

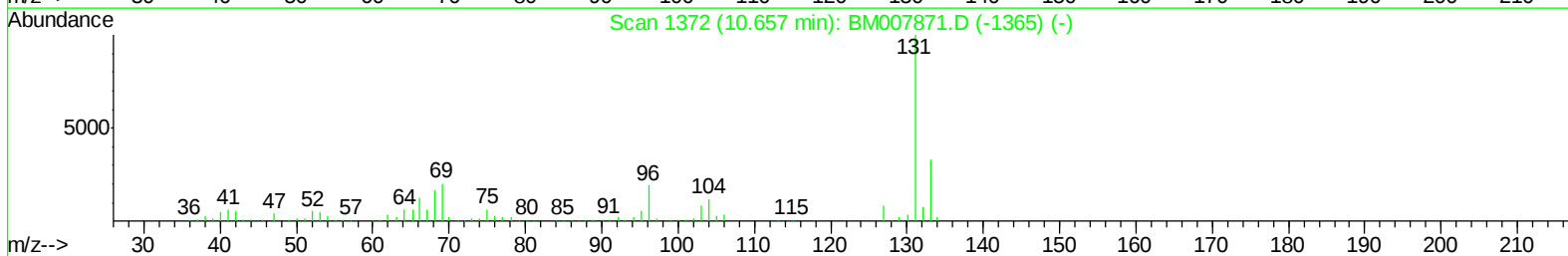
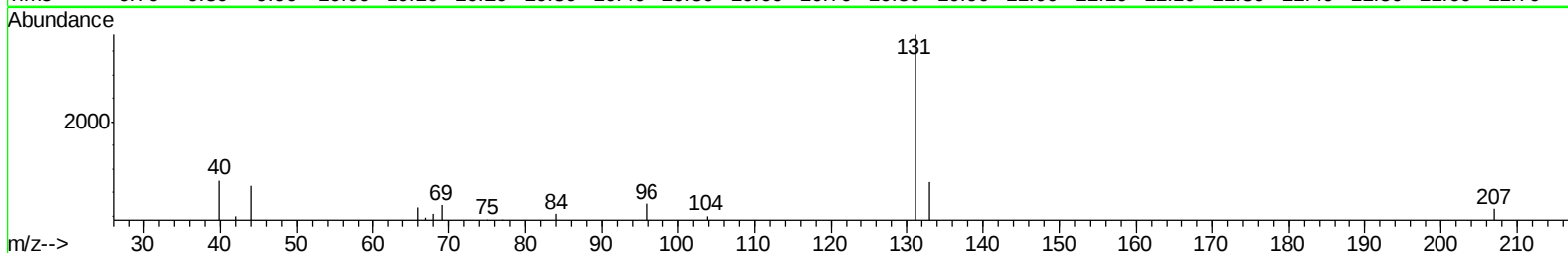
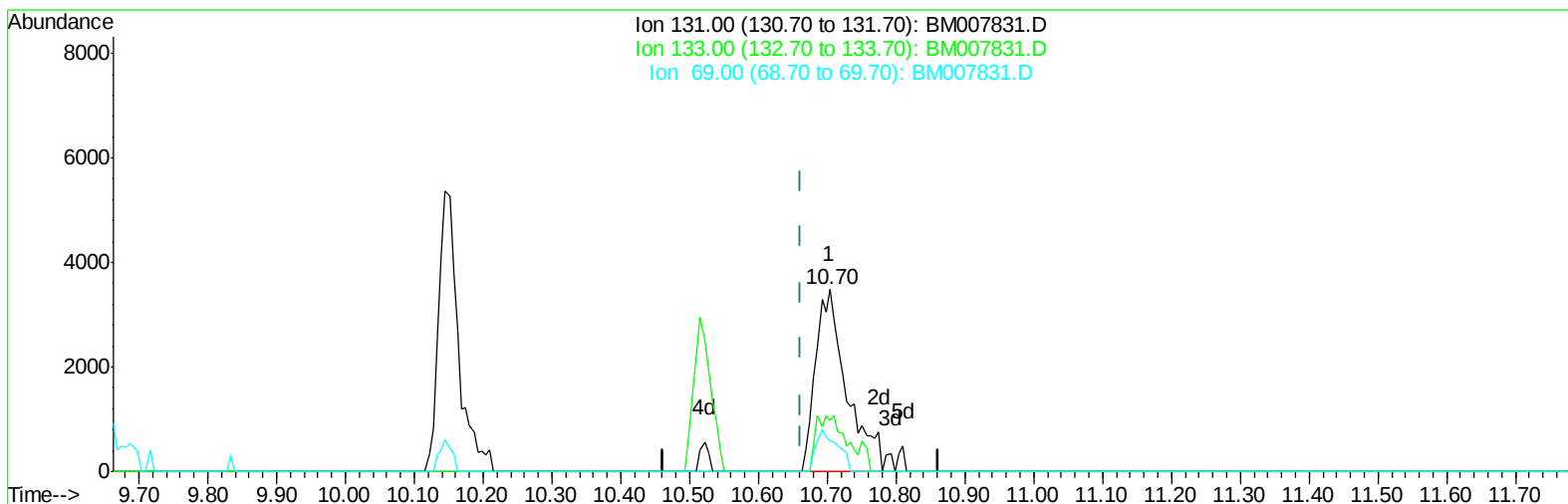
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM111116\
Data File : BM007831.D
Acq On : 11 Nov 2016 18:24
Operator : UM/SJ
Sample : H5612-07
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
F2H10

Quant Time: Nov 12 03:37:39 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM102016.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Nov 12 03:12:55 2016
Response via : Initial Calibration

Manual Integrations
APPROVED

sohil
11/14/2016 6:38:17 PM



TIC: BM007831.D

(29) 4-Chloroaniline-d4 (S)

10.704min (+0.041) 0.96ng/ul m

response 10893

Ion	Exp%	Act%
131.00	100	100
133.00	33.60	28.29
69.00	19.50	16.53
0.00	0.00	0.00

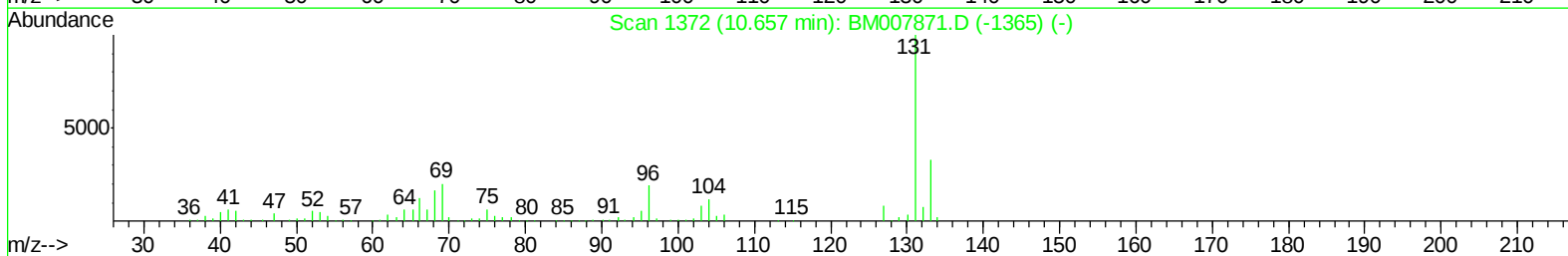
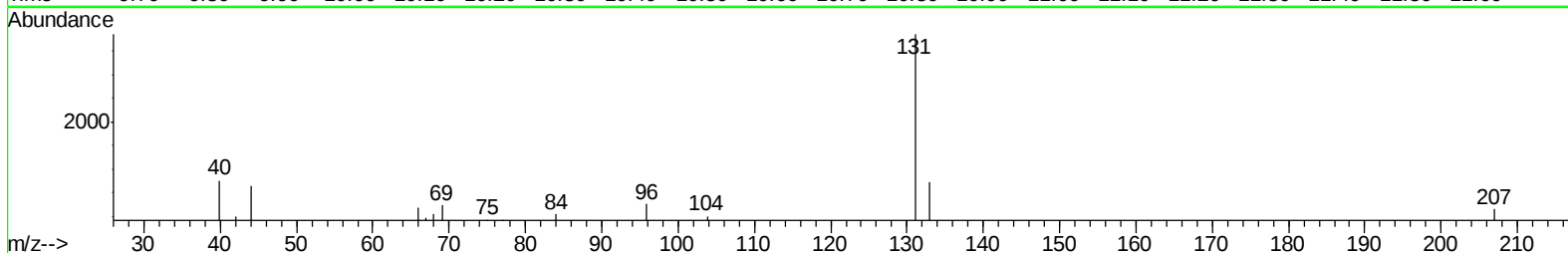
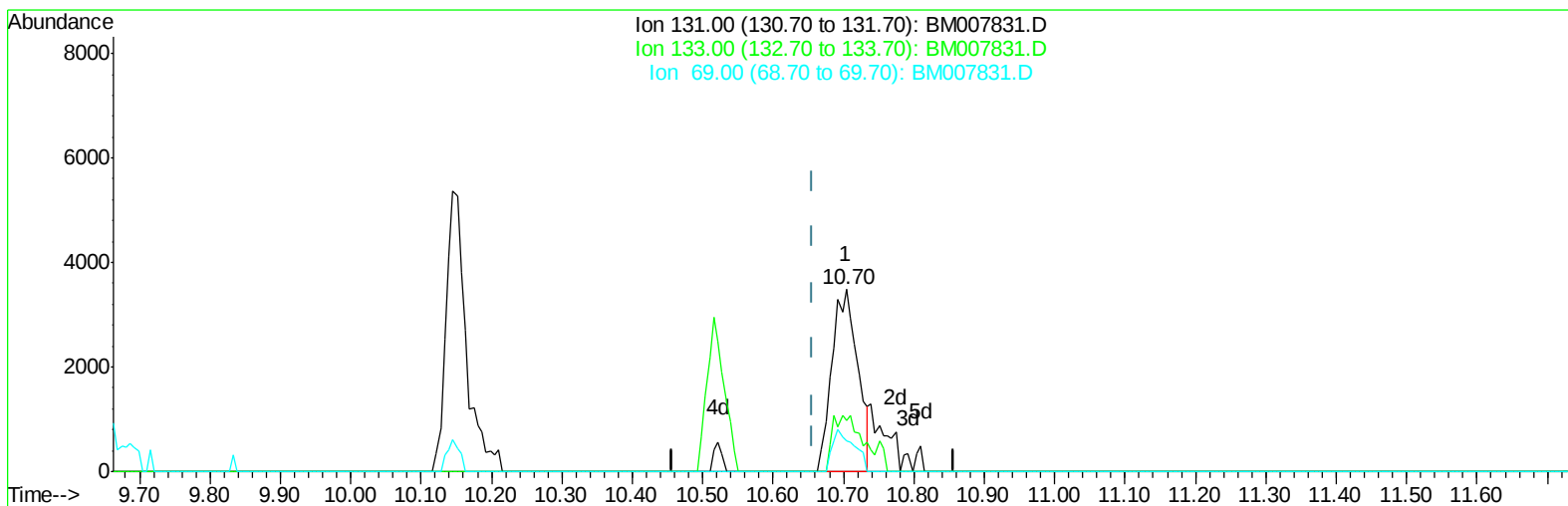
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ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Manual Integrations
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Quant Time: Nov 16 03:45:34 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM102016.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 15 03:45:58 2016
Response via : Initial Calibration



TIC: BM007831.D

(29) 4-Chloroaniline-d4 (S)
10.704min (+0.047) 0.78ng/ul
response 8893

Ion	Exp%	Act%
131.00	100	100
133.00	33.60	28.29
69.00	19.50	16.53
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM111116\
 Data File : BM007831.D
 Acq On : 11 Nov 2016 18:24
 Operator : UM/SJ
 Sample : H5612-07
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F2H10

Manual Integrations
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Quant Time: Nov 12 03:37:39 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM102016.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Nov 12 03:12:55 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	170687	20.00	ng/ul	0.00
18) Naphthalene-d8	10.52	136	662343	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.37	164	422044	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.12	188	854971	20.00	ng/ul	0.00
75) Chrysene-d12	21.31	240	1090416	20.00	ng/ul	0.00
83) Perylene-d12	23.56	264	1106705	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.29	96	6140	1.54	ng/uL	0.00
5) Phenol-d5	6.92	99	92525	7.02	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.08	67	264092	33.78	ng/ul	0.00
9) 2-Chlorophenol-d4	7.27	132	281997	27.91	ng/ul	0.00
13) 4-Methylphenol-d8	8.45	113	175417	17.02	ng/ul	0.00
19) Nitrobenzene-d5	8.89	128	171530	35.99	ng/ul	0.00
22) 2-Nitrophenol-d4	9.61	143	183092	36.23	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.15	165	310409	31.24	ng/ul	0.00
29) 4-Chloroaniline-d4	10.70	131	10893m	0.96	ng/ul	0.04
43) Dimethylphthalate-d6	13.79	166	1054055	36.68	ng/ul	0.00
46) Acenaphthylene-d8	14.06	160	1238248	35.80	ng/ul	0.00
51) 4-Nitrophenol-d4	14.63	143	21012	4.62	ng/ul	0.04
57) Fluorene-d10	15.37	176	913383	36.27	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.51	200	126532	24.91	ng/ul	0.00
70) Anthracene-d10	17.22	188	1399778	36.33	ng/ul	0.00
76) Pyrene-d10	19.52	212	1668508	37.24	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.42	264	1785030	36.94	ng/ul	0.00

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

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