

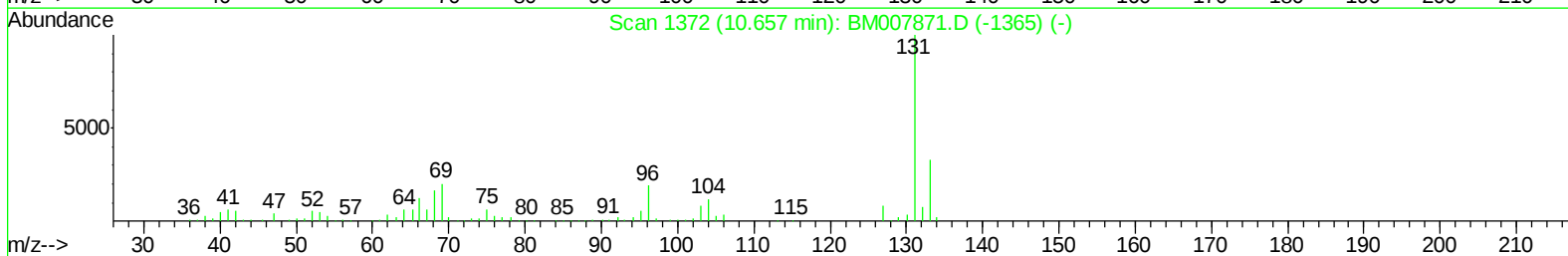
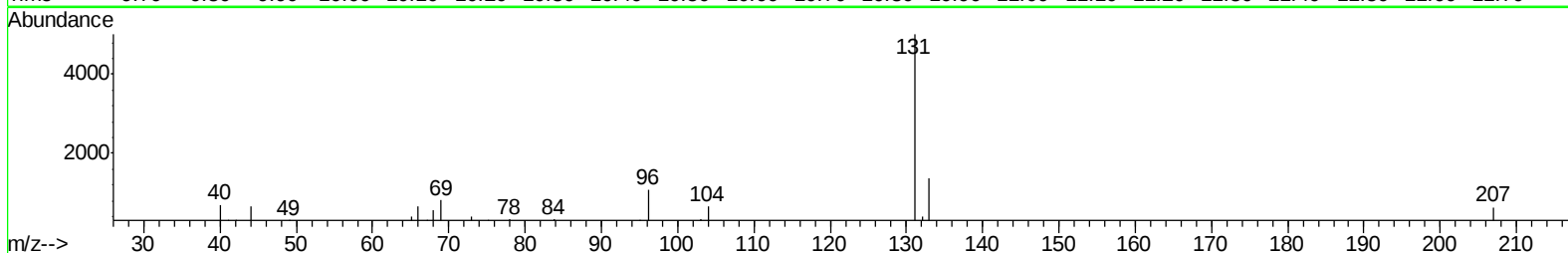
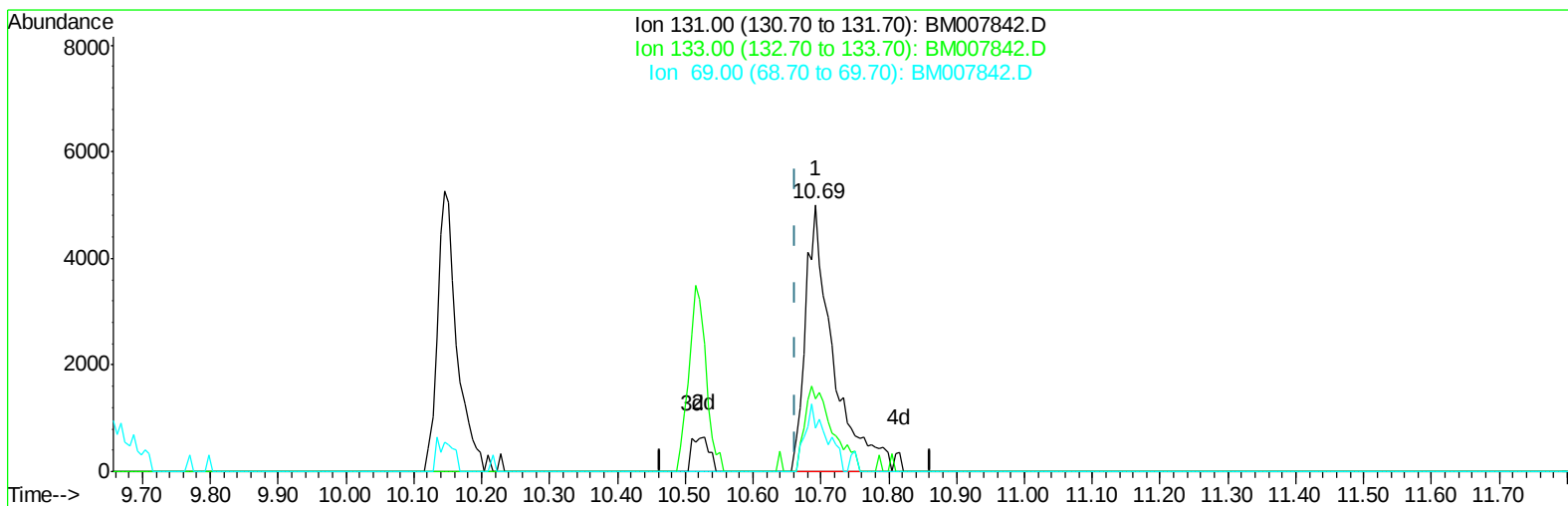
Data Path : Z:\HPCHEM1\BNA_M\Data\BM111116\
Data File : BM007842.D
Acq On : 12 Nov 2016 00:52
Operator : UM/SJ
Sample : H5552-07
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AA0

Manual Integrations
APPROVED

sohil
11/14/2016 6:38:48 PM

Quant Time: Nov 12 04:10:13 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



TIC: BM007842.D

(29) 4-Chloroaniline-d4 (S)

10.692min (+0.029) 1.15ng/ul m

response 14166

Ion	Exp%	Act%
131.00	100	100
133.00	33.60	27.47
69.00	19.50	16.31
0.00	0.00	0.00

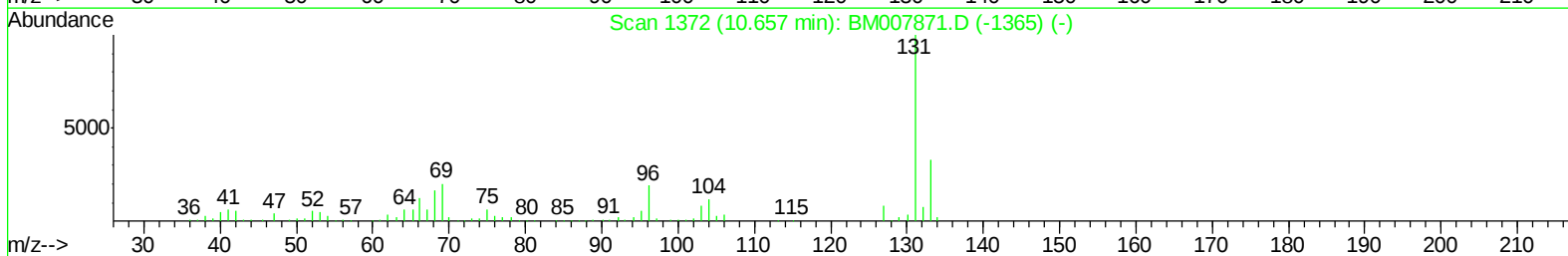
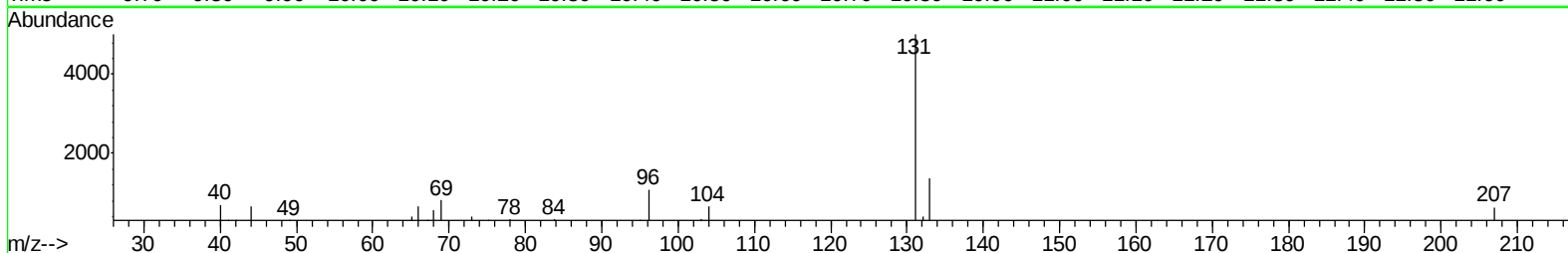
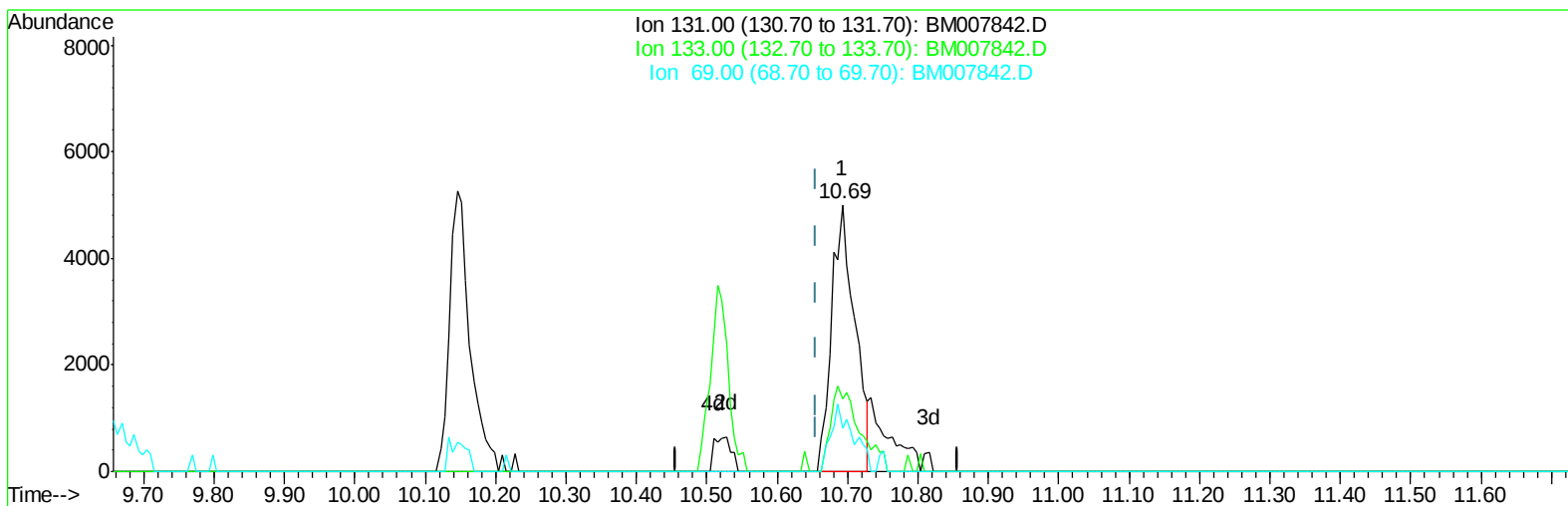
Data Path : Z:\HPCHEM1\BNA_M\Data\BM111116\
Data File : BM007842.D
Acq On : 12 Nov 2016 00:52
Operator : UM/SJ
Sample : H5552-07
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AA0

Manual Integrations
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Quant Time: Nov 16 04:55:53 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: BM007842.D

(29) 4-Chloroaniline-d4 (S)
10.692min (+0.035) 0.93ng/ul
response 11440

Ion	Exp%	Act%
131.00	100	100
133.00	33.60	27.47
69.00	19.50	16.31
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_M\Data\BM111116\
 Data File : BM007842.D
 Acq On : 12 Nov 2016 00:52
 Operator : UM/SJ
 Sample : H5552-07
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
 C0AA0

Manual Integrations
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Quant Time: Nov 12 04:10:13 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	7.73	152	183580	20.00	ng/ul	0.00
18) Naphthalene-d8	10.52	136	719249	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.37	164	466462	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.12	188	944467	20.00	ng/ul	0.00
75) Chrysene-d12	21.31	240	1195204	20.00	ng/ul	0.00
83) Perylene-d12	23.56	264	1226207	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.28	96	6126	1.43	ng/uL	0.00
5) Phenol-d5	6.92	99	92126	6.50	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.07	67	249087	29.62	ng/ul	0.00
9) 2-Chlorophenol-d4	7.27	132	280850	25.84	ng/ul	0.00
13) 4-Methylphenol-d8	8.45	113	174403	15.73	ng/ul	0.00
19) Nitrobenzene-d5	8.89	128	173828	33.58	ng/ul	0.00
22) 2-Nitrophenol-d4	9.61	143	185985	33.89	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.15	165	311994	28.92	ng/ul	0.00
29) 4-Chloroaniline-d4	10.69	131	14166m	1.15	ng/ul	0.03
43) Dimethylphthalate-d6	13.79	166	1074464	33.83	ng/ul	0.00
46) Acenaphthylene-d8	14.06	160	1268614	33.19	ng/ul	0.00
51) 4-Nitrophenol-d4	14.62	143	22564	4.49	ng/ul	0.02
57) Fluorene-d10	15.37	176	935279	33.60	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.50	200	135525	24.15	ng/ul	0.00
70) Anthracene-d10	17.22	188	1517429	35.65	ng/ul	0.00
76) Pyrene-d10	19.52	212	1784349	36.34	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.41	264	1908441	35.65	ng/ul	0.00

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

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