

Method Path : Z:\voasrv\HPCHEM1\MSVOA\_D\Method\  
 Method File : 82D072224S.M  
 Title : SW846 8260  
 Last Update : Tue Jul 23 10:24:07 2024  
 Response Via : Initial Calibration

## Calibration Files

5 =VD079393.D 10 =VD079387.D 20 =VD079388.D 50 =VD079389.D 100 =VD079390.D 150 =VD079391.D

| Compound | 5                                  | 10    | 20    | 50    | 100   | 150   | Avg   | %RSD  |        |
|----------|------------------------------------|-------|-------|-------|-------|-------|-------|-------|--------|
| -----    |                                    |       |       |       |       |       |       |       |        |
| 1) I     | Pentafluorobenzene -----ISTD-----  |       |       |       |       |       |       |       |        |
| 2) T     | Dichlorodifluo...                  | 0.569 | 0.484 | 0.486 | 0.399 | 0.407 | 0.374 | 0.453 | 16.10  |
| 3) P     | Chloromethane                      | 0.476 | 0.489 | 0.480 | 0.472 | 0.447 | 0.383 | 0.458 | 8.52   |
| 4) C     | Vinyl Chloride                     | 0.462 | 0.596 | 0.575 | 0.598 | 0.470 | 0.399 | 0.517 | 16.25# |
| 5) T     | Bromomethane                       | 0.512 | 0.463 | 0.357 | 0.398 | 0.349 | 0.308 | 0.398 | 19.22  |
| 6) T     | Chloroethane                       | 0.264 | 0.295 | 0.229 | 0.281 | 0.253 | 0.199 | 0.254 | 13.85  |
| 7) T     | Trichlorofluor...                  | 1.031 | 1.018 | 0.848 | 0.851 | 0.850 | 0.763 | 0.893 | 11.96  |
| 8) T     | Diethyl Ether                      | 0.303 | 0.326 | 0.255 | 0.302 | 0.290 | 0.259 | 0.289 | 9.55   |
| 9) T     | 1,1,2-Trichlor...                  | 0.722 | 0.676 | 0.590 | 0.594 | 0.590 | 0.516 | 0.615 | 11.85  |
| 10) T    | Methyl Iodide                      | 0.577 | 0.638 | 0.665 | 0.633 | 0.768 | 0.724 | 0.667 | 10.27  |
| 11) T    | Tert butyl alc...                  | 0.040 | 0.038 | 0.034 | 0.036 | 0.034 | 0.032 | 0.035 | 8.47   |
| 12) CM   | 1,1-Dichloroet...                  | 0.664 | 0.628 | 0.549 | 0.555 | 0.574 | 0.526 | 0.583 | 9.06#  |
| 13) T    | Acrolein                           | 0.037 | 0.039 | 0.032 | 0.023 | 0.020 | 0.018 | 0.028 | 31.79  |
| 14) T    | Allyl chloride                     | 0.588 | 0.796 | 0.547 | 0.745 | 0.776 | 0.691 | 0.691 | 16.56  |
| 15) T    | Acrylonitrile                      | 0.149 | 0.150 | 0.122 | 0.144 | 0.142 | 0.133 | 0.140 | 7.55   |
| 16) T    | Acetone                            | 0.093 | 0.126 | 0.087 | 0.122 | 0.098 | 0.075 | 0.100 | 19.89  |
| 17) T    | Carbon Disulfide                   | 1.830 | 1.977 | 1.636 | 1.669 | 1.675 | 1.396 | 1.697 | 11.54  |
| 18) T    | Methyl Acetate                     | 0.327 | 0.373 | 0.265 | 0.343 | 0.345 | 0.331 | 0.331 | 12.15  |
| 19) T    | Methyl tert-bu...                  | 1.336 | 1.323 | 1.200 | 1.365 | 1.375 | 1.311 | 1.318 | 4.78   |
| 20) T    | Methylene Chlo...                  | 2.820 | 1.913 | 1.169 | 0.933 | 0.764 | 0.720 | 1.386 | 59.63  |
| 21) T    | trans-1,2-Dich...                  | 0.736 | 0.667 | 0.654 | 0.631 | 0.631 | 0.602 | 0.653 | 7.05   |
| 22) T    | Diisopropyl ether                  | 1.684 | 1.835 | 1.768 | 1.763 | 1.709 | 1.598 | 1.726 | 4.75   |
| 23) T    | Vinyl Acetate                      | 1.298 | 1.323 | 1.364 | 1.394 | 1.377 | 1.267 | 1.337 | 3.69   |
| 24) P    | 1,1-Dichloroet...                  | 1.437 | 1.238 | 1.161 | 1.120 | 1.123 | 1.055 | 1.189 | 11.39  |
| 25) T    | 2-Butanone                         | 0.190 | 0.172 | 0.174 | 0.169 | 0.165 | 0.154 | 0.171 | 6.92   |
| 26) T    | 2,2-Dichloropr...                  | 1.192 | 1.031 | 0.987 | 0.922 | 0.939 | 0.876 | 0.991 | 11.31  |
| 27) T    | cis-1,2-Dichlo...                  | 0.789 | 0.766 | 0.749 | 0.733 | 0.761 | 0.730 | 0.755 | 2.93   |
| 28) T    | Bromochloromet...                  | 0.574 | 0.488 | 0.499 | 0.469 | 0.449 | 0.419 | 0.483 | 10.99  |
| 29) T    | Tetrahydrofuran                    | 0.105 | 0.099 | 0.107 | 0.107 | 0.105 | 0.097 | 0.103 | 4.05   |
| 30) C    | Chloroform                         | 1.431 | 1.358 | 1.213 | 1.175 | 1.168 | 1.091 | 1.239 | 10.39# |
| 31) T    | Cyclohexane                        | 1.121 | 0.968 | 0.921 | 0.831 | 0.835 | 0.800 | 0.913 | 13.14  |
| 32) T    | 1,1,1-Trichlor...                  | 1.197 | 1.060 | 1.009 | 0.957 | 0.973 | 0.905 | 1.017 | 10.06  |
| 33) S    | 1,2-Dichloroet...                  | 0.606 | 0.560 | 0.600 | 0.576 | 0.553 | 0.514 | 0.568 | 5.96   |
| -----    |                                    |       |       |       |       |       |       |       |        |
| 34) I    | 1,4-Difluorobenzene -----ISTD----- |       |       |       |       |       |       |       |        |
| 35) S    | Dibromofluorom...                  | 0.357 | 0.351 | 0.361 | 0.356 | 0.343 | 0.320 | 0.348 | 4.30   |
| 36) T    | 1,1-Dichloropr...                  | 0.501 | 0.485 | 0.463 | 0.458 | 0.464 | 0.432 | 0.467 | 5.13   |
| 37) T    | Ethyl Acetate                      | 0.212 | 0.194 | 0.199 | 0.218 | 0.206 | 0.187 | 0.203 | 5.64   |
| 38) T    | Carbon Tetrach...                  | 0.545 | 0.515 | 0.505 | 0.473 | 0.476 | 0.444 | 0.493 | 7.28   |
| 39) T    | Methylcyclohexane                  | 0.510 | 0.509 | 0.541 | 0.551 | 0.571 | 0.539 | 0.537 | 4.49   |
| 40) TM   | Benzene                            | 1.587 | 1.496 | 1.458 | 1.433 | 1.474 | 1.371 | 1.470 | 4.87   |
| 41) T    | Methacrylonitrile                  | 0.095 | 0.124 | 0.110 | 0.107 | 0.117 | 0.110 | 0.111 | 8.84   |
| 42) TM   | 1,2-Dichloroet...                  | 0.404 | 0.396 | 0.373 | 0.349 | 0.350 | 0.326 | 0.366 | 8.18   |
| 43) T    | Isopropyl Acetate                  | 0.344 | 0.379 | 0.384 | 0.389 | 0.382 | 0.357 | 0.372 | 4.80   |
| 44) TM   | Trichloroethene                    | 0.383 | 0.378 | 0.362 | 0.338 | 0.344 | 0.325 | 0.355 | 6.51   |
| 45) C    | 1,2-Dichloropr...                  | 0.380 | 0.401 | 0.364 | 0.356 | 0.359 | 0.338 | 0.366 | 5.88#  |
| 46) T    | Dibromomethane                     | 0.235 | 0.234 | 0.214 | 0.208 | 0.208 | 0.189 | 0.214 | 8.11   |
| 47) T    | Bromodichlorom...                  | 0.568 | 0.570 | 0.529 | 0.515 | 0.515 | 0.475 | 0.528 | 6.85   |
| 48) T    | Methyl methacr...                  | 0.161 | 0.157 | 0.172 | 0.174 | 0.180 | 0.166 | 0.168 | 4.99   |
| 49) T    | 1,4-Dioxane                        | 0.002 | 0.002 | 0.002 | 0.002 | 0.003 | 0.002 | 0.002 | 4.92   |
| 50) S    | Toluene-d8                         | 1.226 | 1.256 | 1.371 | 1.375 | 1.355 | 1.310 | 1.315 | 4.76   |
| 51) T    | 4-Methyl-2-Pen...                  | 0.186 | 0.197 | 0.212 | 0.215 | 0.208 | 0.190 | 0.201 | 6.03   |
| 52) CM   | Toluene                            | 0.933 | 0.915 | 0.928 | 0.909 | 0.946 | 0.888 | 0.920 | 2.21#  |
| 53) T    | t-1,3-Dichloro...                  | 0.494 | 0.485 | 0.476 | 0.470 | 0.487 | 0.458 | 0.478 | 2.75   |
| 54) T    | cis-1,3-Dichlo...                  | 0.552 | 0.569 | 0.566 | 0.562 | 0.576 | 0.531 | 0.559 | 2.87   |
| 55) T    | 1,1,2-Trichlor...                  | 0.303 | 0.296 | 0.290 | 0.284 | 0.283 | 0.258 | 0.286 | 5.41   |
| 56) T    | Ethyl methacry...                  | 0.299 | 0.323 | 0.341 | 0.364 | 0.381 | 0.355 | 0.344 | 8.52   |

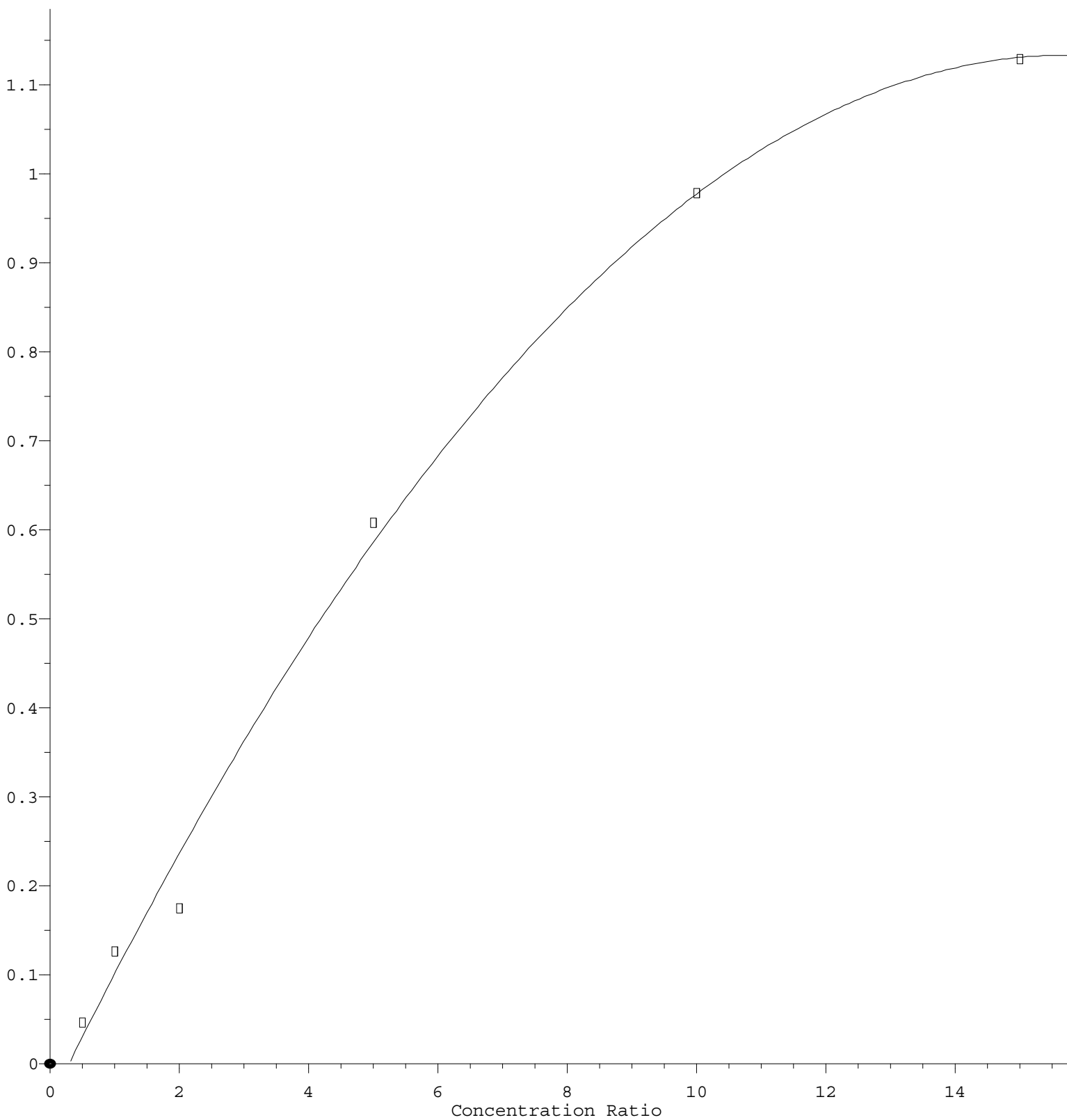
Method Path : Z:\voasrv\HPCHEM1\MSVOA\_D\Method\  
 Method File : 82D072224S.M

|     |    |                       |                |       |       |       |       |       |       |       |
|-----|----|-----------------------|----------------|-------|-------|-------|-------|-------|-------|-------|
| 57) | T  | 1,3-Dichloropr...     | 0.462          | 0.501 | 0.479 | 0.469 | 0.474 | 0.439 | 0.471 | 4.33  |
| 58) | T  | 2-Chloroethyl ...     | 0.130          | 0.146 | 0.150 | 0.167 | 0.168 | 0.158 | 0.153 | 9.23  |
| 59) | T  | 2-Hexanone            | 0.130          | 0.139 | 0.150 | 0.153 | 0.148 | 0.136 | 0.143 | 6.26  |
| 60) | T  | Dibromochlorom...     | 0.387          | 0.379 | 0.375 | 0.366 | 0.359 | 0.336 | 0.367 | 4.90  |
| 61) | T  | 1,2-Dibromoethane     | 0.279          | 0.267 | 0.269 | 0.265 | 0.264 | 0.246 | 0.265 | 4.06  |
| 62) | S  | 4-Bromofluorob...     | 0.372          | 0.379 | 0.415 | 0.420 | 0.417 | 0.412 | 0.402 | 5.26  |
| 63) | I  | Chlorobenzene-d5      | -----ISTD----- |       |       |       |       |       |       |       |
| 64) | T  | Tetrachloroethene     | 0.378          | 0.326 | 0.305 | 0.293 | 0.303 | 0.289 | 0.316 | 10.53 |
| 65) | PM | Chlorobenzene         | 1.220          | 1.102 | 1.097 | 1.064 | 1.068 | 1.039 | 1.098 | 5.84  |
| 66) | T  | 1,1,1,2-Tetrac...     | 0.400          | 0.375 | 0.368 | 0.367 | 0.366 | 0.361 | 0.373 | 3.75  |
| 67) | C  | Ethyl Benzene         | 1.744          | 1.703 | 1.780 | 1.845 | 1.932 | 1.917 | 1.820 | 5.13# |
| 68) | T  | m/p-Xylenes           | 0.649          | 0.669 | 0.707 | 0.723 | 0.743 | 0.745 | 0.706 | 5.59  |
| 69) | T  | o-Xylene              | 0.619          | 0.608 | 0.636 | 0.679 | 0.695 | 0.701 | 0.656 | 6.16  |
| 70) | T  | Styrene               | 1.036          | 1.055 | 1.168 | 1.229 | 1.258 | 1.243 | 1.165 | 8.37  |
| 71) | P  | Bromoform             | 0.235          | 0.222 | 0.223 | 0.216 | 0.211 | 0.210 | 0.219 | 4.14  |
| 72) | I  | 1,4-Dichlorobenzen... | -----ISTD----- |       |       |       |       |       |       |       |
| 73) | T  | Isopropylbenzene      | 3.262          | 3.368 | 3.497 | 3.567 | 3.657 | 3.569 | 3.487 | 4.19  |
| 74) | T  | N-amyl acetate        | 0.737          | 0.769 | 0.838 | 0.834 | 0.814 | 0.764 | 0.793 | 5.29  |
| 75) | P  | 1,1,2,2-Tetrac...     | 0.842          | 0.808 | 0.771 | 0.762 | 0.735 | 0.673 | 0.765 | 7.68  |
| 76) | T  | 1,2,3-Trichlor...     | 0.637          | 0.438 | 0.563 | 0.442 | 0.517 | 0.486 | 0.514 | 14.94 |
| 77) | T  | Bromobenzene          | 0.883          | 0.851 | 0.857 | 0.841 | 0.842 | 0.807 | 0.847 | 2.93  |
| 78) | T  | n-propylbenzene       | 4.281          | 4.160 | 4.421 | 4.514 | 4.533 | 4.402 | 4.385 | 3.25  |
| 79) | T  | 2-Chlorotoluene       | 2.509          | 2.449 | 2.528 | 2.505 | 2.538 | 2.480 | 2.501 | 1.31  |
| 80) | T  | 1,3,5-Trimethy...     | 2.720          | 2.810 | 2.950 | 2.996 | 3.080 | 2.989 | 2.924 | 4.57  |
| 81) | T  | trans-1,4-Dich...     | 0.256          | 0.270 | 0.288 | 0.270 | 0.261 | 0.247 | 0.265 | 5.36  |
| 82) | T  | 4-Chlorotoluene       | 2.753          | 2.642 | 2.775 | 2.621 | 2.677 | 2.553 | 2.670 | 3.13  |
| 83) | T  | tert-Butylbenzene     | 2.340          | 2.324 | 2.463 | 2.528 | 2.666 | 2.609 | 2.488 | 5.60  |
| 84) | T  | 1,2,4-Trimethy...     | 2.675          | 2.780 | 2.990 | 3.070 | 3.185 | 3.105 | 2.968 | 6.70  |
| 85) | T  | sec-Butylbenzene      | 3.824          | 3.707 | 3.872 | 3.869 | 4.076 | 3.944 | 3.882 | 3.17  |
| 86) | T  | p-Isopropyltol...     | 2.905          | 2.916 | 3.174 | 3.250 | 3.474 | 3.369 | 3.181 | 7.34  |
| 87) | T  | 1,3-Dichlorobe...     | 1.890          | 1.803 | 1.762 | 1.717 | 1.746 | 1.679 | 1.766 | 4.16  |
| 88) | T  | 1,4-Dichlorobe...     | 1.947          | 1.854 | 1.772 | 1.689 | 1.708 | 1.621 | 1.765 | 6.73  |
| 89) | T  | n-Butylbenzene        | 3.034          | 2.882 | 3.096 | 3.185 | 3.271 | 3.188 | 3.109 | 4.44  |
| 90) | T  | Hexachloroethane      | 0.807          | 0.744 | 0.713 | 0.672 | 0.683 | 0.664 | 0.714 | 7.60  |
| 91) | T  | 1,2-Dichlorobe...     | 1.591          | 1.558 | 1.555 | 1.488 | 1.519 | 1.459 | 1.528 | 3.20  |
| 92) | T  | 1,2-Dibromo-3-...     | 0.116          | 0.107 | 0.111 | 0.106 | 0.102 | 0.094 | 0.106 | 7.00  |
| 93) | T  | 1,2,4-Trichlor...     | 0.881          | 0.862 | 0.850 | 0.884 | 0.904 | 0.888 | 0.878 | 2.20  |
| 94) | T  | Hexachlorobuta...     | 0.561          | 0.489 | 0.494 | 0.471 | 0.471 | 0.458 | 0.491 | 7.53  |
| 95) | T  | Naphthalene           | 1.700          | 1.589 | 1.716 | 1.875 | 1.940 |       | 1.764 | 8.04  |
| 96) | T  | 1,2,3-Trichlor...     | 0.874          | 0.740 | 0.775 | 0.803 | 0.816 | 0.775 | 0.797 | 5.77  |

(#) = Out of Range

Acetone

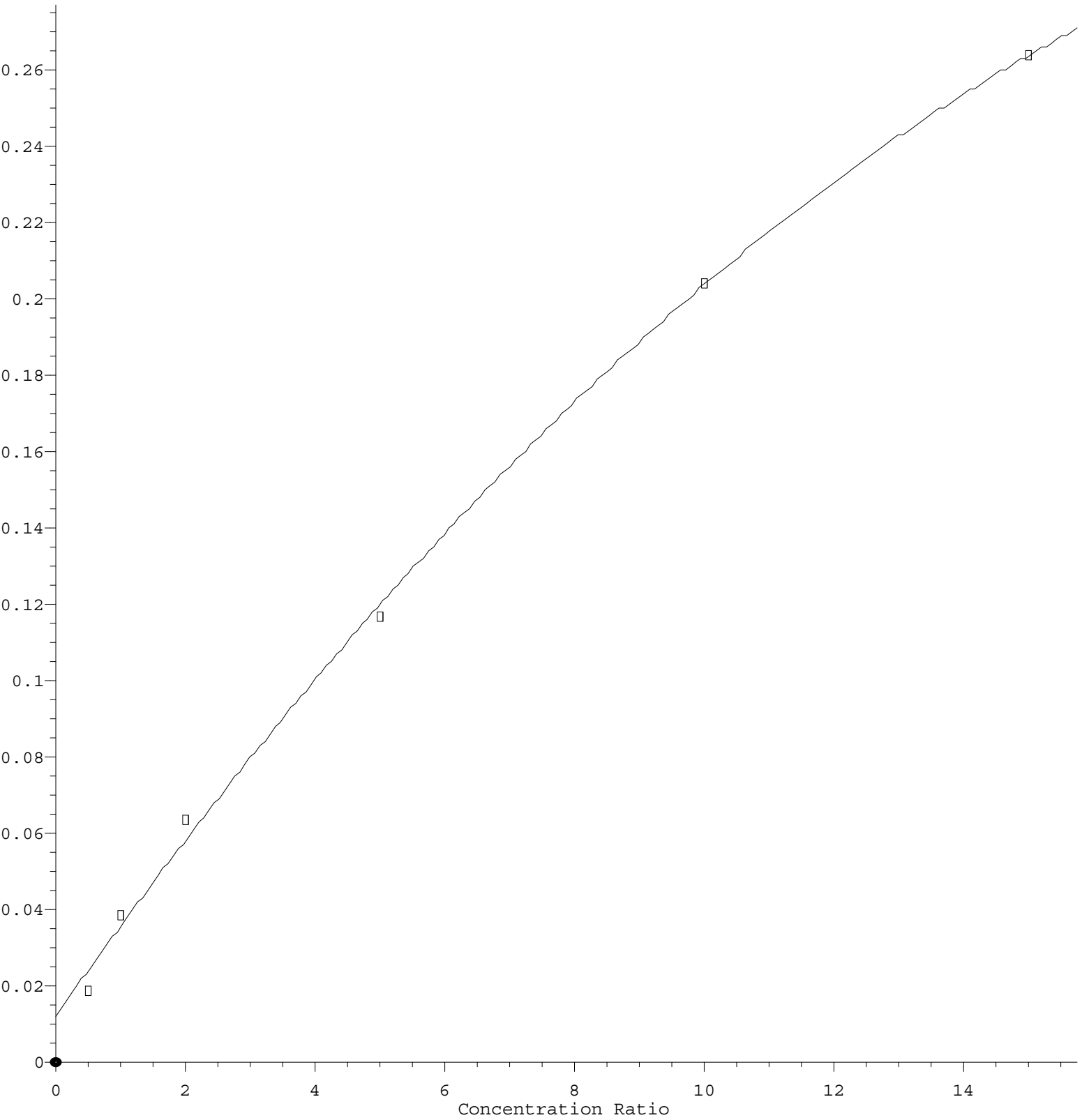
Response Ratio



R = -4.758e-003 A\*A + 1.497e-001 A - 4.355e-002  
Coef of Det (r^2) = 0.995182 Curve Fit: Quadratic  
Method Name: Z:\voasrv\HPCHEM1\MSVOA D\Method\82D072224S.M  
Calibration Table Last Updated: Tue Jul 23 10:24:07 2024

Acrolein

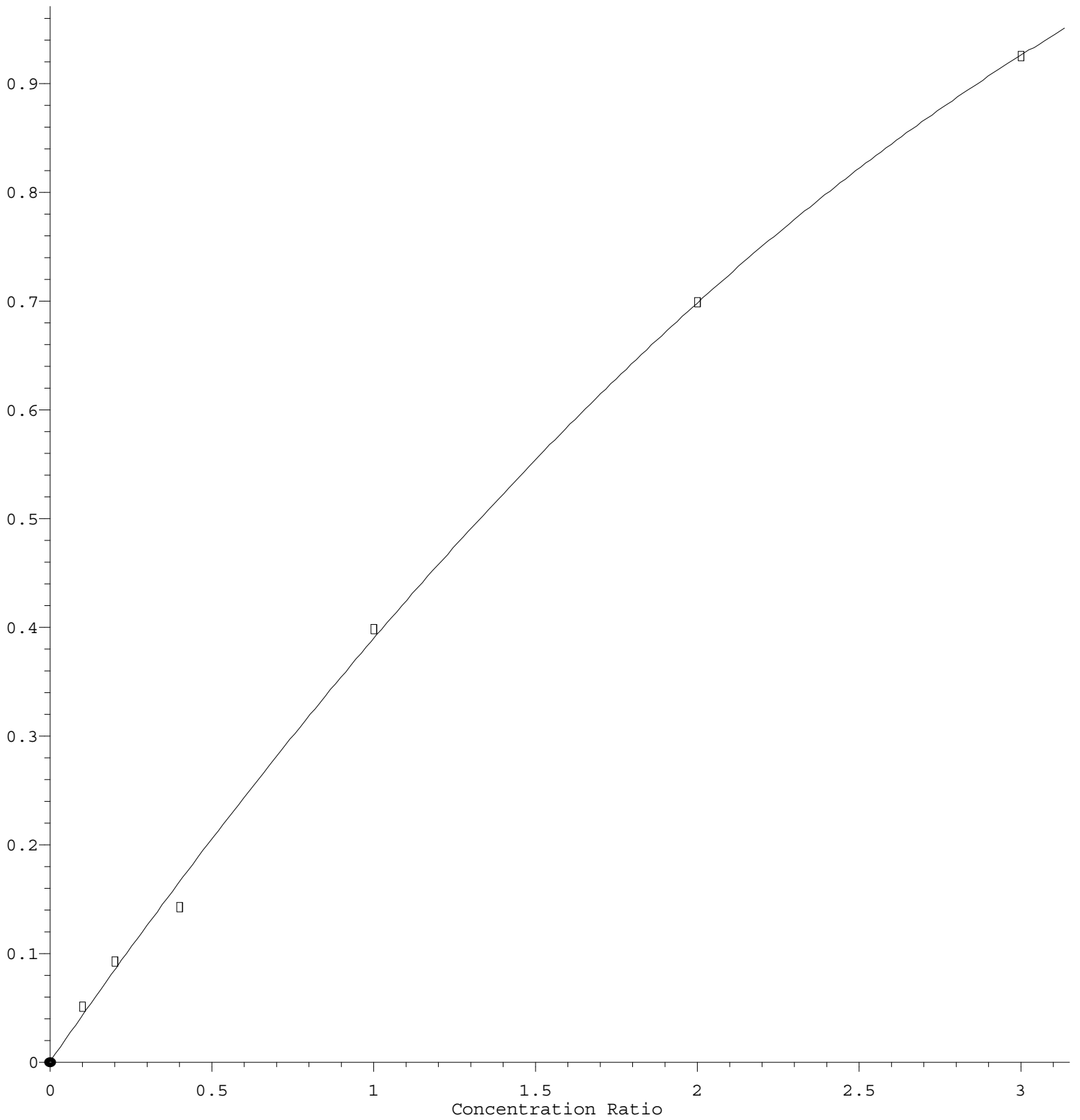
Response Ratio



R = -4.746e-004 A\*A + 2.388e-002 A + 1.232e-002  
Coef of Det (r^2) = 0.998430 Curve Fit: Quadratic  
Method Name: Z:\voasrv\HPCHEM1\MSVOA D\Method\82D072224S.M  
Calibration Table Last Updated: Tue Jul 23 10:24:07 2024

# Bromomethane

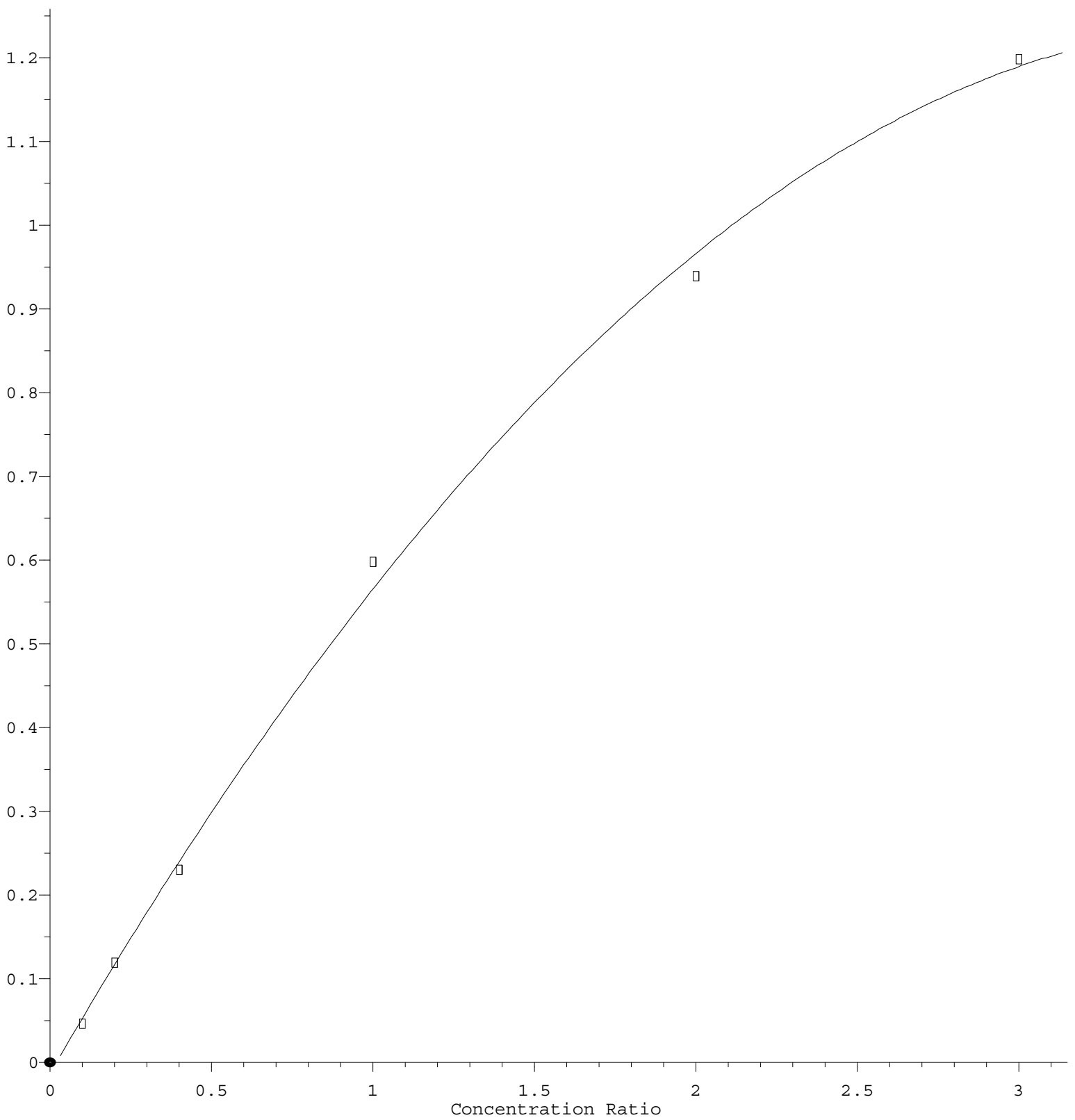
Response Ratio



R =  $-4.031e-002 A^2 + 4.293e-001 A + 8.705e-004$   
Coef of Det ( $r^2$ ) = 0.998856 Curve Fit: Quadratic  
Method Name: Z:\voasrv\HPCHEM1\MSVOA D\Method\82D072224S.M  
Calibration Table Last Updated: Tue Jul 23 10:24:07 2024

# Vinyl Chloride

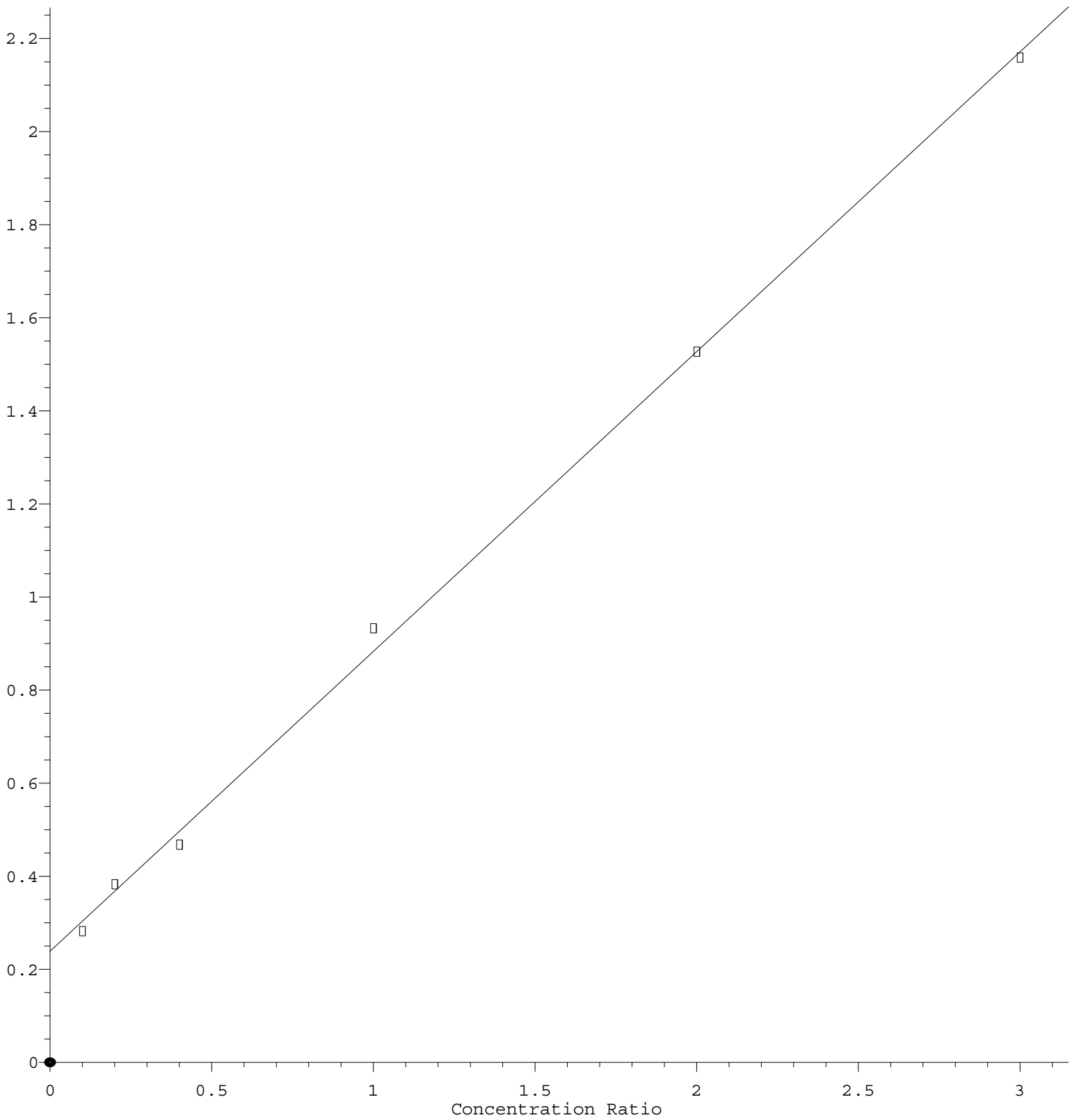
Response Ratio



$R = -8.879e-002 A^2 + 6.672e-001 A - 1.299e-002$   
Coef of Det ( $r^2$ ) = 0.998192 Curve Fit: Quadratic  
Method Name: Z:\voasrv\HPCHEM1\MSVOA D\Method\82D072224S.M  
Calibration Table Last Updated: Tue Jul 23 10:24:07 2024

Methylene Chloride

Response Ratio



Response = 6.441e-001 \* Amt + 2.393e-001

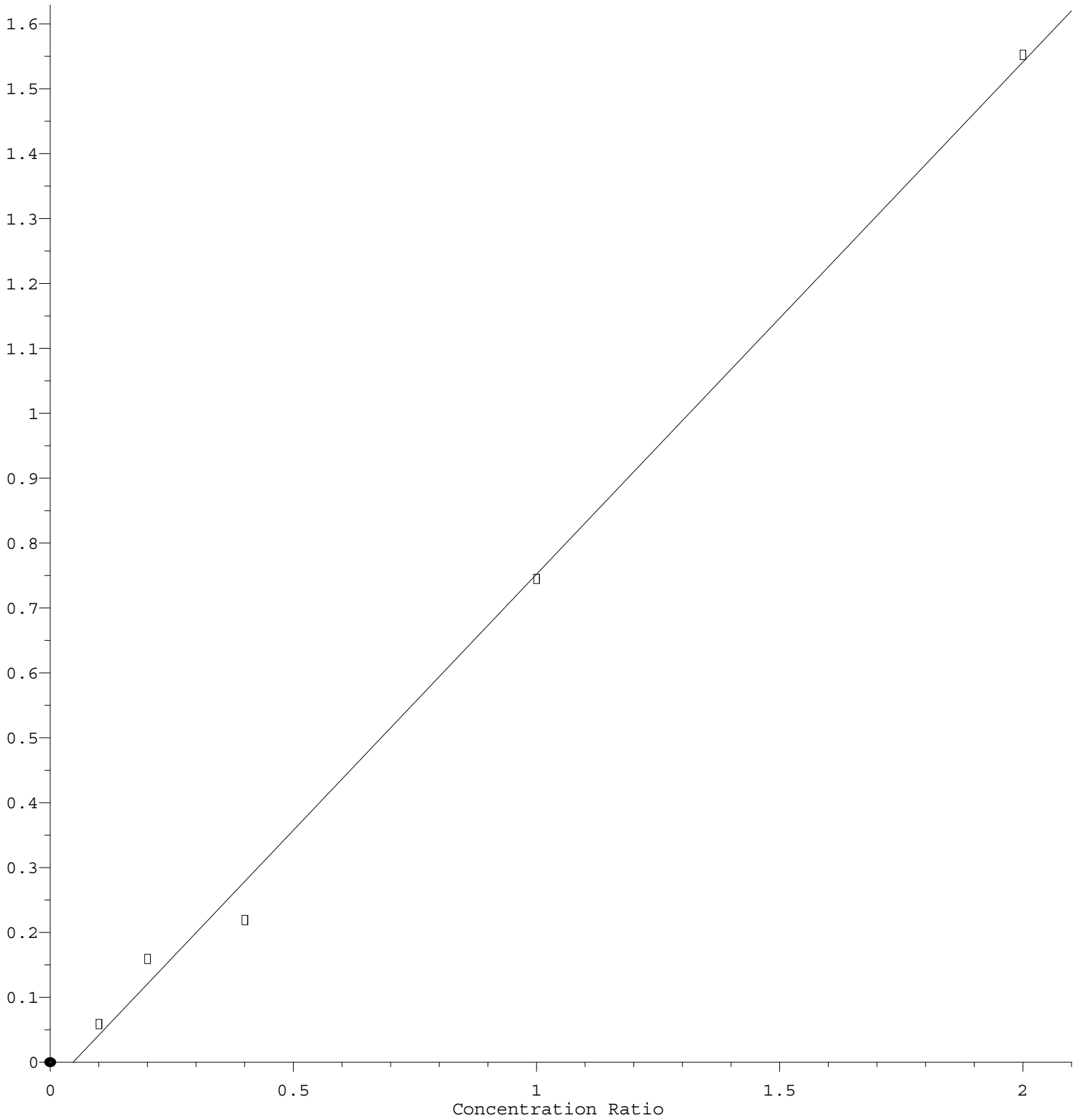
Coef of Det (r^2) = 0.998527 Curve Fit: Linear

Method Name: Z:\voasrv\HPCHEM1\MSVOA D\Method\82D072224S.M

Calibration Table Last Updated: Tue Jul 23 10:24:07 2024

Allyl chloride

Response Ratio



Response = 7.893e-001 \* Amt - 3.732e-002

Coef of Det (r^2) = 0.996445 Curve Fit: Linear

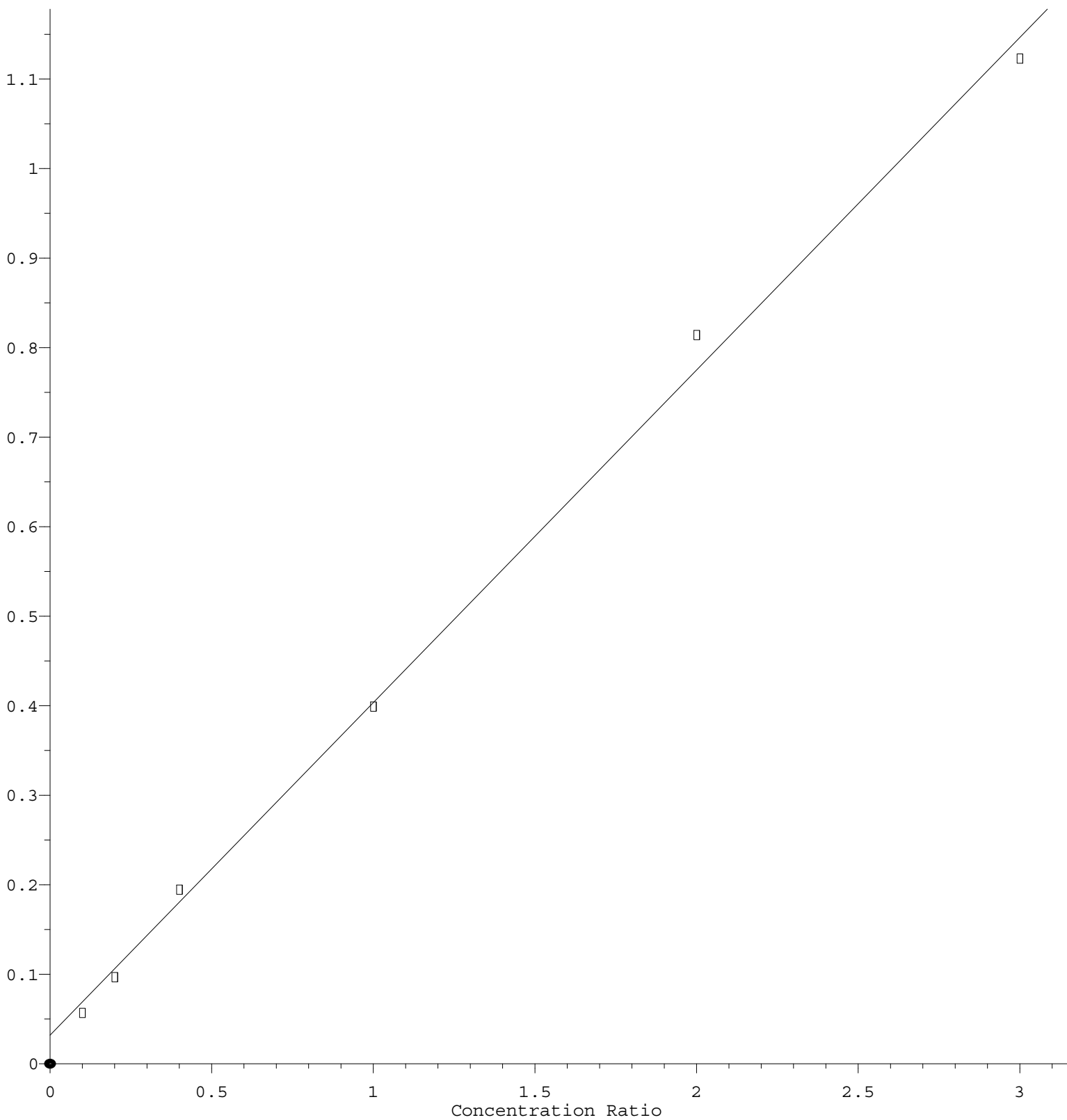
Method Name: Z:\voasrv\HPCHEM1\MSVOA D\Method\82D072224S.M

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Dichlorodifluoromethane

Response Ratio



Response = 3.718e-001 \* Amt + 3.217e-002

Coef of Det (r^2) = 0.997282 Curve Fit: Linear

Method Name: Z:\voasrv\HPCHEM1\MSVOA D\Method\82D072224S.M

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