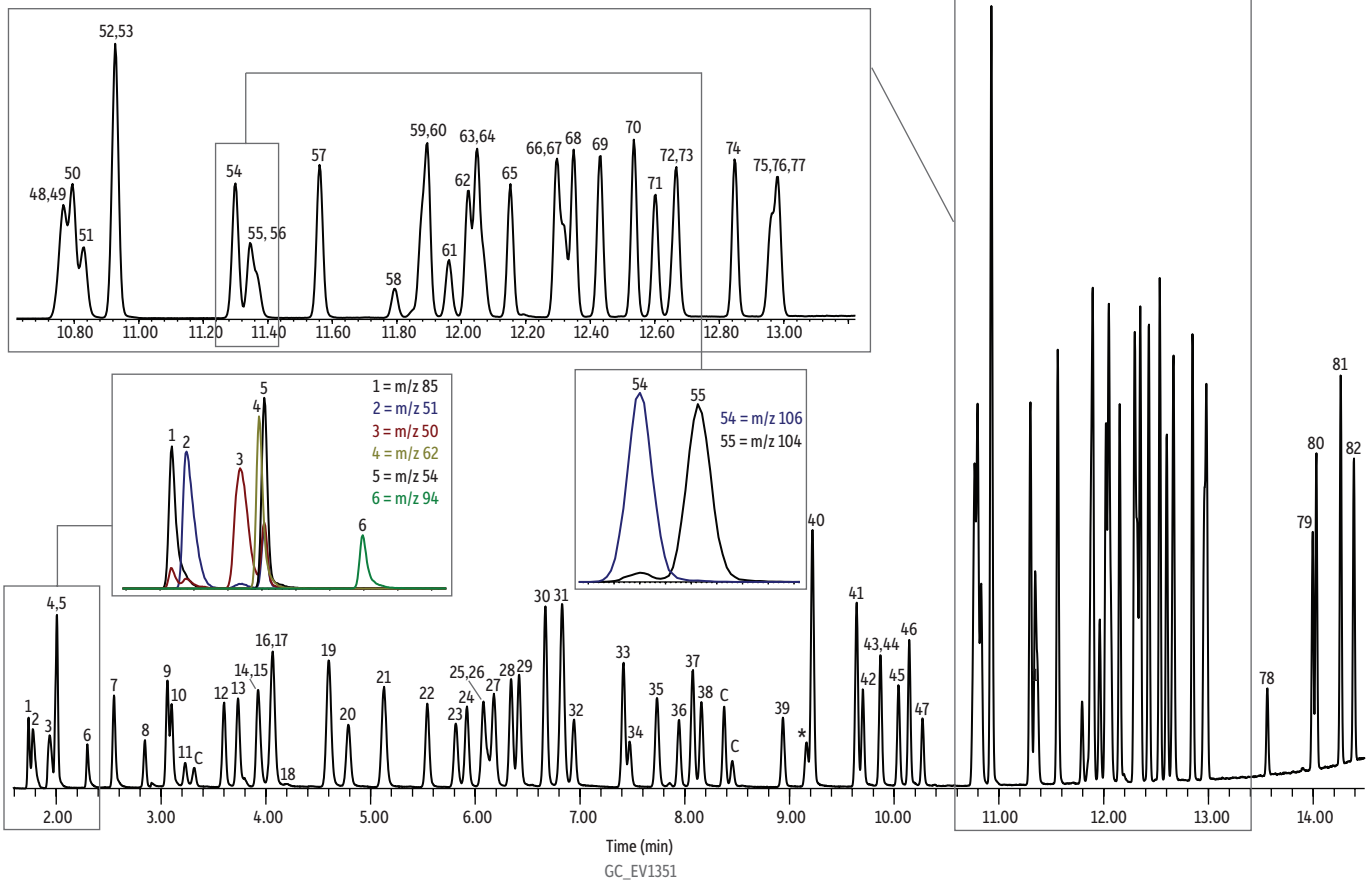


40 ppb Volatiles in Drinking Water on Rtx[®]-VMS by EPA Method 524.4

- | | | | |
|--------------------------------------|---|---------------------------------------|---------------------------------|
| Peaks | Peaks | Peaks | Peaks |
| 1. Dichlorodifluoromethane | 21. <i>tert</i> -Butyl ethyl ether (ETBE) | 41. Tetrachloroethene | 61. 1,1,2,2-Tetrachloroethane |
| 2. Chlorodifluoromethane | 22. <i>cis</i> -1,2-Dichloroethene | 42. <i>trans</i> -1,3-Dichloropropene | 62. 2-Chlorotoluene |
| 3. Chloromethane | 23. Bromochloromethane | 43. 1,1,2-Trichloroethane | 63. 1,3,5-Trimethylbenzene |
| 4. Vinyl chloride | 24. Chloroform | 44. Ethyl methacrylate | 64. 1,2,3-Trichloropropane |
| 5. 1,3-Butadiene | 25. Carbon tetrachloride | 45. Dibromochloromethane | 65. 4-Chlorotoluene |
| 6. Bromomethane | 26. Tetrahydrofuran | 46. 1,3-Dichloropropane | 66. <i>tert</i> -Butylbenzene |
| 7. Trichlorofluoromethane | 27. 1,1,1-Trichloroethane | 47. 1,2-Dibromoethane | 67. Pentachloroethane |
| 8. Diethyl ether | 28. 1,1-Dichloropropene | 48. Chlorobenzene-d5 | 68. 1,2,4-Trimethylbenzene |
| 9. 1,1-Dichloroethene | 29. 1-Chlorobutane | 49. Chlorobenzene | 69. <i>sec</i> -Butylbenzene |
| 10. Carbon disulfide | 30. Benzene | 50. Ethylbenzene | 70. 4-Isopropyltoluene |
| 11. Methyl iodide | 31. <i>tert</i> -Amyl methyl ether (TAME) | 51. 1,1,1,2-Tetrachloroethane | 71. 1,3-Dichlorobenzene |
| 12. Allyl chloride | 32. 1,2-Dichloroethane | 52. <i>m</i> -Xylene | 72. 1,4-Dichlorobenzene-d4 |
| 13. Methylene chloride | 33. Trichloroethene | 53. <i>p</i> -Xylene | 73. 1,4-Dichlorobenzene |
| 14. <i>trans</i> -1,2-Dichloroethene | 34. 1,4-Difluorobenzene | 54. <i>o</i> -Xylene | 74. <i>n</i> -Butylbenzene |
| 15. Methyl acetate | 35. <i>tert</i> -Amyl ethyl ether (TAE) | 55. Styrene | 75. Hexachloroethane |
| 16. MTBE-d3 (SS) | 36. Dibromomethane | 56. Bromoform | 76. 1,2-Dichlorobenzene-d4 (SS) |
| 17. MTBE | 37. 1,2-Dichloropropane | 57. Isopropylbenzene | 77. 1,2-Dichlorobenzene |
| 18. <i>tert</i> -Butyl alcohol (TBA) | 38. Bromodichloromethane | 58. 4-Bromofluorobenzene (SS) | 78. 1,2-Dibromo-3-chloropropane |
| 19. Diisopropyl ether (DIPE) | 39. <i>cis</i> -1,3-Dichloropropene | 59. Bromobenzene | 79. Hexachlorobutadiene |
| 20. 1,1-Dichloroethane | 40. Toluene | 60. <i>n</i> -Propylbenzene | 80. 1,2,4-Trichlorobenzene |

C. Contamination from nitrogen gas line; * Toluene-d8



Column Rtx[®]-VMS, 30 m, 0.25 mm ID, 1.40 μm (cat.# 19915)
Sample 524.3 Internal standard/surrogate mix (cat.# 30017)
 524.3 Gas calibration mix (cat.# 30014)
 524.3 VOA MegaMix[®] standard (cat.# 30013)
Diluent: RO water
Conc.: 40 ng/mL (5 mL sample)
Injection purge and trap split (split ratio 30:1)
Liner: Premium 1.0 mm ID straight inlet liner (cat.# 23333.1)
Inj. Temp.: 200 °C
Purge and Trap
Instrument: EST Encon Evolution
Trap Type: Vocabr 3000
Purge: 11 min, flow 40 mL/min
Dry Purge: 1 min, flow 50 mL/min
Desorb: 1 min @ 260 °C, flow 30.9 mL/min
Bake: 8 min @ 265 °C
Interface
Connection: injection port
Transfer Line
Temp.: 150 °C
Flow Rate: 0.9 mL/min

Oven
Oven Temp.: 45 °C (hold 4.5 min) to 100 °C at 12 °C/min to 240 °C at 25 °C/min (hold 1.32 min)
Carrier Gas He, constant flow
Flow Rate: 0.9 mL/min
Detector MS
Mode: Scan
Scan Program:

Group	Start Time (min)	Scan Range (amu)	Scan Rate (scans/sec)
1	1.5	47-300	5.4
2	2.9	35-300	5.19

Transfer Line
Temp.: 240 °C
Analyzer Type: Quadrupole
Source Temp.: 230 °C
Quad Temp.: 150 °C
Electron Energy: 70 eV
Solvent Delay
Time: 1.5 min
Tune Type: BFB
Ionization Mode: EI
Instrument Agilent 7890A GC & 5975C MSD
Notes Nitrogen was used as the purge gas for the EST Encon Evolution