-0.01

Target Compounds

					Ovalue
6) Phenol	6.96	94	22699	1.039	ng/ul 95
44) Dimethylphthalate	13.86	163	492572	8.025	ng/ul 99

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

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