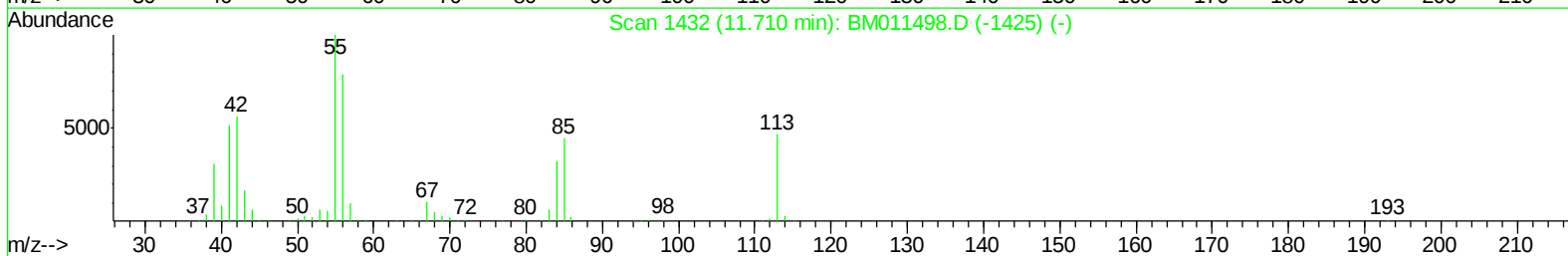
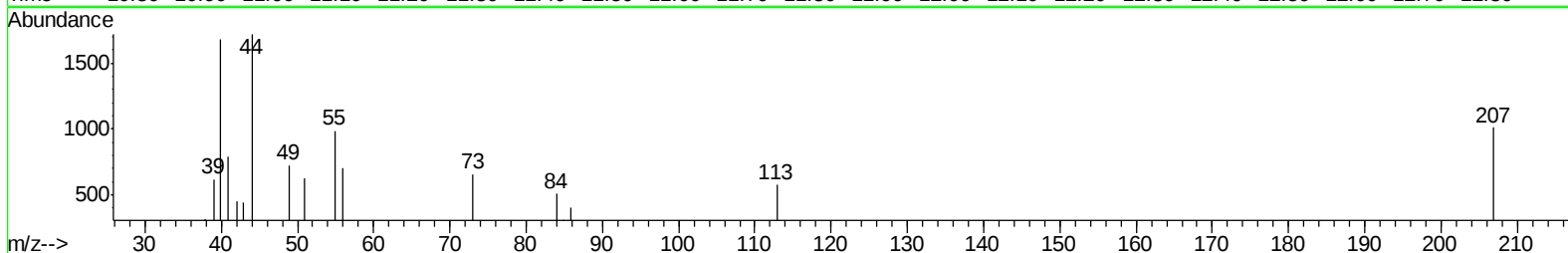
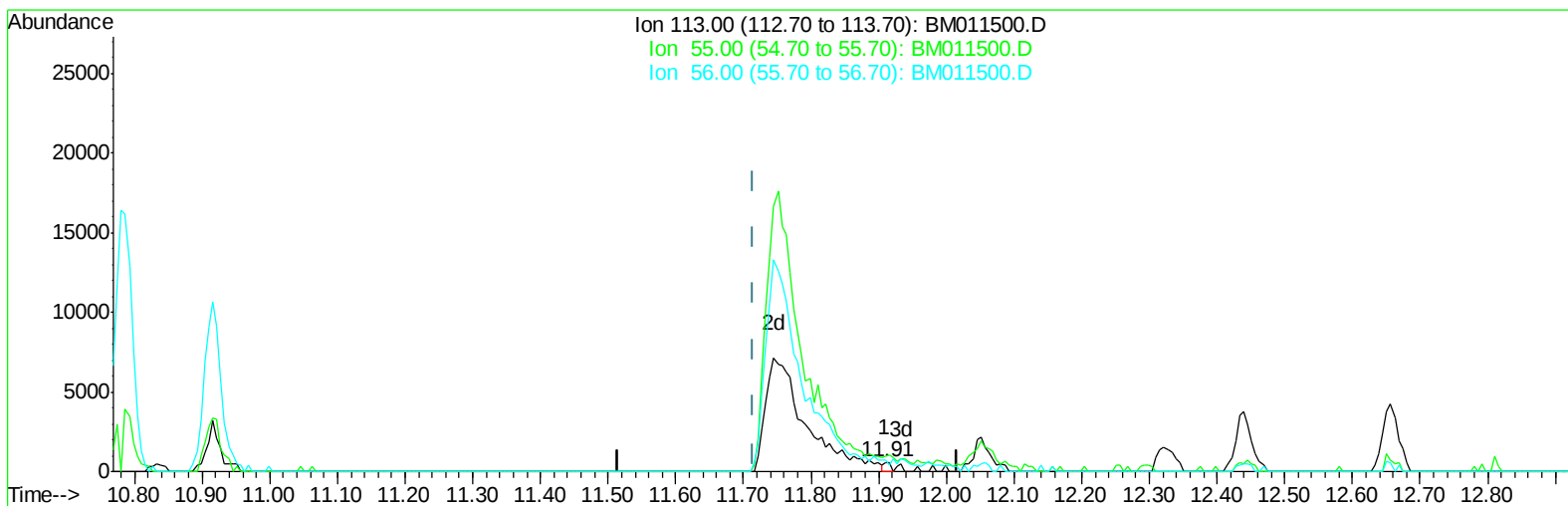


Data Path : Z:\HPCHEM1\BNA\_M\Data\BM091117\  
Data File : BM011500.D  
Acq On : 11 Sep 2017 17:14  
Operator : SJ/JU  
Sample : MDL-S-ME-02  
Misc :  
ALS Vial : 13 Sample Multiplier: 1

Quant Time: Sep 11 18:06:25 2017  
Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM090817MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Sep 11 18:01:51 2017  
Response via : Initial Calibration



TIC: BM011500.D

(32) Caprolactam

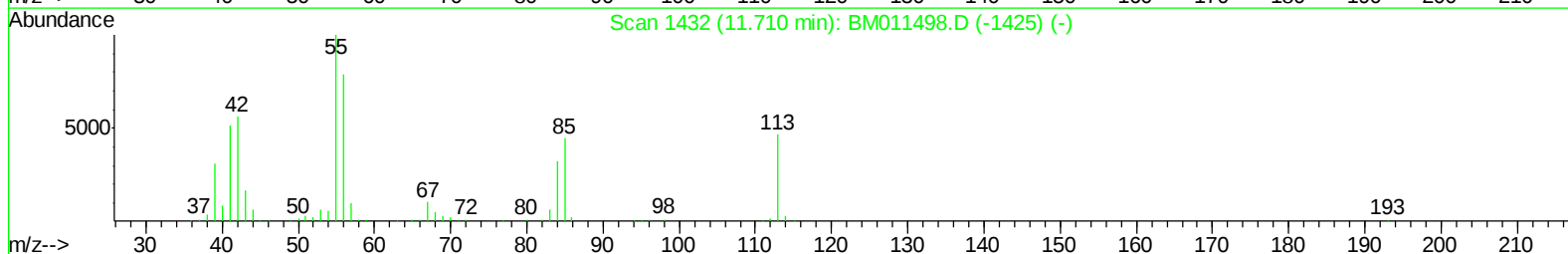
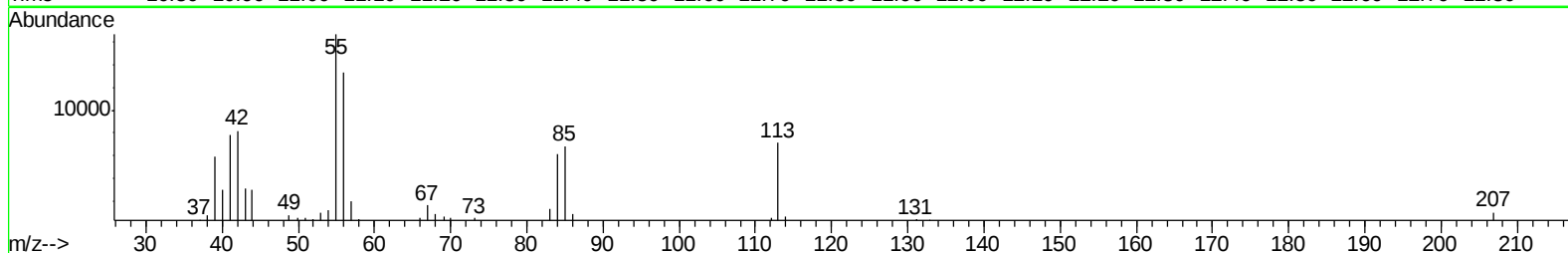
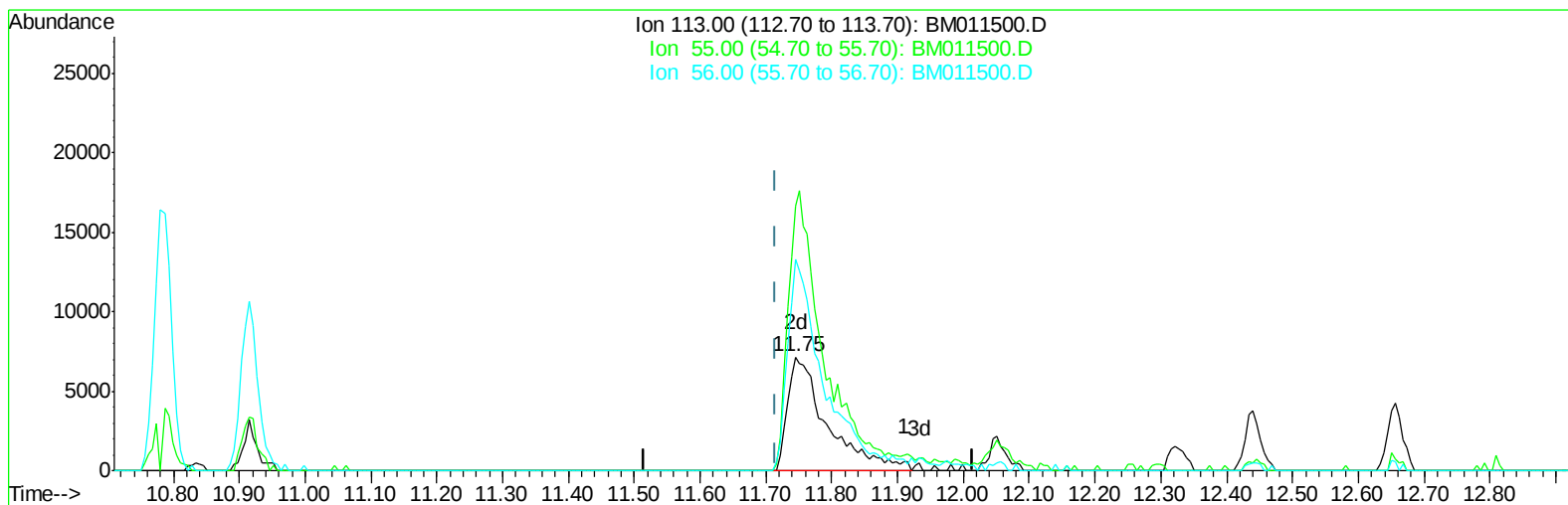
11.910min (+0.194) 0.04ng/ul

response 389

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 164.60 | 171.65 |
| 56.00  | 115.90 | 122.96 |
| 0.00   | 0.00   | 0.00   |

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Quant Title : SVOA CALIBRATION  
QLast Update : Mon Sep 11 18:01:51 2017  
Response via : Initial Calibration



TIC: BM011500.D

(32) Caprolactam

11.745min (+0.029) 3.13ng/ul m

response 29809

| Ion    | Exp%   | Act%    |
|--------|--------|---------|
| 113.00 | 100    | 100     |
| 55.00  | 164.60 | 233.13# |
| 56.00  | 115.90 | 185.73# |
| 0.00   | 0.00   | 0.00    |

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 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Sep 11 18:01:51 2017  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.97  | 152  | 437671   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.79 | 136  | 1978317  | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.61 | 164  | 1181242  | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.34 | 188  | 2676793  | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.51 | 240  | 3035456  | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.88 | 264  | 2842139  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.40  | 96  | 60559   | 6.22  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.12  | 99  | 1304610 | 31.04 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.30  | 67  | 950733  | 33.57 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.50  | 132 | 1028634 | 32.48 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.67  | 113 | 1052108 | 32.07 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.14  | 128 | 518725  | 34.54 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.86  | 143 | 521747  | 33.42 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.40 | 165 | 1000383 | 31.83 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.92 | 131 | 1442775 | 41.60 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 14.01 | 166 | 3445403 | 34.50 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 14.30 | 160 | 4179816 | 34.61 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.79 | 143 | 535031  | 28.51 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.60 | 176 | 2948620 | 34.09 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.70 | 200 | 429333  | 27.56 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 17.44 | 188 | 4537081 | 34.42 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.73 | 212 | 5069695 | 35.67 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 23.73 | 264 | 4727532 | 34.90 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |              | Qvalue |
|--------------------------------|-------|-----|--------|--------------|--------|
| 2) 1,4-Dioxane                 | 3.43  | 88  | 18394  | 1.725 ng/uL# | 76     |
| 4) Benzaldehyde                | 7.12  | 77  | 56419  | 2.349 ng/ul  | 96     |
| 6) Phenol                      | 7.15  | 94  | 151839 | 3.640 ng/ul  | 98     |
| 8) Bis(2-Chloroethyl)ether     | 7.39  | 93  | 130615 | 3.977 ng/ul# | 86     |
| 10) 2-Chlorophenol             | 7.53  | 128 | 116954 | 3.785 ng/ul  | 93     |
| 11) 2-Methylphenol             | 8.41  | 108 | 108661 | 3.435 ng/ul  | 90     |
| 12) 2,2'-oxybis(1-Chloropropan | 8.51  | 45  | 218038 | 4.096 ng/ul  | 93     |
| 14) Acetophenone               | 8.80  | 105 | 187189 | 3.733 ng/ul# | 93     |
| 15) N-Nitroso-di-n-propylamine | 8.78  | 70  | 101215 | 3.761 ng/ul# | 75     |
| 16) 4-Methylphenol             | 8.74  | 108 | 123110 | 3.557 ng/ul  | 87     |
| 17) Hexachloroethane           | 9.06  | 117 | 44054  | 3.730 ng/ul  | 83     |
| 20) Nitrobenzene               | 9.19  | 77  | 139214 | 3.942 ng/ul  | 95     |
| 21) Isophorone                 | 9.70  | 82  | 264303 | 3.717 ng/ul  | 97     |
| 23) 2-Nitrophenol              | 9.89  | 139 | 62574  | 3.815 ng/ul# | 93     |
| 24) 2,4-Dimethylphenol         | 9.95  | 107 | 131264 | 3.740 ng/ul  | 87     |
| 25) Bis(2-Chloroethoxy)methane | 10.19 | 93  | 176933 | 3.816 ng/ul  | 95     |
| 27) 2,4-Dichlorophenol         | 10.42 | 162 | 112441 | 3.711 ng/ul# | 88     |
| 28) Naphthalene                | 10.83 | 128 | 403300 | 3.931 ng/ul  | 98     |
| 30) 4-Chloroaniline            | 10.94 | 127 | 147739 | 4.483 ng/ul  | 96     |
| 31) Hexachlorobutadiene        | 11.12 | 225 | 67727  | 3.938 ng/ul  | 92     |
| 32) Caprolactam                | 11.75 | 113 | 29809m | 3.125 ng/ul  |        |
| 33) 4-Chloro-3-methylphenol    | 12.05 | 107 | 108623 | 3.381 ng/ul  | 96     |
| 34) 2-Methylnaphthalene        | 12.44 | 142 | 288334 | 3.785 ng/ul  | 97     |

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 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Sep 11 18:01:51 2017  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc  | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|-------|--------|----------|
| 35) 1-Methylnaphthalene        | 12.66 | 142  | 276739   | 3.814 | ng/ul# | 95       |
| 37) 1,2,4,5-Tetrachlorobenzene | 12.80 | 216  | 136616   | 3.899 | ng/ul# | 94       |
| 38) Hexachlorocyclopentadiene  | 12.78 | 237  | 52404    | 2.668 | ng/ul# | 97       |
| 39) 2,4,6-Trichlorophenol      | 13.04 | 196  | 78895    | 3.307 | ng/ul  | 98       |
| 40) 2,4,5-Trichlorophenol      | 13.11 | 196  | 77123    | 3.188 | ng/ul  | 97       |
| 41) 1,1'-Biphenyl              | 13.44 | 154  | 379042   | 3.856 | ng/ul# | 93       |
| 42) 2-Chloronaphthalene        | 13.49 | 162  | 283217   | 3.829 | ng/ul  | 99       |
| 43) 2-Nitroaniline             | 13.69 | 65   | 70492    | 3.069 | ng/ul# | 83       |
| 45) Dimethylphthalate          | 14.06 | 163  | 373109   | 3.902 | ng/ul# | 97       |
| 46) 2,6-Dinitrotoluene         | 14.18 | 165  | 52172    | 3.090 | ng/ul  | 89       |
| 48) Acenaphthylene             | 14.33 | 152  | 476146   | 3.917 | ng/ul  | 96       |
| 49) 3-Nitroaniline             | 14.51 | 138  | 53444    | 2.567 | ng/ul# | 99       |
| 50) Acenaphthene               | 14.67 | 153  | 315531   | 3.842 | ng/ul  | 97       |
| 51) 2,4-Dinitrophenol          | 14.71 | 184  | 22206    | 2.129 | ng/ul# | 74       |
| 53) 4-Nitrophenol              | 14.80 | 109  | 41303    | 3.424 | ng/ul# | 34       |
| 54) Dibenzofuran               | 15.00 | 168  | 450500   | 3.921 | ng/ul  | 95       |
| 55) 2,4-Dinitrotoluene         | 14.96 | 165  | 79876    | 3.218 | ng/ul# | 73       |
| 56) 2,3,4,6-Tetrachlorophenol  | 15.23 | 232  | 72908    | 3.307 | ng/ul# | 77       |
| 57) Diethylphthalate           | 15.42 | 149  | 373952   | 3.758 | ng/ul  | 96       |
| 59) Fluorene                   | 15.65 | 166  | 374033   | 3.885 | ng/ul  | 99       |
| 60) 4-Chlorophenyl-phenylether | 15.64 | 204  | 174407   | 3.866 | ng/ul# | 90       |
| 61) 4-Nitroaniline             | 15.67 | 138  | 62665    | 2.810 | ng/ul# | 79       |
| 64) 4,6-Dinitro-2-methylphenol | 15.72 | 198  | 53849    | 3.374 | ng/ul# | 48       |
| 65) N-Nitrosodiphenylamine     | 15.85 | 169  | 302694   | 3.712 | ng/ul  | 96       |
| 66) 4-Bromophenyl-phenylether  | 16.53 | 248  | 100520   | 3.677 | ng/ul  | 93       |
| 67) Hexachlorobenzene          | 16.65 | 284  | 113225   | 3.762 | ng/ul# | 92       |
| 68) Atrazine                   | 16.80 | 200  | 89976    | 3.185 | ng/ul  | 98       |
| 69) Pentachlorophenol          | 16.99 | 266  | 50099    | 2.593 | ng/ul  | 87       |
| 70) Phenanthrene               | 17.39 | 178  | 579545   | 3.884 | ng/ul  | 99       |
| 72) Anthracene                 | 17.48 | 178  | 597300   | 3.901 | ng/ul  | 98       |
| 73) 1,2,3,4-Tetrachlorobenzene | 13.41 | 216  | 136056   | 3.814 | ng/uL  | 96       |
| 74) Pentachlorobenzene         | 14.92 | 250  | 140085   | 3.872 | ng/uL  | 97       |
| 75) Carbazole                  | 17.74 | 167  | 492275   | 3.769 | ng/ul  | 98       |
| 76) Di-n-butylphthalate        | 18.30 | 149  | 598386   | 3.488 | ng/ul  | 99       |
| 77) Fluoranthene               | 19.39 | 202  | 632657   | 4.123 | ng/ul# | 88       |
| 80) Pyrene                     | 19.76 | 202  | 724018   | 4.088 | ng/ul# | 83       |
| 81) Butylbenzylphthalate       | 20.64 | 149  | 256073   | 3.144 | ng/ul  | 92       |
| 82) 3,3'-Dichlorobenzidine     | 21.42 | 252  | 163317   | 2.963 | ng/ul# | 96       |
| 83) Benzo(a)anthracene         | 21.49 | 228  | 653988   | 3.913 | ng/ul  | 99       |
| 84) Bis(2-ethylhexyl)phthalate | 21.40 | 149  | 396540   | 3.170 | ng/ul# | 95       |
| 85) Chrysene                   | 21.54 | 228  | 598055   | 3.947 | ng/ul  | 99       |
| 87) Di-n-octyl phthalate       | 22.32 | 149  | 666426   | 3.304 | ng/ul  | 100      |
| 88) Benzo(b)fluoranthene       | 23.16 | 252  | 614148   | 3.592 | ng/ul# | 96       |
| 89) Benzo(k)fluoranthene       | 23.21 | 252  | 614153   | 3.740 | ng/ul# | 94       |
| 91) Benzo(a)pyrene             | 23.78 | 252  | 617337   | 3.825 | ng/ul# | 95       |
| 92) Indeno(1,2,3-cd)pyrene     | 26.32 | 276  | 643751   | 3.626 | ng/ul# | 85       |
| 93) Dibenzo(a,h)anthracene     | 26.33 | 278  | 531361   | 3.530 | ng/ul# | 94       |
| 94) Benzo(g,h,i)perylene       | 27.06 | 276  | 535317   | 3.723 | ng/ul# | 89       |

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QLast Update : Mon Sep 11 18:01:51 2017  
Response via : Initial Calibration

| Internal Standards                                                         | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------------------------------------------------------|------|------|----------|------|-------|----------|
| -----                                                                      |      |      |          |      |       |          |
| (#) = qualifier out of range (m) = manual integration (+) = signals summed |      |      |          |      |       |          |

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