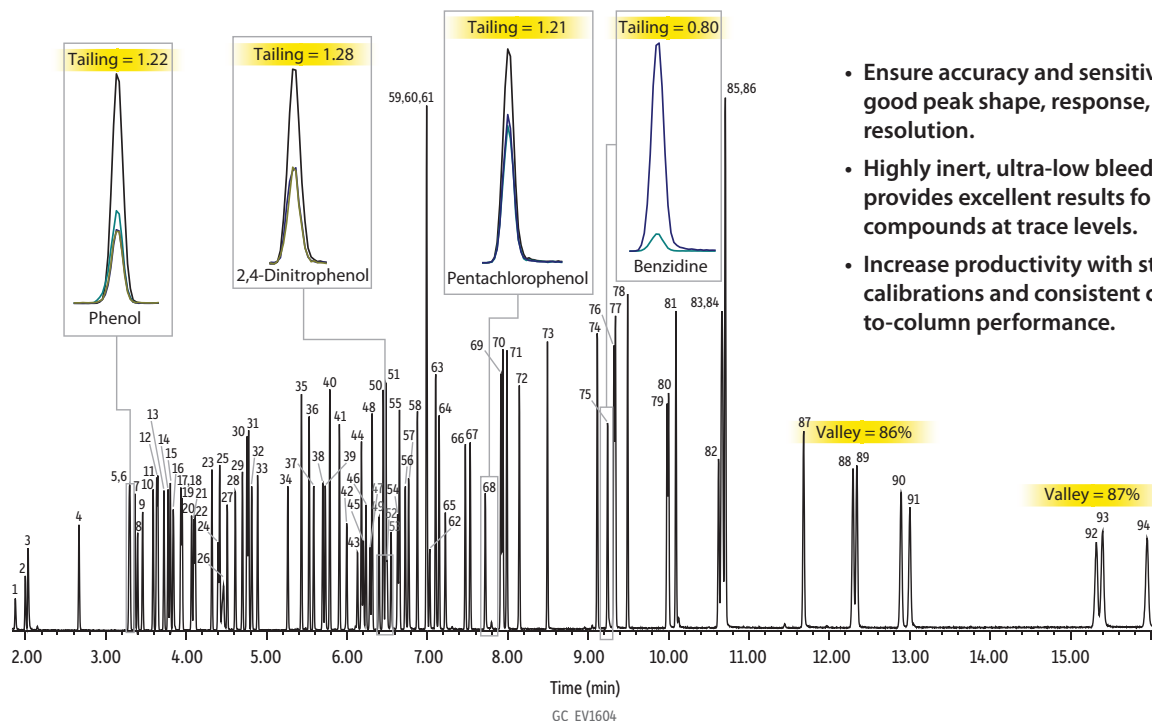


# Semivolatiles Analysis on Rxi-SVOCms



- Ensure accuracy and sensitivity with good peak shape, response, and resolution.
- Highly inert, ultra-low bleed column provides excellent results for active compounds at trace levels.
- Increase productivity with stable calibrations and consistent column-to-column performance.

Peaks	tr (min)	Peaks	tr (min)	Peaks	tr (min)	Peaks	tr (min)
1. (IS) 1,4-Dioxane-d8	1.87	25. 2,4-Dimethylphenol	4.42	49. 3-Nitroaniline	6.40	73. di-n-Butyl phthalate	8.49
2. N-Nitrosodimethylamine	2.00	26. Benzoic acid	4.46	50. (IS) Acenaphthene-D10	6.45	74. Fluoranthene	9.12
3. Pyridine	2.03	27. Bis(2-chloroethoxy)methane	4.51	51. Acenaphthene	6.48	75. Benzidine	9.24
4. (SS) 2-Fluorophenol	2.67	28. 2,4-Dichlorophenol	4.61	52. 2,4-Dinitrophenol	6.50	76. (SS) Pyrene-D10	9.32
5. (SS) Phenol-d6	3.29	29. 1,2,4-Trichlorobenzene	4.70	53. 4-Nitrophenol	6.55	77. Pyrene	9.34
6. Phenol	3.30	30. (IS) Naphthalene-D8	4.76	54. 2,4-Dinitrotoluene	6.63	78. (SS) p-Terphenyl-d14	9.49
7. Aniline	3.36	31. Naphthalene	4.78	55. Dibenzofuran	6.65	79. 3,3'-Dimethylbenzidine	9.98
8. Bis(2-chloroethyl) ether	3.40	32. 4-Chloroaniline	4.82	56. 2,3,5,6-Tetrachlorophenol	6.73	80. Butyl benzyl phthalate	10.00
9. 2-Chlorophenol	3.46	33. Hexachlorobutadiene	4.89	57. 2,3,4,6-Tetrachlorophenol	6.77	81. Bis(2-ethylhexyl) adipate	10.09
10. 1,3-Dichlorobenzene	3.59	34. 4-Chloro-3-methylphenol	5.26	58. Diethyl phthalate	6.88	82. 3,3'-Dichlorobenzidine	10.62
11. (IS) 1,4-Dichlorobenzene-D4	3.63	35. 2-Methylnaphthalene	5.43	59. 4-Chlorophenyl phenyl ether	6.99	83. Benz[a]anthracene	10.66
12. 1,4-Dichlorobenzene	3.65	36. 1-Methylnaphthalene	5.53	60. Fluorene	6.99	84. (IS) Chrysene-D12	10.67
13. Benzyl alcohol	3.72	37. Hexachlorocyclopentadiene	5.59	61. 4-Nitroaniline	7.00	85. Chrysene	10.71
14. 1,2-Dichlorobenzene	3.78	38. 2,4,6-Trichlorophenol	5.70	62. 4,6-Dinitro-2-methylphenol	7.03	86. Bis(2-ethylhexyl) phthalate	10.71
15. 2-Methylphenol	3.80	39. 2,4,5-Trichlorophenol	5.73	63. N-Nitrosodiphenylamine	7.10	87. Di-n-octyl phthalate	11.68
16. Bis(2-Chloroisopropyl)ether	3.84	40. (SS) 2-Fluorobiphenyl	5.79	64. N,N-Diphenylhydrazine	7.15	88. Benzo[b]fluoranthene	12.30
17. 4-Methylphenol	3.93	41. 2-Chloronaphthalene	5.91	65. (SS) 2,4,6-Tribromophenol	7.23	89. Benzo[k]fluoranthene	12.34
18. 3-Methylphenol	3.93	42. 2-Nitroaniline	6.00	66. 4-Bromophenyl phenyl ether	7.47	90. Benzo[a]pyrene	12.89
19. N-Nitrosodi-N-propylamine	3.95	43. 1,4-Dinitrobenzene	6.13	67. Hexachlorobenzene	7.53	91. (IS) Perylene-D12	13.00
20. Hexachloroethane	4.07	44. Dimethyl phthalate	6.18	68. Pentachlorophenol	7.72	92. Indeno[1,2,3-cd]pyrene	15.32
21. (SS) Nitrobenzene-D5	4.10	45. 1,3-Dinitrobenzene	6.20	69. (IS) Phenanthrene-D10	7.92	93. Dibenzo[a,h]anthracene	15.40
22. Nitrobenzene	4.11	46. 2,6-Dinitrotoluene	6.24	70. Phenanthrene	7.94	94. Benzo[ghi]perylene	15.95
23. Isophorone	4.32	47. 1,2-Dinitrobenzene	6.29	71. Anthracene	7.99		
24. 2-Nitrophenol	4.40	48. Acenaphthylene	6.31	72. Carbazole	8.15		

All compounds are 2 ng on column.

**Column** Rxi-SVOCms, 30 m, 0.25 mm ID, 0.25 µm (cat.# 16623)  
**Standard/Sample** 8270 MegaMix standard (cat.# 31850)  
 8270 Benzidines mix (cat.# 31852)  
 Benzoic acid (cat.# 31879)  
 Revised SV internal standard mix (cat.# 31886)  
 Revised B/N surrogate mix (cat.# 31888)  
 Acid surrogate mix (cat.# 31063)  
**Diluent:** Dichloromethane  
**Conc.:** 20 µg/mL  
**Injection**  
**Inj. Vol.:** 1 µL split (split ratio 10:1)  
**Liner:** Topaz 4.0 mm ID single taper inlet liner with wool (cat.# 23303)  
**Inj. Temp.:** 250 °C  
**Split Vent Flow Rate:** 12 mL/min  
**Oven**  
**Oven Temp.:** 60 °C (hold 0.5 min) to 285 °C at 25 °C/min to 305 °C at 3 °C/min to 330 °C at 20 °C/min (hold 5 min)  
**Carrier Gas**  
**Flow Rate:** He, constant flow  
 1.2 mL/min  
**Detector**  
 MS  
**Mode:** Scan

## Scan Program:

Group	Start Time (min)	Scan Range (amu)	Scan Rate (scans/sec)
1	1.55	35-500	5.9

**Transfer Line Temp.:** 280 °C  
**Analyzer Type:** Quadrupole  
**Source Type:** Inert  
**Drawout Plate:** 6 mm ID  
**Source Temp.:** 330 °C  
**Quad Temp.:** 180 °C  
**Electron Energy:** 70 eV  
**Tune Type:** DFTPP  
**Ionization Mode:** EI  
**Instrument** Agilent 7890A GC & 5975C MSD  
**Sample Preparation** Samples were aliquoted into amber 2 mL, 9 mm short-cap, screw-thread vials (cat.# 21143) containing glass Big Mouth inserts (cat.# 21782) and sealed with 2.0 mL, 9 mm short-cap, screw-vial closures (cat.# 23842).

