

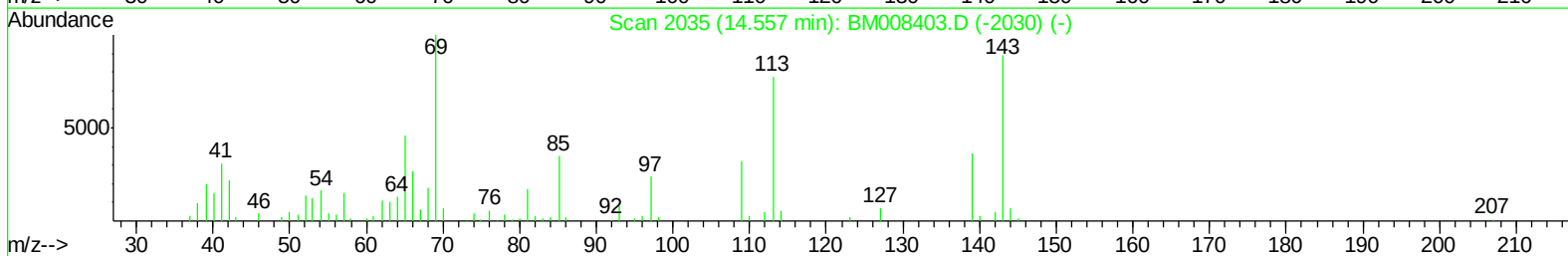
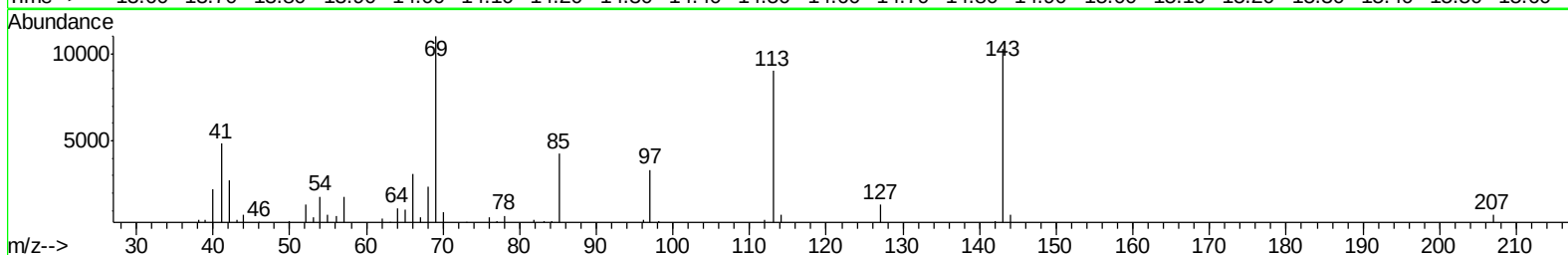
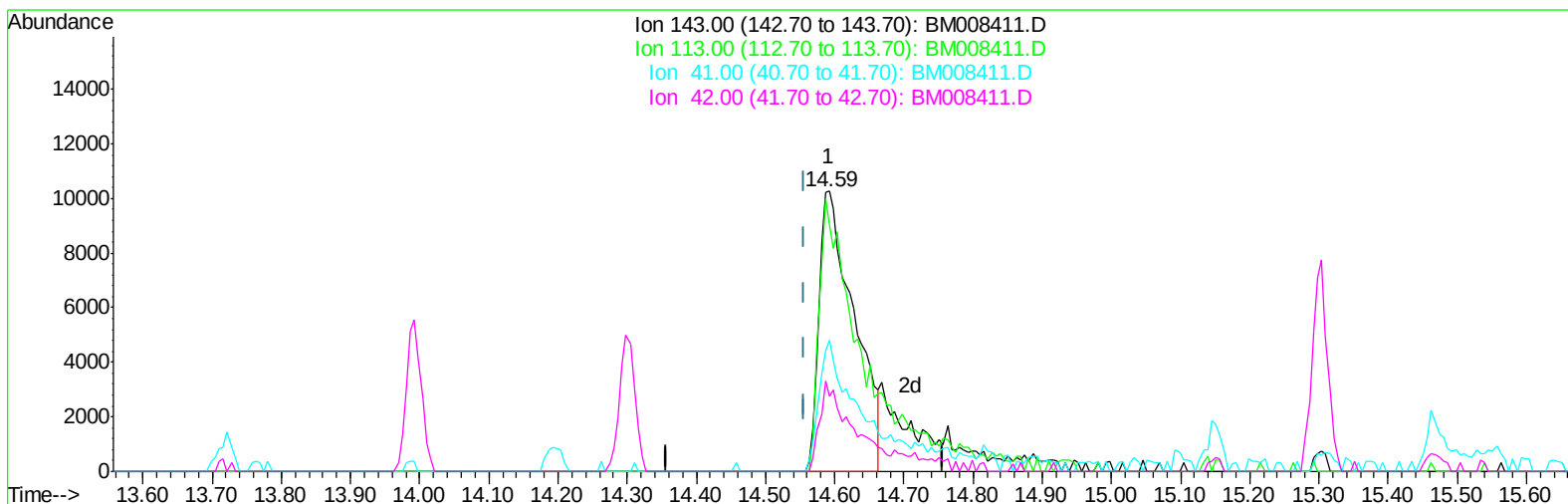
Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM121816\  
Data File : BM008411.D  
Acq On : 17 Dec 2016 14:45  
Operator : UM/SJ  
Sample : H6067-16  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
DA7G7

Quant Time: Dec 18 02:36:34 2016  
Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-2016\SOM02.2-EPA-BM12  
Quant Title : SVOA CALIBRATION  
QLast Update : Sun Dec 18 02:31:56 2016  
Response via : Initial Calibration

Manual Integrations  
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TIC: BM008411.D

(51) 4-Nitrophenol-d4 (S)

14.592min (+0.036) 13.80ng/ul

response 36356

| Ion    | Exp%   | Act%  |
|--------|--------|-------|
| 143.00 | 100    | 100   |
| 113.00 | 101.20 | 87.63 |
| 41.00  | 39.40  | 46.99 |
| 42.00  | 27.40  | 26.86 |

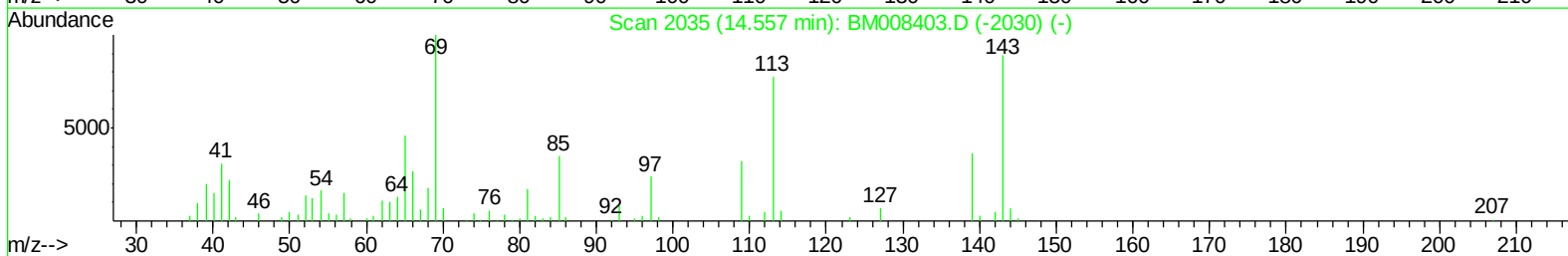
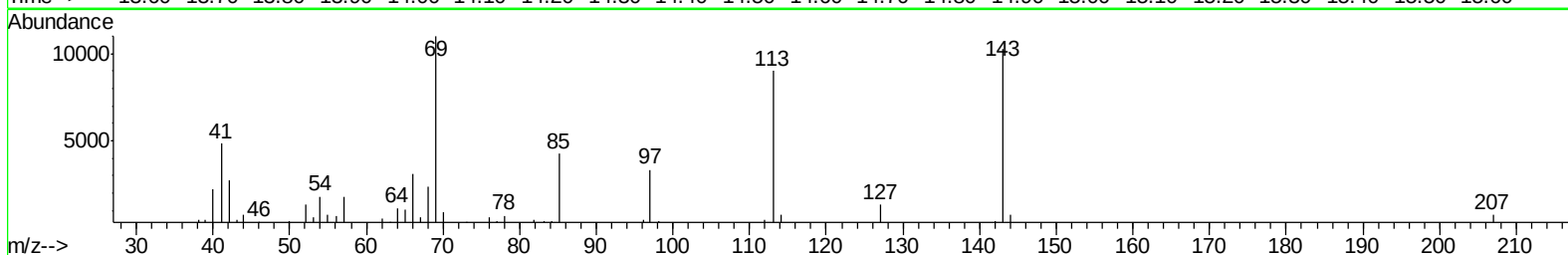
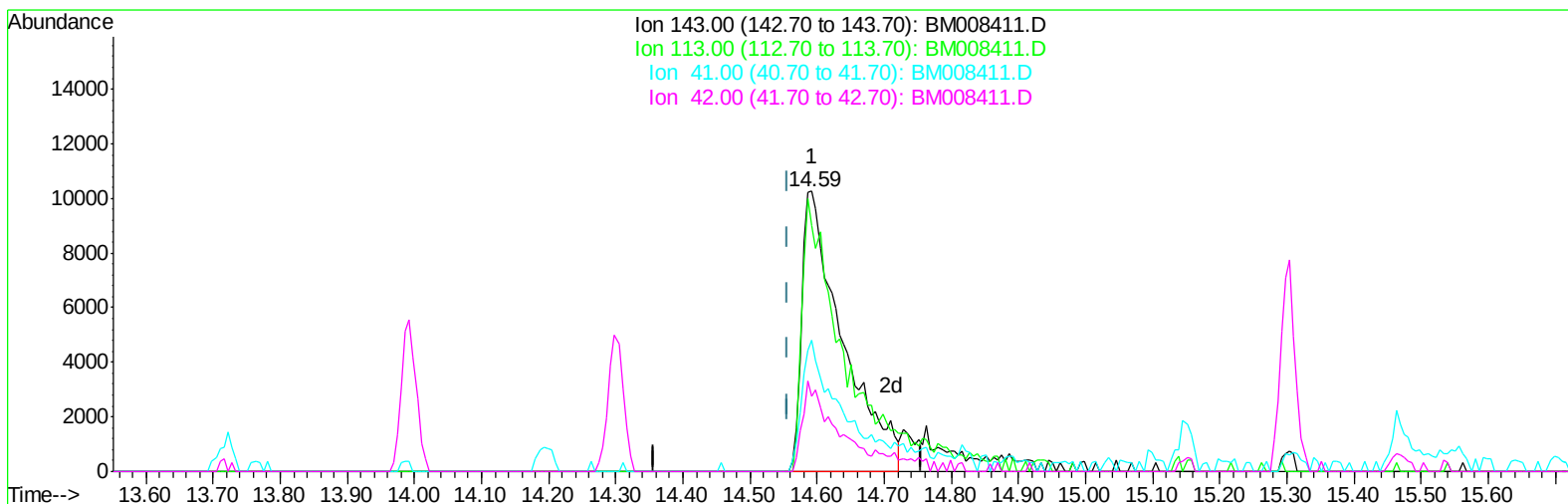
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TIC: BM008411.D

(51) 4-Nitrophenol-d4 (S)

14.592min (+0.036) 16.37ng/ul m

response 43101

| Ion    | Exp%   | Act%  |
|--------|--------|-------|
| 143.00 | 100    | 100   |
| 113.00 | 101.20 | 87.63 |
| 41.00  | 39.40  | 46.99 |
| 42.00  | 27.40  | 26.86 |

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 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 DA7G7

Manual Integrations  
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Quant Time: Dec 18 03:59:46 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-2016\SOM02.2-EPA-BM121  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sun Dec 18 02:31:56 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.66  | 152  | 108250   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.43 | 136  | 424402   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.30 | 164  | 268928   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.06 | 188  | 542531   | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.26 | 240  | 678624   | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 23.49 | 264  | 687691   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.23  | 96  | 15093  | 6.56  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.85  | 99  | 251423 | 29.94 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.00  | 67  | 156954 | 32.52 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.20  | 132 | 204474 | 32.36 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.38  | 113 | 200882 | 30.88 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.82  | 128 | 97253  | 31.71 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.54  | 143 | 103038 | 30.32 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.08 | 165 | 191385 | 30.73 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.60 | 131 | 246543 | 34.53 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 13.72 | 166 | 592430 | 33.14 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 13.99 | 160 | 738982 | 34.98 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.59 | 143 | 43101m | 16.37 | ng/ul | 0.04 |
| 57) Fluorene-d10               | 15.30 | 176 | 511909 | 32.94 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.47 | 200 | 49174  | 15.23 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.16 | 188 | 797811 | 33.75 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.46 | 212 | 951010 | 36.29 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 23.34 | 264 | 990880 | 34.31 | ng/ul | 0.00 |

| Target Compounds      |       |     |        |      |       | Qvalue |
|-----------------------|-------|-----|--------|------|-------|--------|
| 6) Phenol             | 6.88  | 94  | 52441  | 6.01 | ng/ul | 95     |
| 44) Dimethylphthalate | 13.77 | 163 | 153370 | 8.75 | ng/ul | 100    |

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

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