

## Product Information

### 2-(4-(2,5-dioxo-2H-pyrrol-1(5H)-yl)phenyl)acetic acid

**Cat. No.:** X24-09-YYX146

**Size:** 1 g; 5 g

**CAS Number:** 91574-45-7

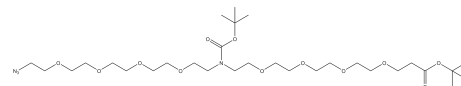
**PubChem CID:** 16780847

**Synonym:** 91574-45-7; 2-[4-(2,5-dioxo-2,5-dihydro-1H-pyrrol-1-yl)phenyl]acetic acid;

4-Maleimidophenylacetic Acid;

2-(4-(2,5-dioxo-2,5-dihydro-1H-pyrrol-1-yl)phenyl)acetic acid;

2-[4-(2,5-dioxopyrrol-1-yl)phenyl]acetic acid



**This product is for research use only and is not intended for diagnostic use.**

#### Product Information

<b>Description</b>	2-(4-(2,5-dioxo-2H-pyrrol-1(5H)-yl)phenyl)acetic acid contains a phenyl group attached to a pyrrolone moiety and an acetic acid group. It can be used as an intermediate in organic synthesis or as a building block for the synthesis of more complex molecules with potential biological activities. It may find applications in the development of drugs or in materials science.
<b>Molecular Weight</b>	231.2
<b>Molecular Formula</b>	C <sub>12</sub> H <sub>9</sub> NO <sub>4</sub>
<b>Functional Group 1</b>	Acid
<b>Functional Group 2</b>	Ketone
<b>Functional Group 3</b>	None
<b>Reactive Group 1</b>	Amine
<b>Reactive Group 2</b>	Hydrazine
<b>IUPAC Name</b>	2-[4-(2,5-Dioxopyrrol-1-yl)phenyl]acetic acid
<b>InChI</b>	InChI=1S/C12H9NO4/c14-10-5-6-11(15)13(10)9-3-1-8(2-4-9)7-12(16)17/h1-6H,7H2,(H,16,17)
<b>InChI Key</b>	KAZUDNXXBQPEPGA-UHFFFAOYSA-N
<b>Canonical SMILES</b>	C1=CC(=CC=C1CC(=O)O)N2C(=O)C=CC2=O
<b>Form</b>	Solid
<b>Identity</b>	Confirmed by NMR.
<b>Applications</b>	This compound serves as a building block in medicinal chemistry for the development of pharmaceuticals. Its structure could be used in the synthesis of various bioactive compounds or as

intermediates in the synthesis of more complex molecules.

**Storage**

Store at -20°C.