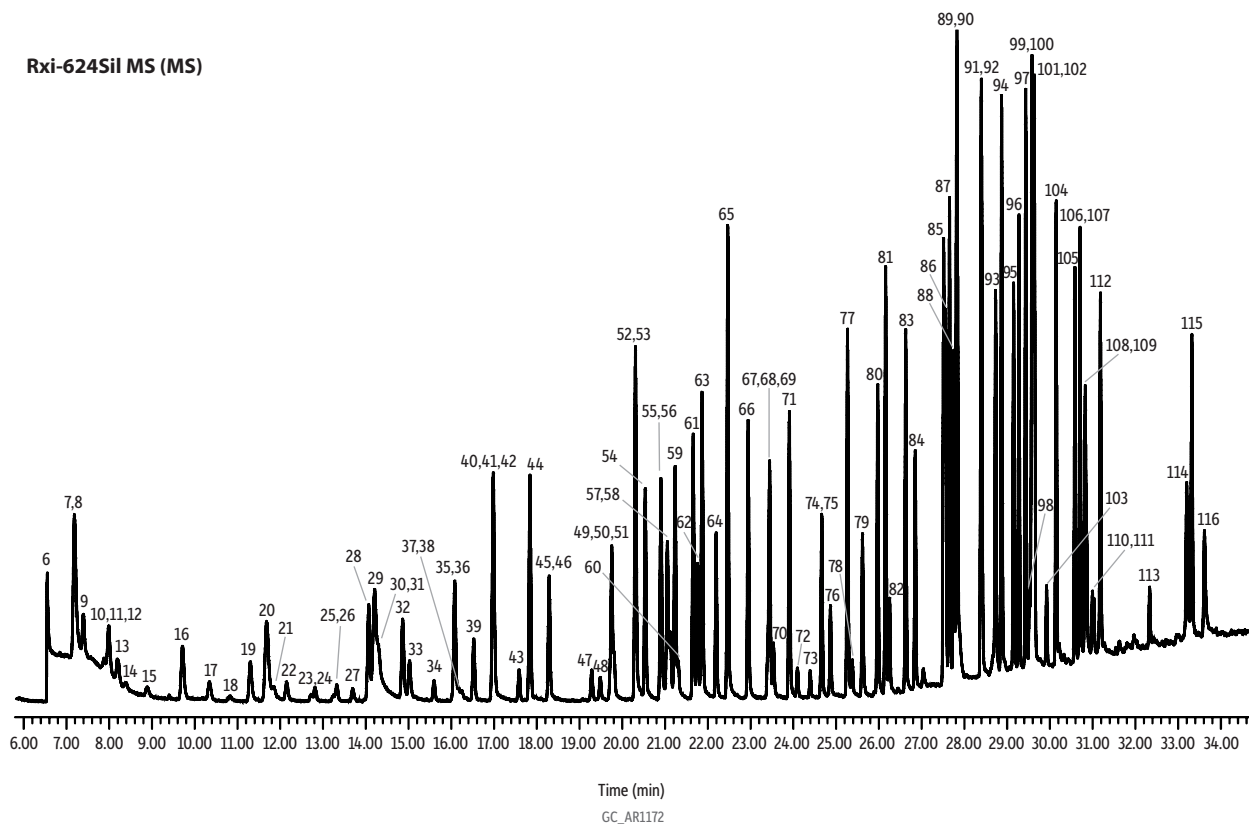
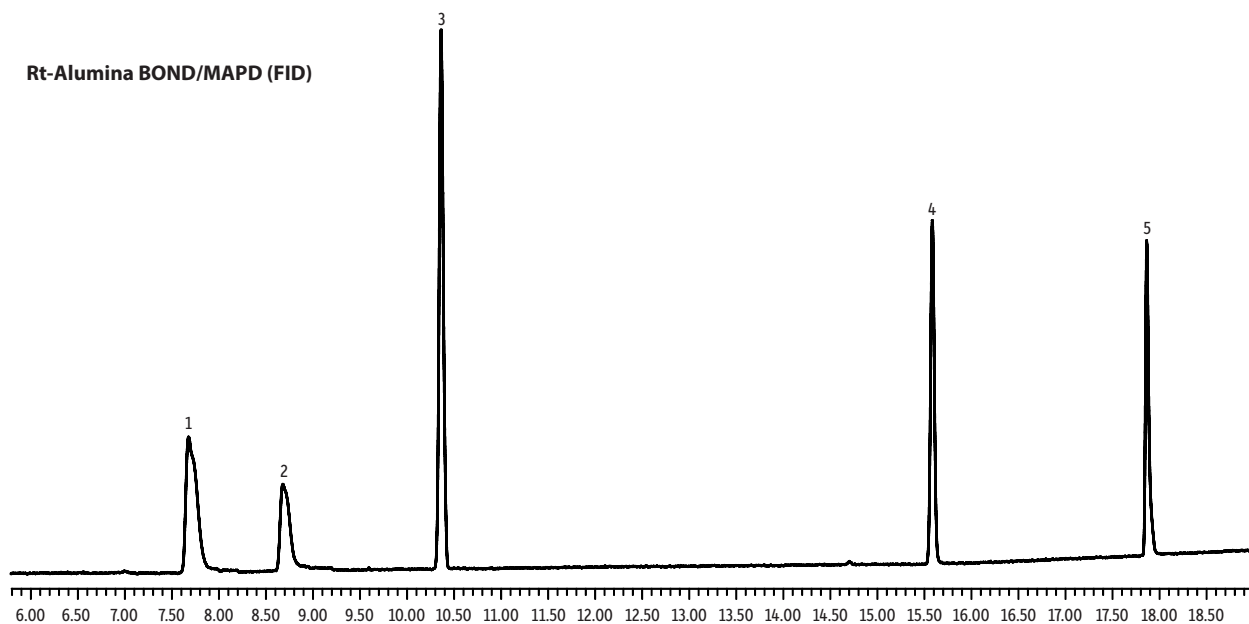


TO-15 (HJ759) + PAMS on Rt-Alumina BOND/MAPD (by FID) and Rxi-624Sil MS (by MS) using Deans Switch

- Combined TO-15 (HJ759) and PAMS analysis.
- No cryogenic cooling.



Time (min)
GC_AR11172

Peaks	tr (min)	Peaks	tr (min)
1. Ethane	7.68	59. Carbon tetrachloride	21.21
2. Ethylene	8.68	60. 3-Methylhexane	21.28
3. Propane	10.36	61. Benzene	21.63
4. Propylene	15.58	62. 1,2-Dichloroethane	21.74
5. Acetylene	17.86	63. 2,2,4-Trimethylpentane	21.84
6. Dichlorodifluoromethane	6.5	64. <i>n</i> -Heptane	22.17
7. 1,2-Dichlorotetrafluoroethane	7.14	65. 1,4-Difluorobenzene	22.45
8. Isobutane	7.19	66. Trichloroethylene	22.93
9. Chloromethane	7.35	67. Methylcyclohexane	23.38
10. <i>trans</i> -2-Butene	7.84	68. 1,2-Dichloropropane	23.42
11. <i>n</i> -Butane	7.94	69. Methyl methacrylate	23.43
12. Vinyl chloride	7.94	70. 1,4-Dioxane	23.52
13. 1,3-Butadiene	8.16	71. Bromodichloromethane	23.89
14. <i>cis</i> -2-Butene	8.35	72. 2,3,4-Trimethylpentane	24.08
15. 1-Butene	8.84	73. 2-Methylheptane	24.37
16. Bromomethane	9.67	74. 3-Methylheptane	24.64
17. Chloroethane	10.31	75. <i>cis</i> -1,3-Dichloropropene	24.65
18. Isopentane	10.77	76. Methyl isobutyl ketone	24.85
19. Vinyl bromide	11.26	77. Toluene	25.25
20. Trichlorofluoromethane	11.64	78. <i>n</i> -Octane	25.36
21. 1-Pentene	11.81	79. <i>trans</i> -1,3-Dichloropropene	25.6
22. <i>n</i> -Pentane	12.11	80. 1,1,2-Trichloroethane	25.95
23. Ethanol	12.68	81. Tetrachloroethene	26.14
24. <i>trans</i> -2-Pentene	12.77	82. 2-Hexanone	26.23
25. Isoprene	13.21	83. Dibromochloromethane	26.61
26. <i>cis</i> -2-Pentene	13.29	84. 1,2-Dibromoethane	26.84
27. Acrolein	13.65	85. Chlorobenzene-d5	27.5
28. 1,1-Dichloroethene	14.03	86. Chlorobenzene	27.55
29. 1,1,2-Trichlorotrifluoroethane	14.17	87. Ethylbenzene	27.64
30. Acetone	14.27	88. <i>n</i> -Nonane	27.74
31. 2,2-Dimethylbutane	14.27	89. <i>m</i> -Xylene	27.81
32. Carbon disulfide	14.83	90. <i>p</i> -Xylene	27.81
33. Isopropyl alcohol	14.99	91. <i>o</i> -Xylene	28.38
34. Allyl chloride	15.56	92. Styrene	28.4
35. 2,3-Dimethylbutane	16.01	93. Bromoform	28.72
36. Methylene chloride	16.05	94. Isopropylbenzene	28.86
37. 2-Methylpentane	16.14	95. 4-Bromofluorobenzene	29.14
38. Cyclopentane	16.21	96. 1,1,2,2-Tetrachloroethane	29.27
39. <i>tert</i> -Butyl alcohol	16.49	97. <i>n</i> -Propylbenzene	29.42
40. Methyl <i>tert</i> -butyl ether	16.94	98. 1,2,3-Trimethylbenzene	29.51
41. <i>trans</i> -1,2-Dichloroethene	16.95	99. <i>n</i> -Decane	29.55
42. 3-Methylpentane	16.97	100. <i>p</i> -Ethyltoluene	29.57
43. 1-Hexene	17.56	101. 2-Chlorotoluene	29.6
44. <i>n</i> -Hexane	17.81	102. 1,3,5-Trimethylbenzene	29.63
45. 1,1-Dichloroethane	18.26	103. <i>m</i> -Ethyltoluene	29.91
46. Vinyl acetate	18.28	104. 1,2,4-Trimethylbenzene	30.14
47. 2,4-Dimethylpentane	19.26	105. 1,3-Dichlorobenzene	30.58
48. Methylcyclopentane	19.46	106. <i>o</i> -Ethyltoluene	30.68
49. 2-Butanone	19.69	107. 1,4-Dichlorobenzene	30.70
50. <i>cis</i> -1,2-Dichloroethene	19.73	108. Benzyl chloride	30.82
51. Ethyl acetate	19.78	109. <i>m</i> -Diethylbenzene	30.86
52. Bromochloromethane	20.28	110. <i>p</i> -Diethylbenzene	30.99
53. Tetrahydrofuran	20.3	111. <i>n</i> -Undecane	31.02
54. Chloroform	20.51	112. 1,2-Dichlorobenzene	31.18
55. 1,1,1-Trichloroethane	20.88	113. <i>n</i> -Dodecane	32.33
56. 2-Methylhexane	20.89	114. 1,2,4-Trichlorobenzene	33.2
57. Cyclohexane	21.03	115. Hexachlorobutadiene	33.32
58. 2,3-Dimethylpentane	21.1	116. Naphthalene	33.62

Column Rxi-624Sil MS, 60 m, 0.25 mm ID, 1.40 µm (cat.# 13869)
with zero dead volume Valco internal union (cat.# 20150)
Sample Ozone precursor/PAMS mix (cat.# 26370)
75 Comp TO15 + NJ mix (cat.# 34392)
Conc.: 10 ppbv
Injection on-column
Oven
Oven Temp.: 35 °C (hold 7 min) to 60 °C at 3 °C/min to 170 °C at 10 °C/min to 250 °C at 30 °C/min (hold 6 min)
Carrier Gas He, constant flow
Flow Rate: 2.0 mL/min
Detector MS
Mode: Scan

Scan Program:	Group	Start Time (min)	Scan Range (amu)	Scan Rate (scans/sec)
	1	5.7	29-226	3.7

Transfer Line Temp.: 250 °C
Analyzer Type: Quadrupole
Source Type: Extractor
Source Temp.: 230 °C
Quad Temp.: 150 °C
Electron Energy: 70 eV
Solvent Delay Time: 5.7 min
Tune Type: BFB
Ionization Mode: EI
Preconcentrator Markes Unity 2 + CIA

Trap 1 Settings
Cooling temp.: 5 °C
Desorb temp.: 300 °C
Desorb flow: 6 mL/min
Desorb time: 180 sec

Internal Standard
Purge flow: 50 mL/min
Purge time: 60 sec
Vol.: 50 mL
ISTD flow: 50 mL/min

Standard
Size: 400 mL
Purge flow: 50 mL/min
Purge time: 60 sec
Sample flow: 100 mL/min

Instrument Notes
Agilent 7890B GC & 5977A MSD
Deans switch time: 6.489 min
Primary column to Deans switch: Rxi-624Sil MS 60 m x 0.25 mm x 1.40 µm (cat.# 13869), 2 mL/min He
Secondary column to FID: Rt-Alumina BOND/MAPD 30 m x 0.25 mm x 4 µm (cat.# 19781), 3 mL/min He
Transfer line to MS: intermediate-polarity deactivated 1 m x 0.1 mm (cat.# 10100), 3 mL/min He

FID Parameters
Temp.: 250 °C
Hz flow: 30 mL/min
Air flow: 400 mL/min
Makeup (N₂) flow: 22 mL/min
Data rate: 5 Hz