

(QT Reviewed)

**Instrument :**  
BNA\_G  
**LabSampleId :**  
SSTD02009

UMANGI  
12/16/2016 6:21:15 PM

[illegible]

# Quantitation Report (Qedit)

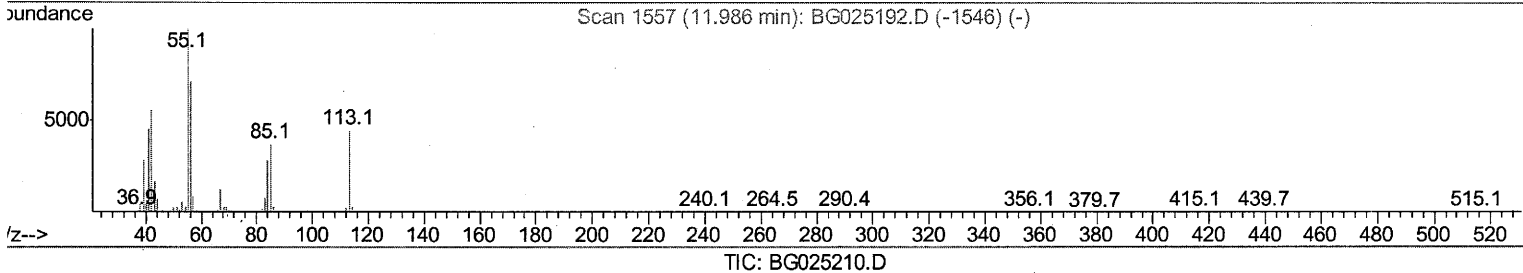
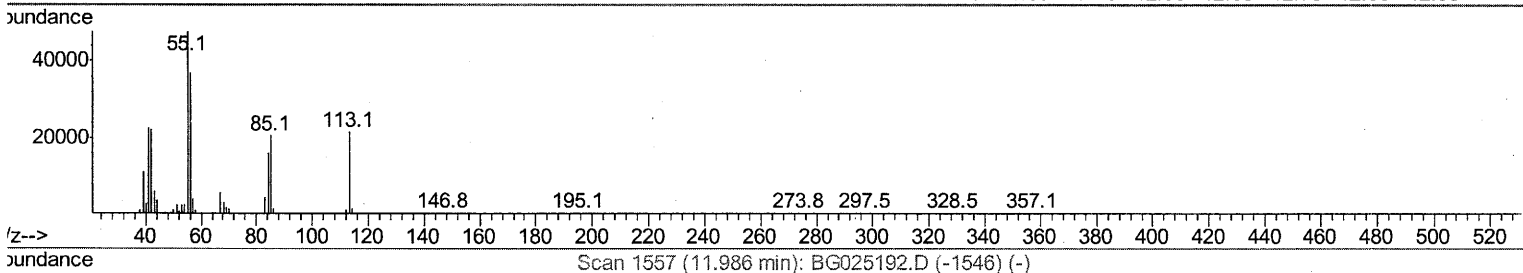
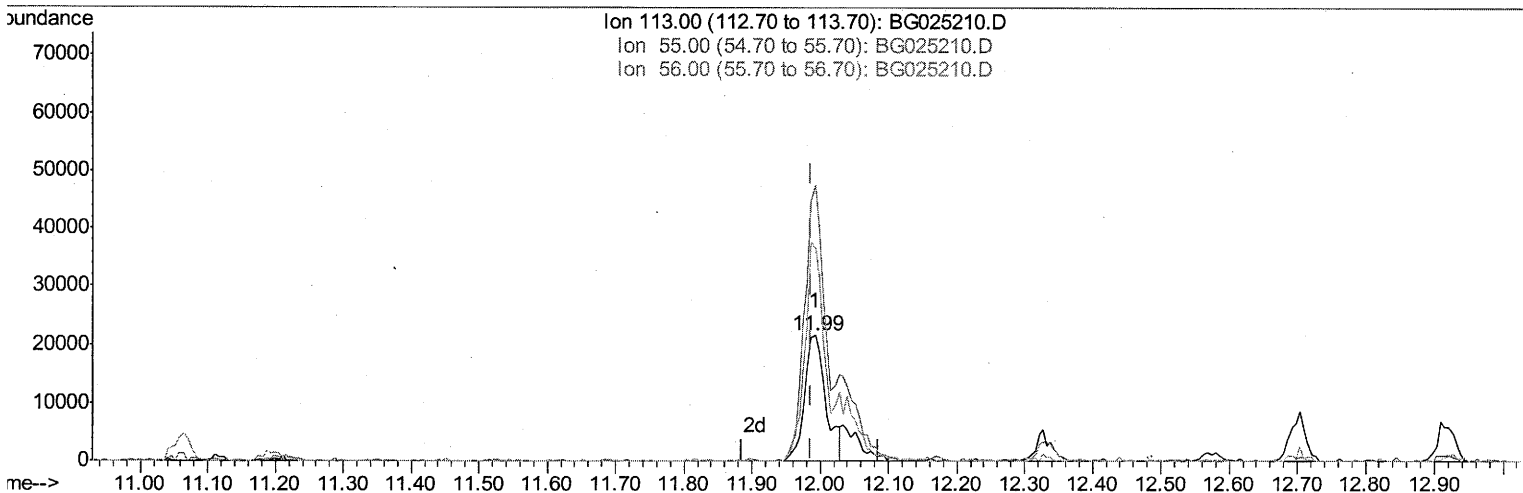
Data Path : Z:\HPCHEM1\BNA G\Data\BG121516\  
 Data File : BG025210.D  
 Acq On : 15 Dec 2016 14:03  
 Operator : UM/SJ  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 LabSampled :  
 SSTD02009

Manual Integrations  
 APPROVED

UMANGI  
 12/16/2016 6:21:15 PM

Quant Time: Dec 16 03:07:08 2016  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG113016.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Dec 15 04:58:54 2016  
 Response via : Initial Calibration



TIC: BG025210.D

(32) Caprolactam

11.992min (+0.006) 14.05ng/ul

response 46679

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 210.70 | 220.59 |
| 56.00  | 160.60 | 170.77 |
| 0.00   | 0.00   | 0.00   |

# Quantitation Report (Qedit)

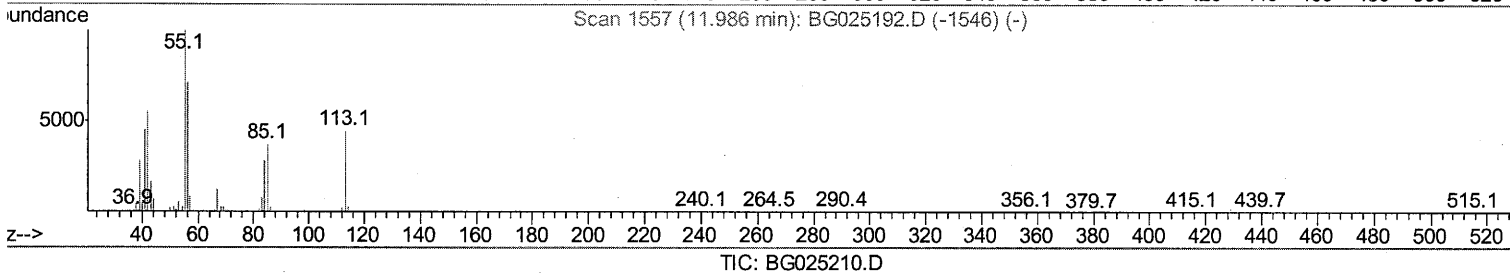
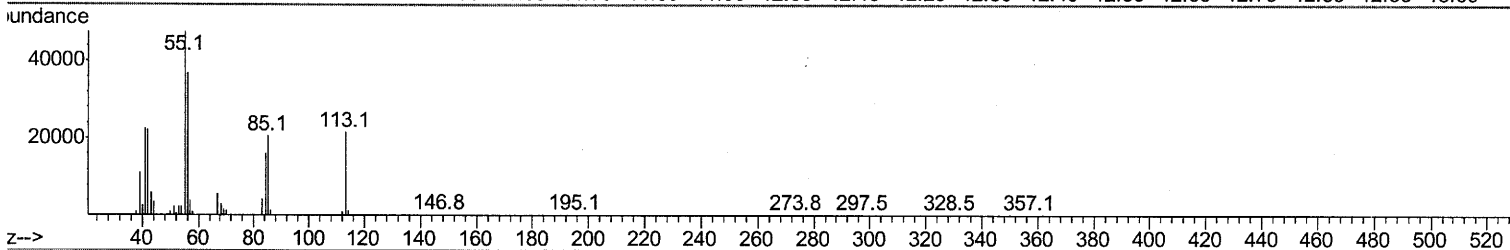
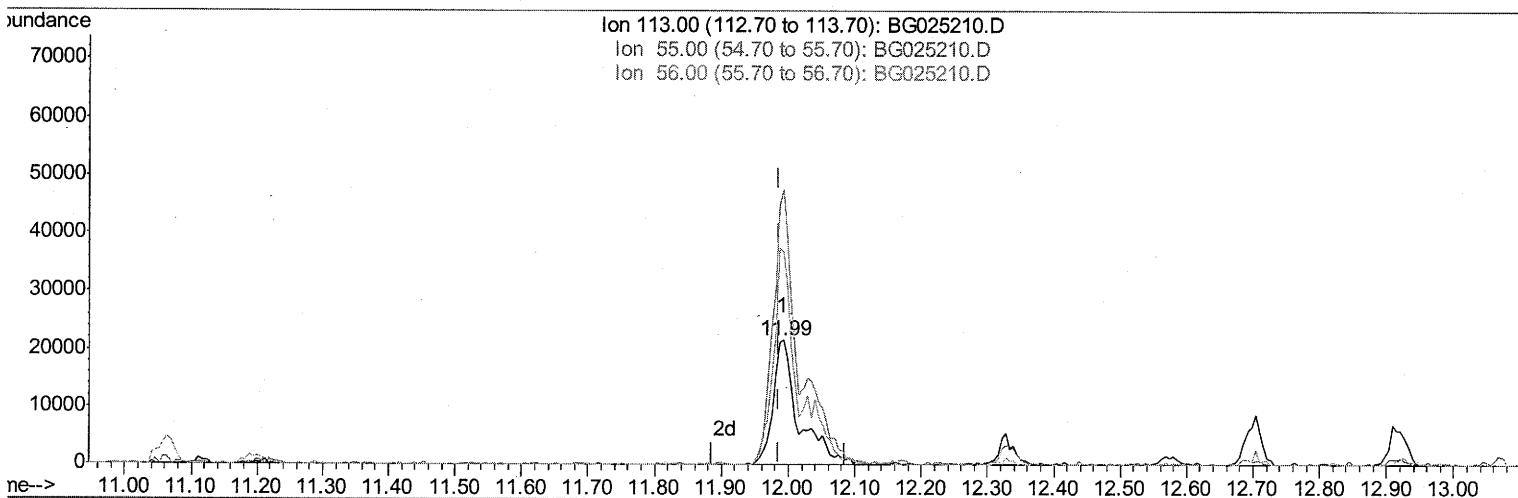
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(32) Caprolactam

11.992min (+0.006) 17.40ng/ul m

response 57804

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 210.70 | 220.59 |
| 56.00  | 160.60 | 170.77 |
| 0.00   | 0.00   | 0.00   |

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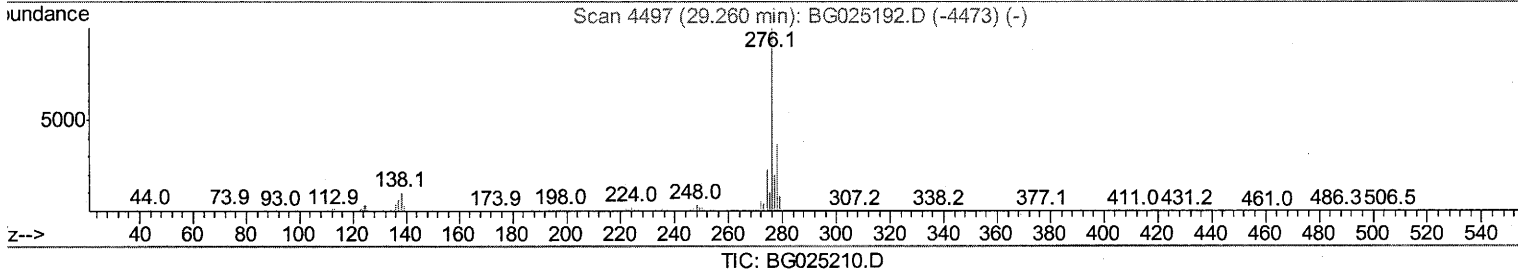
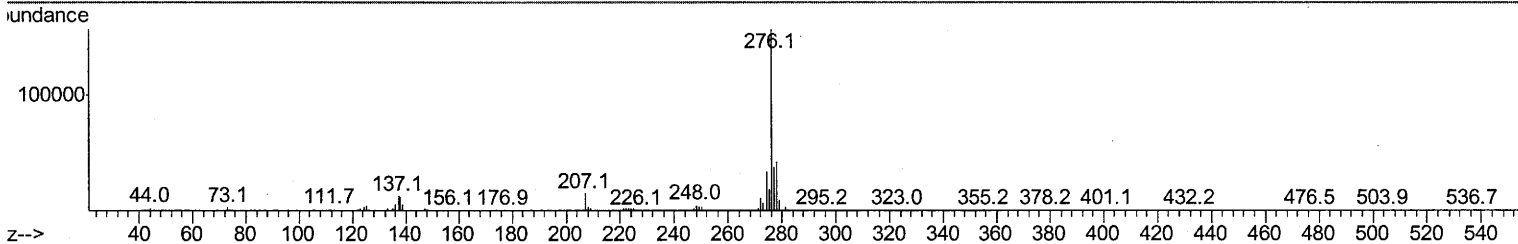
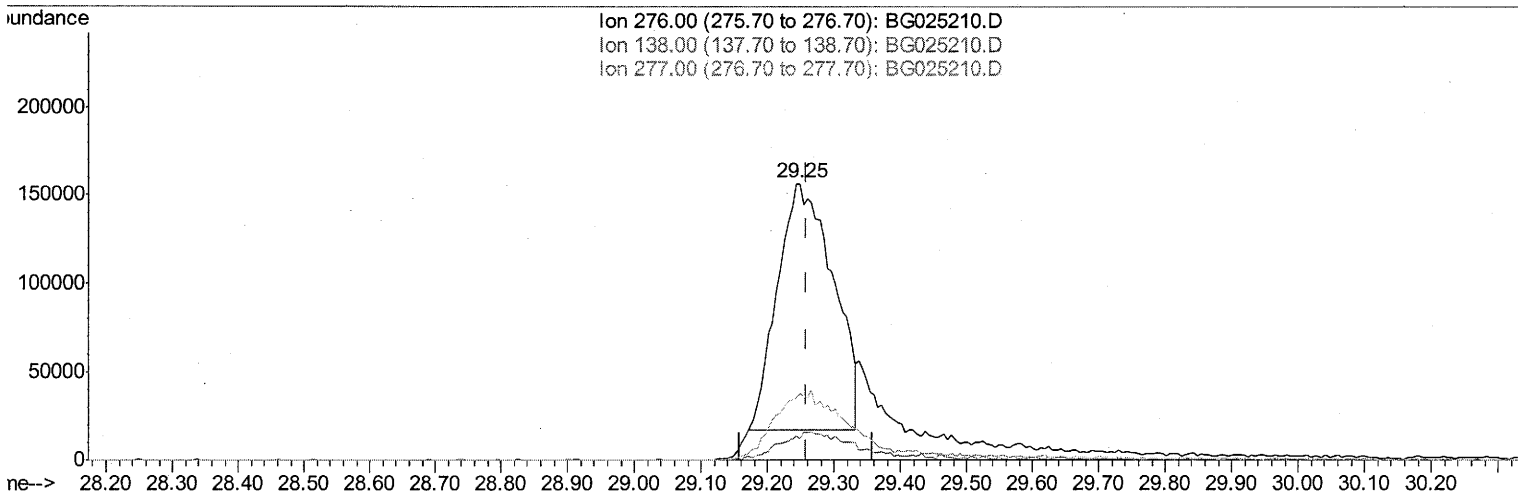
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(89) Indeno(1,2,3-cd)pyrene

29.249min (-0.011) 12.75ng/ul

response 810485

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 276.00 | 100   | 100   |
| 138.00 | 10.30 | 7.74# |
| 277.00 | 23.30 | 24.25 |
| 0.00   | 0.00  | 0.00  |

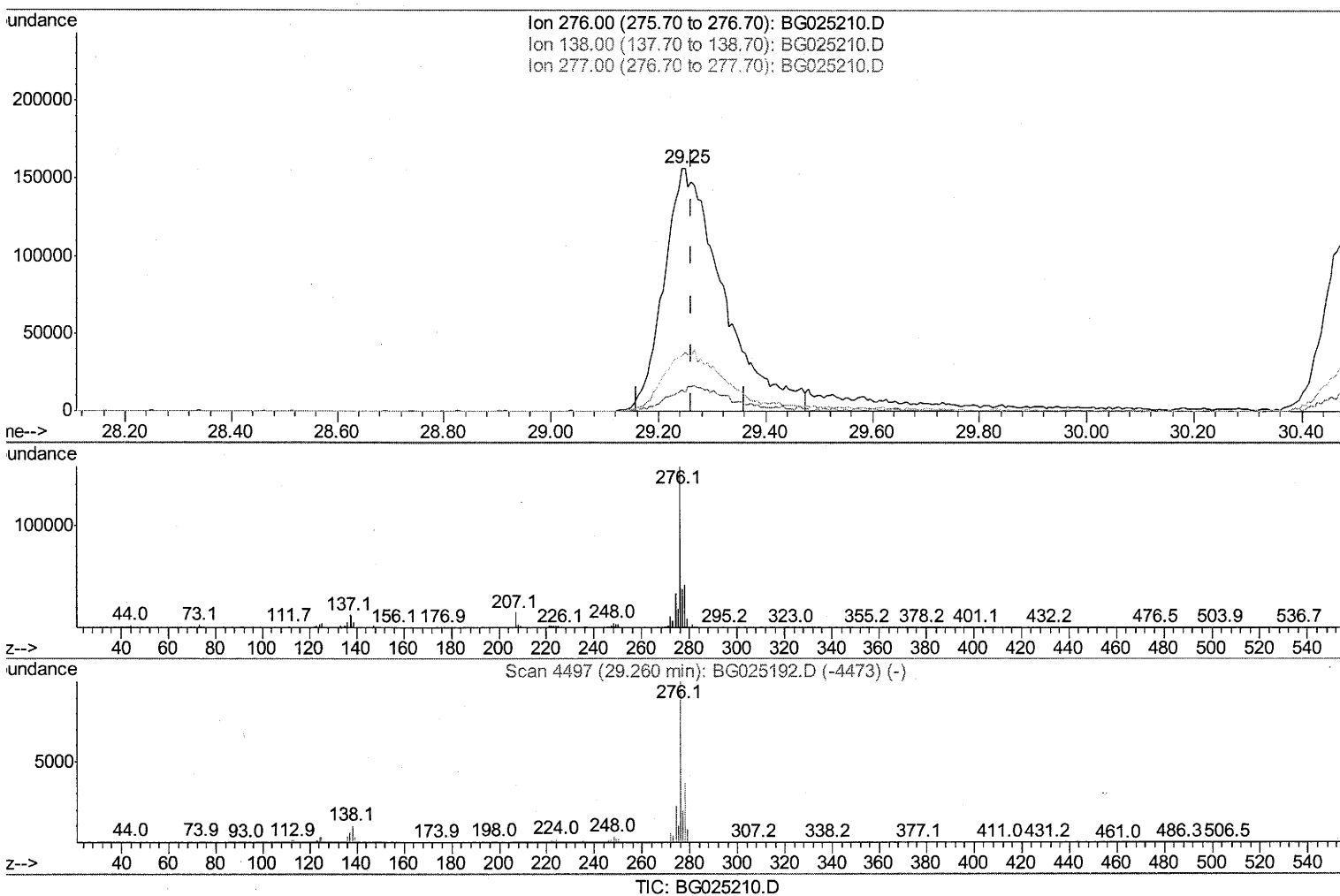
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Response via : Initial Calibration



(89) Indeno(1,2,3-cd)pyrene

29.249min (-0.011) 18.79ng/ul m

response 1194692

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 276.00 | 100   | 100   |
| 138.00 | 10.30 | 7.74# |
| 277.00 | 23.30 | 24.25 |
| 0.00   | 0.00  | 0.00  |

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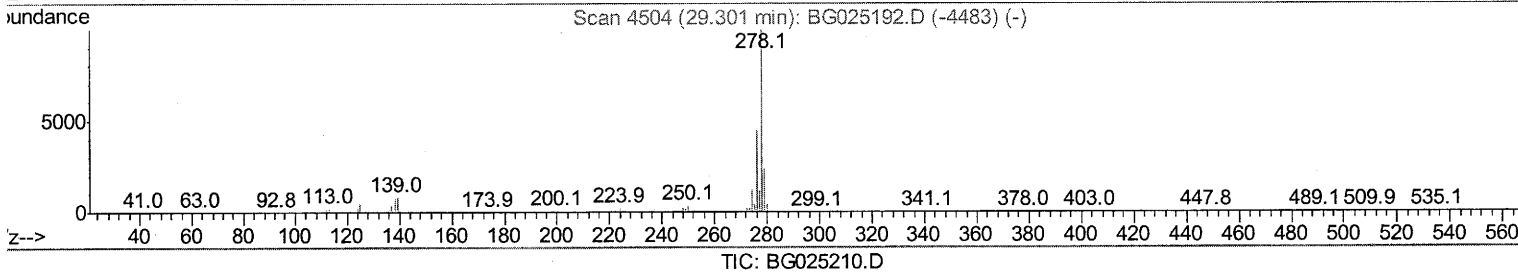
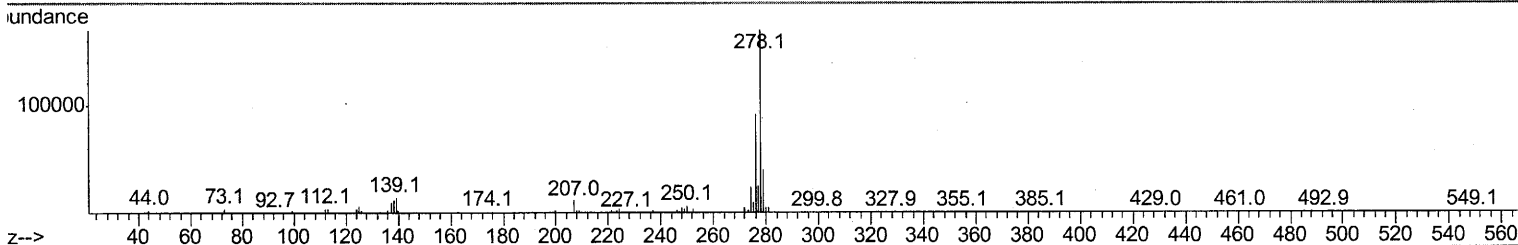
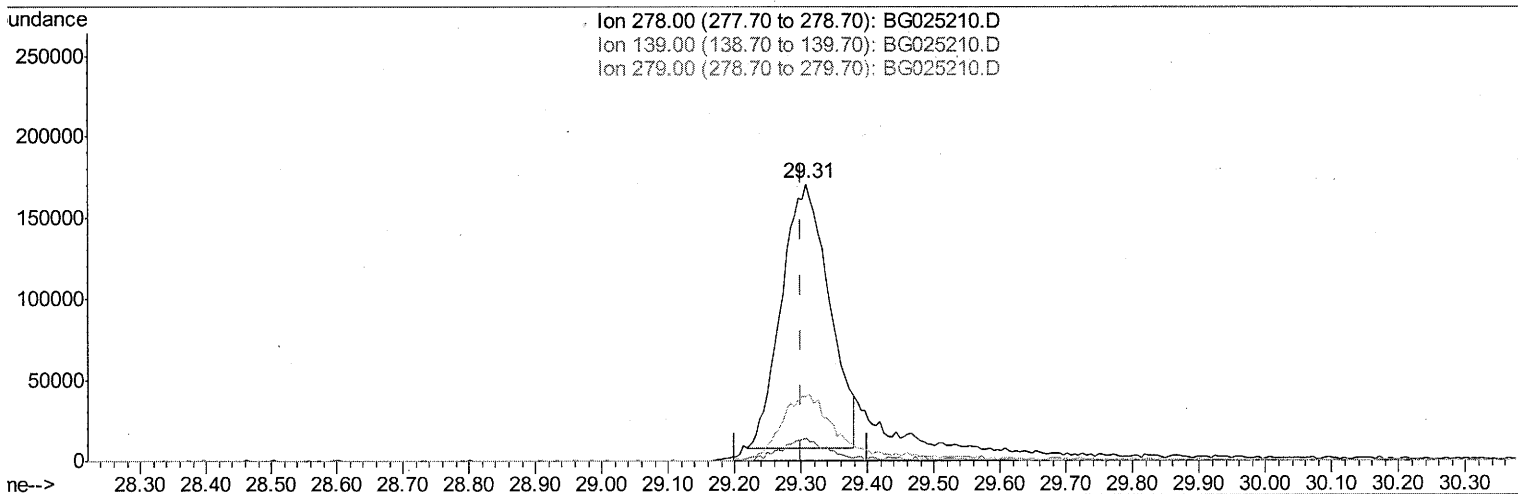
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 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

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 Response via : Initial Calibration



TIC: BG025210.D

(90) Dibenzo(a,h)anthracene

29.307min (+0.006) 15.15ng/ul

response 811804

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 278.00 | 100   | 100   |
| 139.00 | 8.40  | 10.00 |
| 279.00 | 26.10 | 23.12 |
| 0.00   | 0.00  | 0.00  |

# Quantitation Report (Qedit)

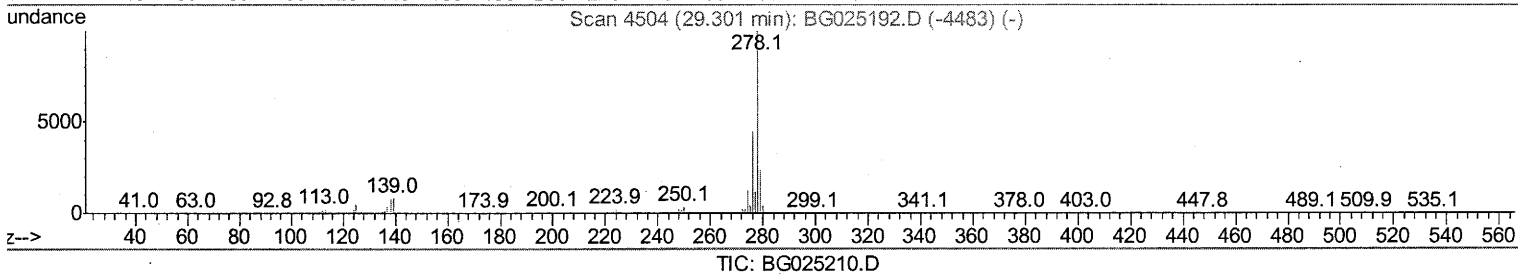
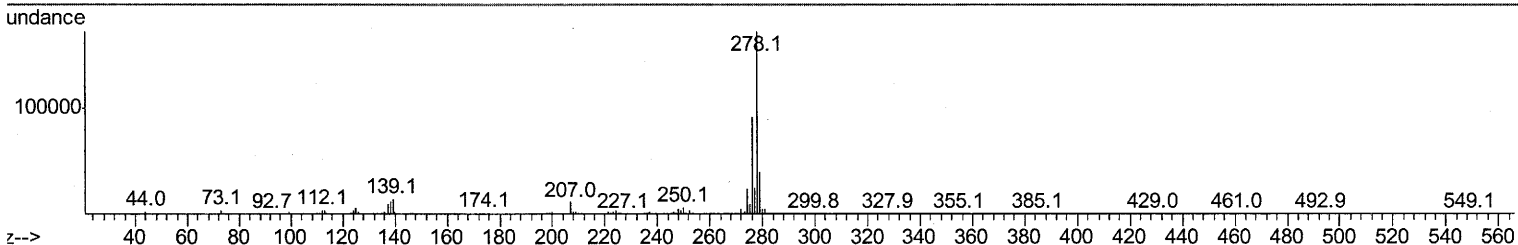
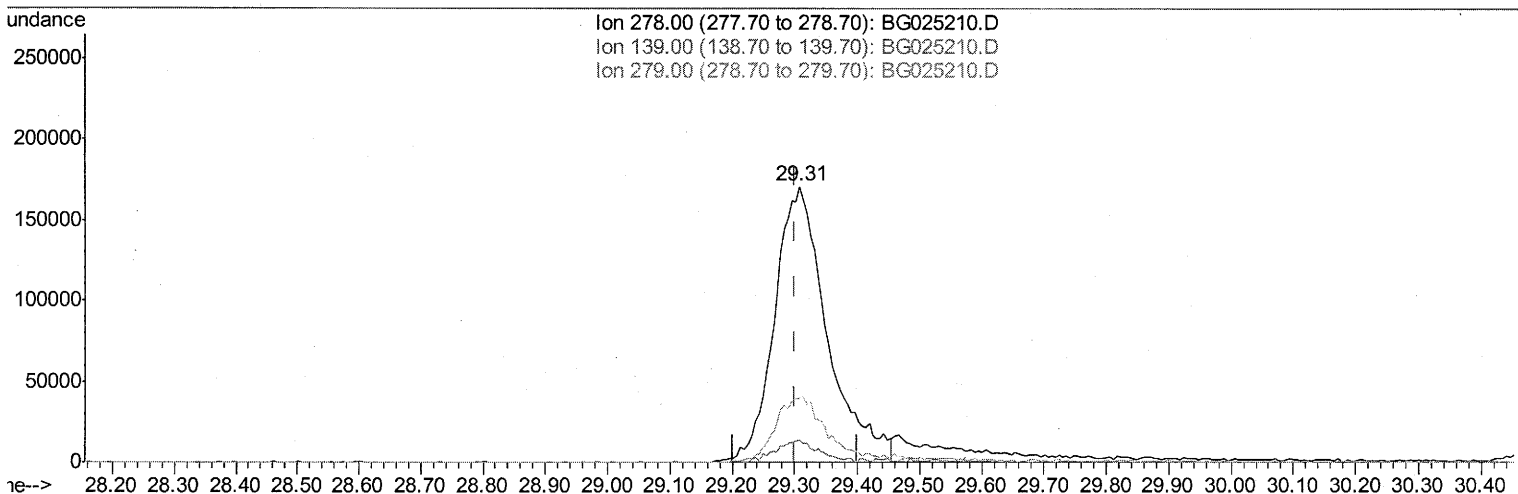
Data Path : Z:\HPCHEM1\BNA G\Data\BG121516\  
 Data File : BG025210.D  
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 Operator : UM/SJ  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

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Manual Integrations  
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 Response via : Initial Calibration



(90) Dibenzo(a,h)anthracene

29.307min (+0.006) 18.61ng/ul m

response 996899

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 278.00 | 100   | 100   |
| 139.00 | 8.40  | 8.35  |
| 279.00 | 26.10 | 23.12 |
| 0.00   | 0.00  | 0.00  |

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## Quantitation Report (Qedit)

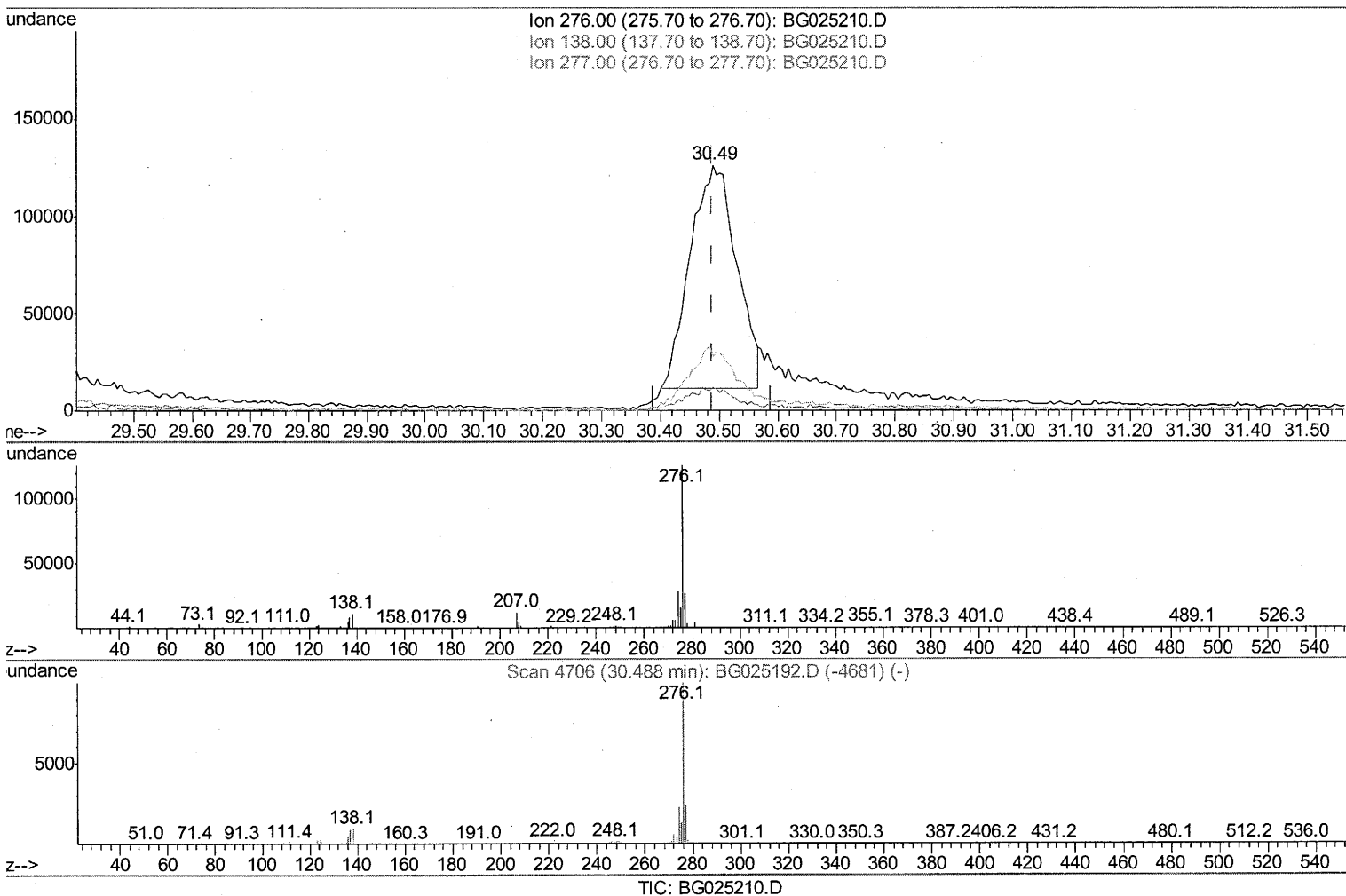
Data Path : Z:\HPCHEM1\BNA G\Data\BG121516\  
Data File : BG025210.D  
Acq On : 15 Dec 2016 14:03  
Operator : UM/SJ  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 2 Sample Multiplier: 1

Instrument :  
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Manual Integrations  
APPROVED

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Response via : Initial Calibration



(91) Benzo(g,h,i)perylene

30.488min (+0.000) 11.81ng/ul

response 626810

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 276.00 | 100   | 100   |
| 138.00 | 10.70 | 9.35  |
| 277.00 | 22.00 | 21.69 |
| 0.00   | 0.00  | 0.00  |

# Quantitation Report (Qedit)

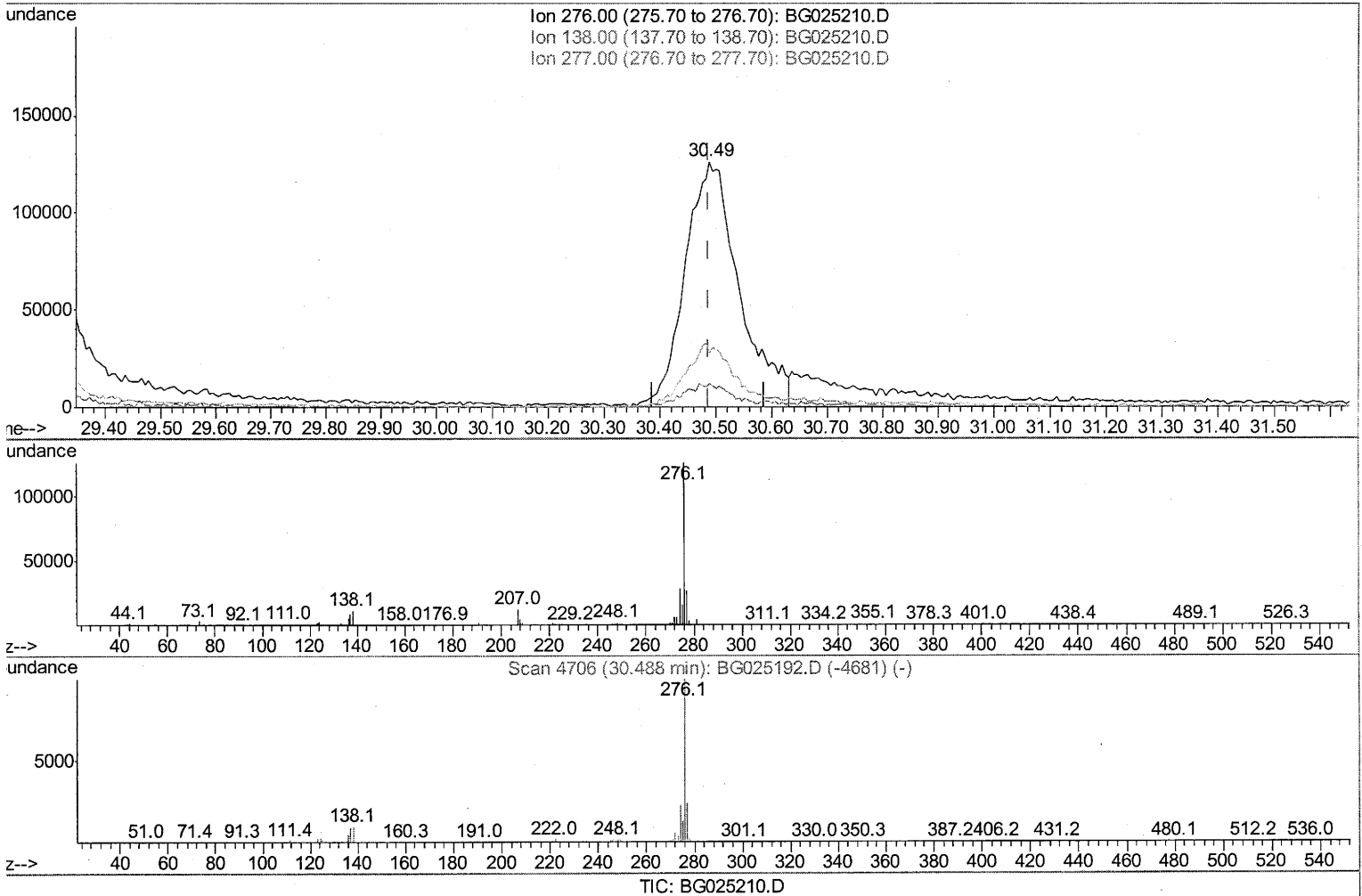
Data Path : Z:\HPCHEM1\BNA G\Data\BG121516\  
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(91) Benzo(g,h,i)perylene

30.488min (+0.000) 15.90ng/ul m

response 843885

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 276.00 | 100   | 100   |
| 138.00 | 10.70 | 9.35  |
| 277.00 | 22.00 | 21.69 |
| 0.00   | 0.00  | 0.00  |

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 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.23  | 152  | 110453   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 11.06 | 136  | 485392   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.86 | 164  | 397162   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.60 | 188  | 867886   | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.90 | 240  | 1085003  | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 25.32 | 264  | 1073081  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.58  | 96  | 14982  | 7.01  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.39  | 99  | 188357 | 19.77 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.55  | 67  | 114108 | 19.36 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.76  | 132 | 129212 | 19.39 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.94  | 113 | 161774 | 20.32 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.41  | 128 | 67633  | 19.68 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 10.14 | 143 | 85558  | 19.97 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.68 | 165 | 168139 | 20.40 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 11.20 | 131 | 192202 | 23.73 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 14.25 | 166 | 543117 | 19.88 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.55 | 160 | 632324 | 21.13 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 15.07 | 143 | 88302  | 18.81 | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.85 | 176 | 536340 | 20.86 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.97 | 200 | 115030 | 21.44 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.70 | 188 | 740944 | 20.23 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.97 | 212 | 898221 | 20.10 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 25.08 | 264 | 904510 | 20.80 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |              | Ovalue |
|--------------------------------|-------|-----|--------|--------------|--------|
| 2) 1,4-Dioxane                 | 3.63  | 88  | 19138  | 6.87 ng/uL#  | 88     |
| 4) Benzaldehyde                | 7.37  | 77  | 146476 | 20.42 ng/ul  | 91     |
| 6) Phenol                      | 7.42  | 94  | 200001 | 20.45 ng/ul  | 95     |
| 8) Bis(2-Chloroethyl)ether     | 7.64  | 93  | 138375 | 19.33 ng/ul  | 93     |
| 10) 2-Chlorophenol             | 7.80  | 128 | 136190 | 20.33 ng/ul  | 93     |
| 11) 2-Methylphenol             | 8.68  | 108 | 138497 | 19.23 ng/ul  | 89     |
| 12) 2,2'-oxybis(1-Chloropropan | 8.76  | 45  | 237978 | 19.55 ng/ul  | 99     |
| 14) Acetophenone               | 9.07  | 105 | 234107 | 19.32 ng/ul  | 96     |
| 15) N-Nitroso-di-n-propylamine | 9.04  | 70  | 132338 | 18.91 ng/ul# | 96     |
| 16) 4-Methylphenol             | 9.00  | 108 | 159361 | 19.89 ng/ul  | 90     |
| 17) Hexachloroethane           | 9.32  | 117 | 57960  | 19.53 ng/ul  | 95     |
| 20) Nitrobenzene               | 9.45  | 77  | 214900 | 20.35 ng/ul  | 96     |
| 21) Isophorone                 | 9.97  | 82  | 377781 | 18.90 ng/ul  | 97     |
| 23) 2-Nitrophenol              | 10.17 | 139 | 90545  | 20.83 ng/ul  | 87     |
| 24) 2,4-Dimethylphenol         | 10.22 | 107 | 198372 | 19.96 ng/ul  | 96     |
| 25) Bis(2-Chloroethoxy)methane | 10.45 | 93  | 205212 | 20.15 ng/ul# | 91     |
| 27) 2,4-Dichlorophenol         | 10.71 | 162 | 166868 | 20.85 ng/ul  | 95     |
| 28) Naphthalene                | 11.11 | 128 | 435965 | 19.32 ng/ul  | 98     |
| 30) 4-Chloroaniline            | 11.22 | 127 | 194121 | 23.45 ng/ul  | 95     |
| 31) Hexachlorobutadiene        | 11.38 | 225 | 133871 | 19.46 ng/ul  | 97     |
| 32) Caprolactam                | 11.99 | 113 | 57804m | 17.40 ng/ul  | 94     |
| 33) 4-Chloro-3-methylphenol    | 12.33 | 107 | 198963 | 20.17 ng/ul  | 94     |
| 34) 2-Methylnaphthalene        | 12.70 | 142 | 347883 | 19.54 ng/ul  | 96     |

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| Internal Standards             | R.T.  | QIon | Response | Conc  | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|-------|--------|----------|
| 36) 1,2,4,5-Tetrachlorobenzene | 13.06 | 216  | 247875   | 20.71 | nq/ul  | 98       |
| 37) Hexachlorocyclopentadiene  | 13.03 | 237  | 126724   | 18.03 | nq/ul# | 97       |
| 38) 2,4,6-Trichlorophenol      | 13.30 | 196  | 166398   | 22.13 | nq/ul  | 95       |
| 39) 2,4,5-Trichlorophenol      | 13.38 | 196  | 174695   | 21.20 | nq/ul  | 93       |
| 40) 1,1'-Biphenyl              | 13.69 | 154  | 494311   | 20.53 | nq/ul  | 99       |
| 41) 2-Chloronaphthalene        | 13.74 | 162  | 390933   | 20.39 | nq/ul  | 97       |
| 42) 2-Nitroaniline             | 13.95 | 65   | 164708   | 21.95 | nq/ul  | 95       |
| 44) Dimethylphthalate          | 14.30 | 163  | 526917   | 19.63 | nq/ul# | 98       |
| 45) 2,6-Dinitrotoluene         | 14.44 | 165  | 113819   | 19.73 | nq/ul  | 92       |
| 47) Acenaphthylene             | 14.58 | 152  | 605258   | 20.81 | nq/ul  | 99       |
| 48) 3-Nitroaniline             | 14.77 | 138  | 107839   | 21.58 | nq/ul# | 80       |
| 49) Acenaphthene               | 14.92 | 153  | 411680   | 20.66 | nq/ul  | 98       |
| 50) 2,4-Dinitrophenol          | 14.98 | 184  | 71428    | 19.06 | nq/ul  | 96       |
| 52) 4-Nitrophenol              | 15.08 | 109  | 107097   | 18.96 | nq/ul  | 88       |
| 53) Dibenzofuran               | 15.25 | 168  | 644533   | 20.45 | nq/ul  | 99       |
| 54) 2,4-Dinitrotoluene         | 15.22 | 165  | 166368   | 19.14 | nq/ul# | 99       |
| 55) 2,3,4,6-Tetrachlorophenol  | 15.48 | 232  | 175388   | 20.14 | nq/ul# | 99       |
| 56) Diethylphthalate           | 15.65 | 149  | 539357   | 18.97 | nq/ul  | 97       |
| 58) Fluorene                   | 15.90 | 166  | 534820   | 20.84 | nq/ul  | 98       |
| 59) 4-Chlorophenyl-phenylether | 15.88 | 204  | 294079   | 20.35 | nq/ul  | 89       |
| 60) 4-Nitroaniline             | 15.93 | 138  | 114161   | 18.95 | nq/ul  | 99       |
| 63) 4,6-Dinitro-2-methylphenol | 15.99 | 198  | 115714   | 20.82 | nq/ul# | 97       |
| 64) N-Nitrosodiphenylamine     | 16.10 | 169  | 477197   | 21.37 | nq/ul  | 96       |
| 65) 4-Bromophenyl-phenylether  | 16.78 | 248  | 211596   | 21.58 | nq/ul  | 93       |
| 66) Hexachlorobenzene          | 16.90 | 284  | 239219   | 21.67 | nq/ul  | 97       |
| 67) Atrazine                   | 17.03 | 200  | 210866   | 19.87 | nq/ul  | 97       |
| 68) Pentachlorophenol          | 17.25 | 266  | 124178   | 18.74 | nq/ul# | 89       |
| 69) Phenanthrene               | 17.64 | 178  | 852240   | 20.57 | nq/ul  | 99       |
| 71) Anthracene                 | 17.73 | 178  | 861806   | 20.24 | nq/ul  | 99       |
| 72) Carbazole                  | 18.00 | 167  | 748663   | 21.07 | nq/ul  | 97       |
| 73) Di-n-butylphthalate        | 18.53 | 149  | 900667   | 20.08 | nq/ul  | 99       |
| 74) Fluoranthene               | 19.64 | 202  | 1066525  | 21.67 | nq/ul  | 99       |
| 77) Pvrene                     | 20.00 | 202  | 1072153  | 20.56 | nq/ul  | 97       |
| 78) Butylbenzylphthalate       | 20.85 | 149  | 407724   | 19.97 | nq/ul  | 96       |
| 79) 3,3'-Dichlorobenzidine     | 21.78 | 252  | 427254   | 22.59 | nq/ul  | 95       |
| 80) Benzo(a)anthracene         | 21.87 | 228  | 1143948  | 20.29 | nq/ul  | 97       |
| 81) Bis(2-ethylhexyl)phthalate | 21.72 | 149  | 545033   | 19.39 | nq/ul  | 99       |
| 82) Chrysene                   | 21.95 | 228  | 1074466  | 20.38 | nq/ul  | 99       |
| 84) Di-n-octyl phthalate       | 22.99 | 149  | 979765   | 22.49 | nq/ul  | 100      |
| 85) Benzo(b)fluoranthene       | 24.21 | 252  | 1084301  | 20.38 | nq/ul  | 100      |
| 86) Benzo(k)fluoranthene       | 24.29 | 252  | 1097754  | 21.70 | nq/ul  | 97       |
| 88) Benzo(a)pyrene             | 25.15 | 252  | 1067169  | 20.86 | nq/ul# | 99       |
| 89) Indeno(1,2,3-cd)pyrene     | 29.25 | 276  | 1194692m | 18.79 | nq/ul  |          |
| 90) Dibenzo(a,h)anthracene     | 29.31 | 278  | 996899m  | 18.61 | nq/ul  |          |
| 91) Benzo(q,h,i)perylene       | 30.49 | 276  | 843885m  | 15.90 | nq/ul  |          |

UM  
 12/16/16

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