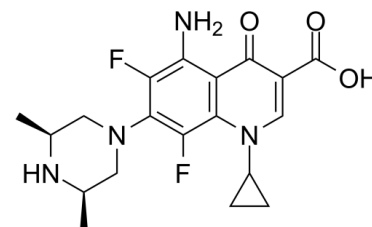


## Data Sheet

<b>Product Name:</b>	Sparfloxacin
<b>Cat. No.:</b>	CS-2347
<b>CAS No.:</b>	110871-86-8
<b>Molecular Formula:</b>	C <sub>19</sub> H <sub>22</sub> F <sub>2</sub> N <sub>4</sub> O <sub>3</sub>
<b>Molecular Weight:</b>	392.40
<b>Target:</b>	Bacterial
<b>Pathway:</b>	Anti-infection
<b>Solubility:</b>	H <sub>2</sub> O : < 0.1 mg/mL (insoluble); 0.1 M NaOH : 50 mg/mL (127.42 mM; ultrasonic and adjust pH to 11 with NaOH); DMSO : 3.33 mg/mL (8.49 mM; Need ultrasonic)



### BIOLOGICAL ACTIVITY:

Sparfloxacin (CI-978) is a fluoroquinolone antibiotic, shows broad and potent antibacterial activity. IC<sub>50</sub> & Target: Antibacterial<sup>[1]</sup>. **In Vitro:** Sparfloxacin (CI-978) shows broad and potent antibacterial activity. Its MICs for 90% of the strains tested are 0.1 to 0.78 µg/ml against gram-positive organisms, such as members of the genera *Staphylococcus*, *Streptococcus* and *Enterococcus*, and 0.0125 to 1.56 µg/ml against gram-negative organisms, such as members of the family Enterobacteriaceae and the genera *Pseudomonas*. Its MICs are 0.025 to 0.78 µg/ml against glucose nonfermenters, 0.2 to 0.78 µg/ml against anaerobes, 0.0125 to 0.05 µg/ml against *Legionella*. Sparfloxacin (CI-978) showed good oral efficacy against systemic infections with *Staphylococcus aureus*, *Streptococcus pyogenes*, *Streptococcus pneumoniae*, *Escherichia coli*, and *Pseudomonas aeruginosa* in mice<sup>[1]</sup>. Sparfloxacin targets DNA gyrase and inhibits DNA synthesis<sup>[2]</sup>.

### References:

- [1]. Nakamura, S., et al., In vitro and in vivo antibacterial activities of AT-4140, a new broad-spectrum quinolone. *Antimicrob Agents Chemother*, 1989. 33(8): p. 1167-73.
- [2]. Pan, X.S. and L.M. Fisher, Targeting of DNA gyrase in *Streptococcus pneumoniae* by sparfloxacin: selective targeting of gyrase or topoisomerase IV by quinolones. *Antimicrob Agents Chemother*, 1997. 41(2): p. 471-4.

### CAIndexNames:

3-Quinolincarboxylic acid, 5-amino-1-cyclopropyl-7-[(3R,5S)-3,5-dimethyl-1-piperazinyl]-6,8-difluoro-1,4-dihydro-4-oxo-, rel-

### SMILES:

O=C(C1=CN(C2CC2)C3=C(C(N)=C(F)C(N4C[C@H](C)N[C@H](C)C4)=C3F)C1=O)O

**Caution: Product has not been fully validated for medical applications. For research use only.**

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