



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 12:32 PM GMT

PDB ID : 3E8R  
Title : Crystal structure of catalytic domain of TACE with hydroxamate inhibitor  
Authors : Orth, P.  
Deposited on : 2008-08-20  
Resolution : 1.90 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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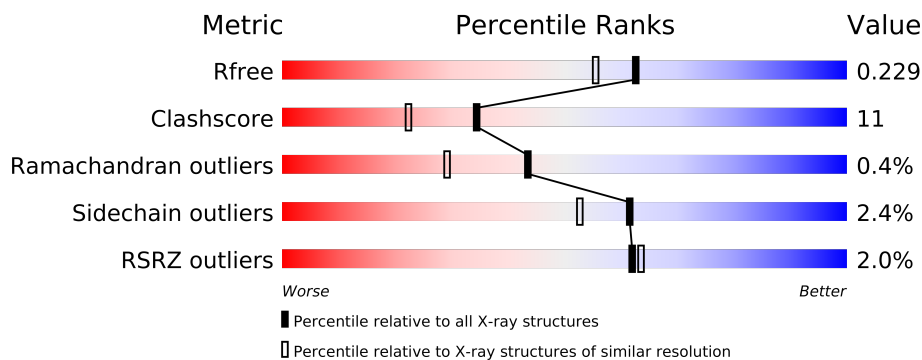
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 1.90 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	66092	3684 (1.90-1.90)
Clashscore	79885	4465 (1.90-1.90)
Ramachandran outliers	78287	4413 (1.90-1.90)
Sidechain outliers	78261	4414 (1.90-1.90)
RSRZ outliers	66119	3686 (1.90-1.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	271	
1	B	271	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	615	A	486[A]	-	X
4	INN	A	4[B]	-	X
5	CIT	B	3	-	X

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 4541 atoms, of which 0 are hydrogen and 0 are deuterium.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called ADAM 17.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	251	Total	C	N	O	S	0	2	0
			1990	1254	332	390	14			
1	B	254	Total	C	N	O	S	0	1	0
			1988	1255	333	386	14			

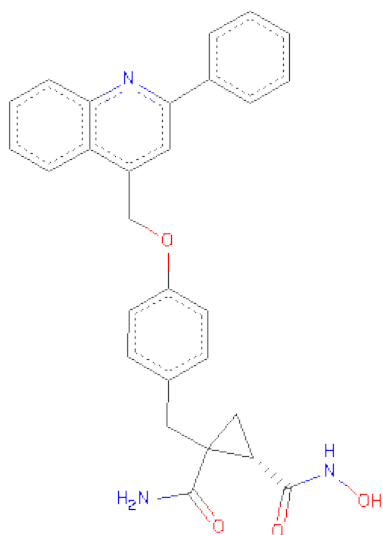
There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	266	ALA	SER	ENGINEERED	UNP P78536
A	353	GLY	VAL	ENGINEERED	UNP P78536
A	452	GLN	ASN	ENGINEERED	UNP P78536
A	478	GLY	-	EXPRESSION TAG	UNP P78536
A	479	SER	-	EXPRESSION TAG	UNP P78536
A	480	HIS	-	EXPRESSION TAG	UNP P78536
A	481	HIS	-	EXPRESSION TAG	UNP P78536
A	482	HIS	-	EXPRESSION TAG	UNP P78536
A	483	HIS	-	EXPRESSION TAG	UNP P78536
A	484	HIS	-	EXPRESSION TAG	UNP P78536
A	485	HIS	-	EXPRESSION TAG	UNP P78536
B	266	ALA	SER	ENGINEERED	UNP P78536
B	353	GLY	VAL	ENGINEERED	UNP P78536
B	452	GLN	ASN	ENGINEERED	UNP P78536
B	478	GLY	-	EXPRESSION TAG	UNP P78536
B	479	SER	-	EXPRESSION TAG	UNP P78536
B	480	HIS	-	EXPRESSION TAG	UNP P78536
B	481	HIS	-	EXPRESSION TAG	UNP P78536
B	482	HIS	-	EXPRESSION TAG	UNP P78536
B	483	HIS	-	EXPRESSION TAG	UNP P78536
B	484	HIS	-	EXPRESSION TAG	UNP P78536
B	485	HIS	-	EXPRESSION TAG	UNP P78536

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

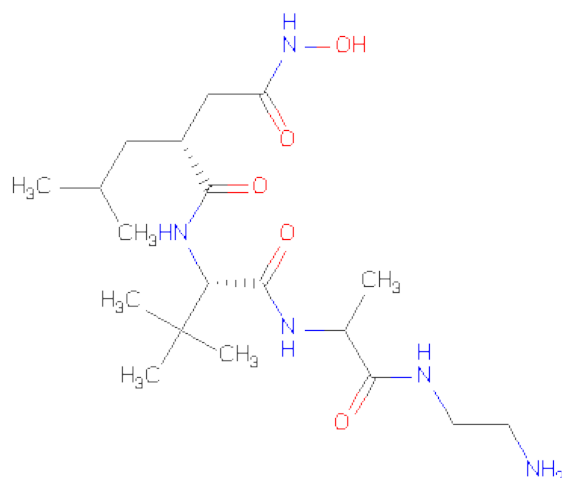
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Zn	0	0
			1	1		
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is (1R,2S)-N 2 -HYDROXY-1-{4-[(2-PHENYLQUINOLIN-4-YL)METHOXY] BENZYL}CYCLOPROPANE-1,2-DICARBOXAMIDE (three-letter code: 615) (formula: C<sub>28</sub>H<sub>25</sub>N<sub>3</sub>O<sub>4</sub>).



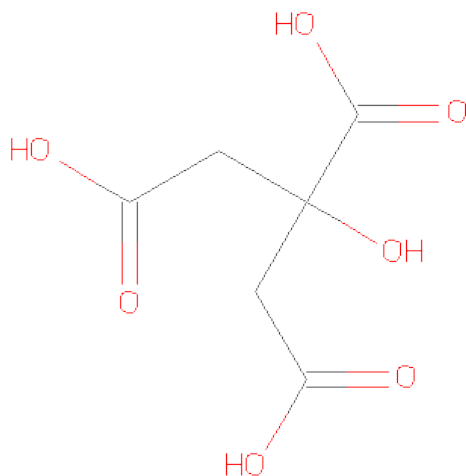
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	1
			35	28	3	4		
3	B	1	Total	C	N	O	0	0
			35	28	3	4		

- Molecule 4 is N-{(2R)-2-[2-(HYDROXYAMINO)-2-OXOETHYL]-4-METHYLPENTANOYL}-3-METHYL-L-VALYL-N-(2-AMINOETHYL)-L-ALANINAMIDE (three-letter code: INN) (formula: C<sub>19</sub>H<sub>37</sub>N<sub>5</sub>O<sub>5</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	1
			29	19	5	5		

- Molecule 5 is CITRIC ACID (three-letter code: CIT) (formula: C<sub>6</sub>H<sub>8</sub>O<sub>7</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			13	6	7		

- Molecule 6 is water.

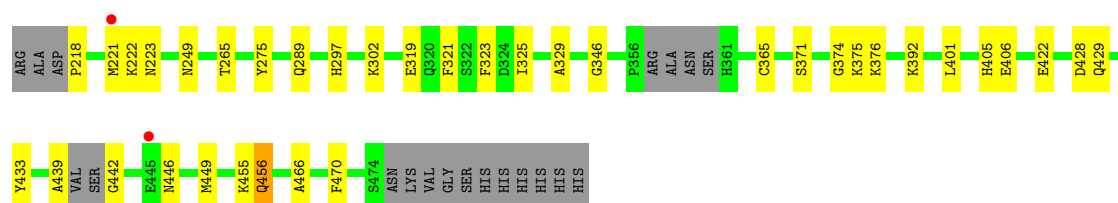
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	211	Total 211	O 211	0	0
6	B	238	Total 238	O 238	0	0

### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

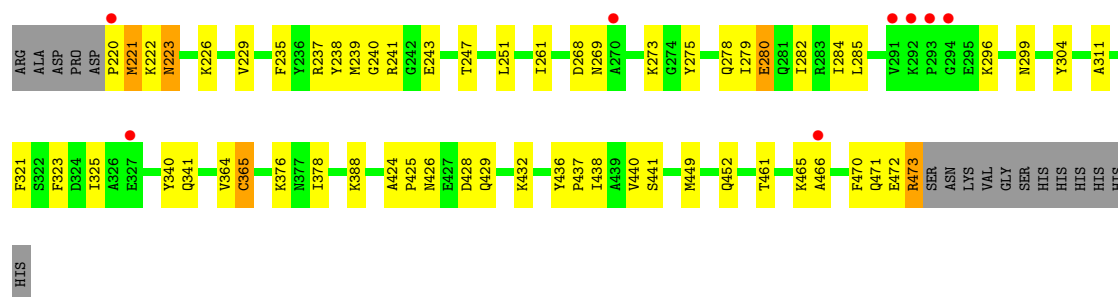
#### • Molecule 1: ADAM 17

Chain A: 



#### • Molecule 1: ADAM 17

Chain B: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	74.72Å 75.76Å 102.98Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.41 – 1.90 42.40 – 1.90	Depositor EDS
% Data completeness (in resolution range)	94.3 (42.41-1.90) 94.6 (42.40-1.90)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.04 (at 1.89Å)	Xtriage
Refinement program	BUSTER-TNT 2.1.1	Depositor
R, $R_{free}$	0.189 , 0.226 0.190 , 0.229	Depositor DCC
$R_{free}$ test set	853 reflections (1.97%)	DCC
Wilson B-factor (Å <sup>2</sup> )	26.4	Xtriage
Anisotropy	0.252	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 45.4	EDS
Estimated twinning fraction	0.015 for k,h,-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 44191 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	4541	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.08% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.



## 5 Model quality i

### 5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, 615, INN, CIT

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	0.65	0/2040	0.70	0/2751
1	B	0.60	0/2039	0.68	1/2753 (0.0%)
All	All	0.63	0/4079	0.69	1/5504 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	473	ARG	NE-CZ-NH2	-5.70	117.45	120.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Close contacts i

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1990	0	1886	29	0
1	B	1988	0	1889	44	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	35	0	23	7	0
3	B	35	0	23	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	29	0	36	11	0
5	B	13	0	6	0	0
6	A	211	0	0	10	0
6	B	238	0	0	4	0
All	All	4541	0	3863	86	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 11.

All (86) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:A:4[B]:INN:H73	4:A:4[B]:INN:HN2	1.31	0.95
1:A:446:ASN:HA	1:A:449[A]:MET:HE2	1.53	0.90
1:B:321:PHE:CZ	1:B:325:ILE:HD13	2.13	0.83
1:B:237:ARG:HD2	1:B:238:TYR:CZ	2.15	0.81
1:A:323:PHE:CE1	1:A:376:LYS:HE2	2.22	0.74
1:A:422:GLU:HG3	6:A:679:HOH:O	1.87	0.74
1:B:241:ARG:O	1:B:243:GLU:HG3	1.88	0.73
3:A:486[A]:615:H2A	6:A:634:HOH:O	1.91	0.71
1:A:302:LYS:HG2	6:A:507:HOH:O	1.90	0.70
1:B:473:ARG:HD2	6:B:576:HOH:O	1.91	0.70
1:B:247:THR:CG2	1:B:284:ILE:HD12	2.23	0.69
1:A:321:PHE:O	1:A:325:ILE:HG13	1.93	0.68
1:B:240:GLY:HA3	1:B:247:THR:OG1	1.92	0.68
1:A:439:ALA:HB1	3:A:486[A]:615:C20	2.28	0.65
1:A:319:GLU:HG3	6:A:642:HOH:O	1.98	0.63
1:A:429:GLN:OE1	6:A:686:HOH:O	2.15	0.63
1:A:374:GLY:O	6:A:629:HOH:O	2.16	0.61
1:A:371:SER:O	1:A:375:LYS:N	2.34	0.61
1:B:226:LYS:HE3	1:B:278:GLN:NE2	2.16	0.60
1:B:296:LYS:O	6:B:518:HOH:O	2.17	0.60
1:A:446:ASN:HA	1:A:449[A]:MET:CE	2.31	0.60
1:A:433:TYR:HE1	1:A:449[A]:MET:HE3	1.68	0.59
1:A:401:LEU:HD11	1:A:442:GLY:HA3	1.86	0.58
1:B:221:MET:SD	1:B:222:LYS:HG3	2.45	0.57
1:B:226:LYS:NZ	1:B:472:GLU:O	2.37	0.57
1:B:449[B]:MET:HE1	6:B:671:HOH:O	2.04	0.57
1:B:296:LYS:HD3	1:B:304:TYR:CE2	2.40	0.56
1:A:323:PHE:CZ	1:A:376:LYS:HE2	2.41	0.56
1:A:439:ALA:HB3	4:A:4[B]:INN:O2	2.06	0.55
4:A:4[B]:INN:H73	4:A:4[B]:INN:N2	2.05	0.55
4:A:4[B]:INN:H72	4:A:4[B]:INN:C4	2.36	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:325:ILE:HD13	1:A:329:ALA:HB2	1.89	0.54
1:B:247:THR:HG22	1:B:284:ILE:HD12	1.89	0.54
1:B:247:THR:HG21	1:B:284:ILE:HD12	1.89	0.54
1:B:279:ILE:HG21	1:B:282:ILE:HG13	1.91	0.53
4:A:4[B]:INN:H123	4:A:4[B]:INN:O2	2.08	0.53
1:B:311:ALA:HB2	1:B:341:GLN:HB2	1.93	0.51
1:A:249[B]:ASN:ND2	6:A:570:HOH:O	2.43	0.51
1:B:237:ARG:HD2	1:B:238:TYR:CE2	2.44	0.50
1:B:220:PRO:HG2	1:B:275:TYR:OH	2.11	0.50
1:B:235:PHE:O	1:B:239:MET:HB2	2.11	0.50
1:B:364:VAL:HG23	1:B:465:LYS:HD3	1.93	0.50
1:B:278:GLN:NE2	1:B:473:ARG:O	2.43	0.49
1:A:289:GLN:HG2	1:A:297:HIS:CG	2.48	0.49
1:B:261:ILE:HD13	6:B:650:HOH:O	2.12	0.48
4:A:4[B]:INN:H71	6:A:694:HOH:O	2.13	0.48
1:B:426:ASN:OD1	1:B:429:GLN:HG2	2.13	0.48
1:B:364:VAL:O	1:B:365:CYS:HB2	2.13	0.48
1:A:218:PRO:HB3	1:A:275:TYR:CE1	2.49	0.48
1:A:446:ASN:OD1	1:A:449[A]:MET:HE1	2.14	0.48
1:B:237:ARG:CD	1:B:238:TYR:CZ	2.93	0.48
1:B:247:THR:HG22	1:B:284:ILE:CD1	2.43	0.48
1:B:378:ILE:C	1:B:378:ILE:HD12	2.33	0.48
4:A:4[B]:INN:O3	4:A:4[B]:INN:C15	2.62	0.47
1:B:229:VAL:CG1	1:B:251:LEU:HD13	2.44	0.47
3:A:486[A]:615:H26	6:A:557:HOH:O	2.15	0.46
1:A:221:MET:O	1:A:222:LYS:HG2	2.16	0.45
1:B:223:ASN:C	1:B:223:ASN:HD22	2.19	0.45
1:B:364:VAL:CG2	1:B:465:LYS:HD3	2.47	0.45
1:B:466:ALA:HA	1:B:470:PHE:CG	2.52	0.45
1:B:438:ILE:HG22	1:B:440:VAL:HG23	1.98	0.45
1:A:325:ILE:C	1:A:325:ILE:HD12	2.37	0.45
1:B:436:TYR:CG	1:B:437:PRO:HD2	2.52	0.44
4:A:4[B]:INN:C12	4:A:4[B]:INN:O2	2.65	0.44
1:A:346:GLY:O	4:A:4[B]:INN:N2	2.51	0.44
1:B:226:LYS:HB3	1:B:280:GLU:HB2	2.00	0.43
1:B:424:ALA:N	1:B:425:PRO:HD3	2.33	0.43
1:A:466:ALA:HA	1:A:470:PHE:CG	2.53	0.43
3:A:486[A]:615:H2	3:A:486[A]:615:HN3A	1.44	0.43
1:B:436:TYR:CD1	1:B:437:PRO:HD2	2.53	0.43
1:B:323:PHE:CE1	1:B:376:LYS:HE2	2.54	0.43
1:A:405:HIS:CE1	4:A:4[B]:INN:H1	2.54	0.43
1:A:401:LEU:HB3	3:A:486[A]:615:H10	2.01	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:340:TYR:CE1	1:B:388:LYS:HB2	2.53	0.42
1:B:269:ASN:ND2	1:B:452:GLN:OE1	2.48	0.42
1:B:285:LEU:HD13	1:B:299:ASN:ND2	2.34	0.42
1:B:226:LYS:HG3	1:B:471:GLN:NE2	2.35	0.42
3:A:486[A]:615:H23A	6:A:696:HOH:O	2.19	0.42
1:A:456:GLN:HB3	1:A:456:GLN:HE21	1.66	0.42
1:B:268:ASP:O	1:B:269:ASN:HB2	2.19	0.42
1:A:406:GLU:OE1	3:A:486[A]:615:O4	2.37	0.42
1:A:265:THR:HG23	1:A:455:LYS:HE2	2.02	0.42
3:B:486:615:H2	3:B:486:615:HN3A	1.71	0.41
1:B:432:LYS:HE2	1:B:436:TYR:CZ	2.55	0.40
4:A:4[B]:INN:O1	4:A:4[B]:INN:H72	2.20	0.40
1:B:461:THR:HG23	1:B:465:LYS:HD2	2.03	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	247/271 (91%)	241 (98%)	5 (2%)	1 (0%)	43	29
1	B	253/271 (93%)	248 (98%)	4 (2%)	1 (0%)	43	29
All	All	500/542 (92%)	489 (98%)	9 (2%)	2 (0%)	43	29

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	365	CYS
1	B	365	CYS

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of

similar resolution. The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	213/231 (92%)	209 (98%)	4 (2%)	69	63
1	B	210/231 (91%)	204 (97%)	6 (3%)	55	44
All	All	423/462 (92%)	413 (98%)	10 (2%)	61	53

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	223	ASN
1	A	392	LYS
1	A	428	ASP
1	A	456	GLN
1	B	221	MET
1	B	223	ASN
1	B	273	LYS
1	B	280	GLU
1	B	428	ASP
1	B	441	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	223	ASN
1	A	456	GLN
1	B	223	ASN
1	B	281	GLN
1	B	471	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates

There are no carbohydrates in this entry.

## 5.6 Ligand geometry

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z  > 2$	Counts	RMSZ	# $ Z  > 2$
4	INN	A	4[B]	2	28,28,28	1.18	3 (10%)	38,38,38	3.83	17 (44%)
3	615	A	486[A]	2	39,39,39	1.30	5 (12%)	56,56,56	3.35	16 (28%)
5	CIT	B	3	-	12,12,12	1.50	3 (25%)	17,17,17	1.91	4 (23%)
3	615	B	486	2	39,39,39	1.33	2 (5%)	56,56,56	3.03	16 (28%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	INN	A	4[B]	2	3/3/9/13	0/40/40/40	0/0/0/0
3	615	A	486[A]	2	1/1/4/5	0/26/34/34	0/2/5/5
5	CIT	B	3	-	-	0/16/16/16	0/0/0/0
3	615	B	486	2	1/1/4/5	0/26/34/34	0/2/5/5

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	486	615	C23-C21	-3.90	1.43	1.51
3	A	486[A]	615	C22-C21	3.54	1.61	1.54
5	B	3	CIT	C3-C6	2.86	1.56	1.53
5	B	3	CIT	O6-C6	2.72	1.40	1.30
3	B	486	615	C23-C22	2.70	1.55	1.51
4	A	4[B]	INN	C6-C5	2.30	1.59	1.55

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	486[A]	615	C23-C22	-2.30	1.47	1.51
4	A	4[B]	INN	CA-C4	2.27	1.55	1.51
3	A	486[A]	615	C2-C22	-2.23	1.51	1.55
3	A	486[A]	615	C17-C16	2.21	1.41	1.36
4	A	4[B]	INN	C5-N1	2.20	1.49	1.45
5	B	3	CIT	C2-C3	2.20	1.56	1.53
3	A	486[A]	615	C18-C19	2.08	1.41	1.36

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	486[A]	615	C23-C21-C3	17.50	153.67	118.26
3	B	486	615	C23-C21-C3	16.70	152.05	118.26
4	A	4[B]	INN	C7-C6-C8	-11.08	86.86	108.77
4	A	4[B]	INN	C7-C6-C9	-10.64	87.73	108.77
4	A	4[B]	INN	C7-C6-C5	-9.46	90.93	109.85
3	A	486[A]	615	C22-C21-C3	-9.13	100.85	119.89
3	A	486[A]	615	C23-C22-C1	-6.38	94.61	116.67
4	A	4[B]	INN	C8-C6-C5	6.36	122.56	109.85
3	B	486	615	O4-N2-C3	-6.15	111.53	119.80
4	A	4[B]	INN	C10-C5-N1	6.11	123.40	109.16
4	A	4[B]	INN	C5-N1-C4	5.69	134.62	122.04
5	B	3	CIT	O6-C6-C3	5.66	121.12	112.89
3	A	486[A]	615	C22-C23-C21	5.14	65.91	61.95
3	B	486	615	C23-C21-C22	5.06	62.86	59.40
4	A	4[B]	INN	C9-C6-C5	5.06	119.97	109.85
3	A	486[A]	615	C10-O3-C7	4.88	130.71	117.66
3	B	486	615	C23-C22-C1	-4.83	99.96	116.67
3	A	486[A]	615	C23-C22-C2	4.49	126.05	118.34
4	A	4[B]	INN	C9-C6-C8	4.30	117.27	108.77
3	B	486	615	C10-O3-C7	4.29	129.13	117.66
3	A	486[A]	615	C2-C22-C1	-4.20	103.93	117.33
3	B	486	615	C10-C11-C15	4.19	125.93	120.27
3	B	486	615	C10-C11-C13	-3.89	115.16	121.33
3	A	486[A]	615	C23-C21-C22	-3.77	56.83	59.40
3	B	486	615	C23-C22-C21	-3.63	54.83	58.92
3	A	486[A]	615	C10-C11-C15	3.46	124.95	120.27
3	A	486[A]	615	O4-N2-C3	-3.40	115.22	119.80
3	A	486[A]	615	C10-C11-C13	-3.33	116.04	121.33
3	A	486[A]	615	O3-C10-C11	3.30	117.95	109.29
3	B	486	615	C22-C2-C4	-3.18	111.08	116.22
4	A	4[B]	INN	C11-N2-C10	3.09	128.20	121.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	486[A]	615	C2-C22-C21	3.06	130.98	119.23
3	B	486	615	O3-C10-C11	3.04	117.26	109.29
3	B	486	615	C21-C3-N2	2.99	119.45	115.51
3	B	486	615	C22-C21-C3	-2.96	113.72	119.89
4	A	4[B]	INN	CA-CB-C1	2.91	119.50	115.95
4	A	4[B]	INN	O4-N-C	-2.75	115.49	119.57
4	A	4[B]	INN	O1-C4-CA	-2.67	119.09	121.89
3	B	486	615	C2-C22-C21	2.67	129.48	119.23
4	A	4[B]	INN	C6-C5-N1	2.63	114.81	112.29
4	A	4[B]	INN	C6-C5-C10	2.58	115.49	112.85
3	B	486	615	C2-C22-C1	-2.57	109.13	117.33
3	A	486[A]	615	C21-C22-C1	2.49	125.06	119.28
5	B	3	CIT	C3-C4-C5	-2.46	107.82	113.77
3	A	486[A]	615	C12-N1-C14	2.36	119.86	118.05
3	A	486[A]	615	C22-C1-N3	-2.30	114.95	117.37
3	B	486	615	C23-C22-C2	2.22	122.16	118.34
4	A	4[B]	INN	C13-C11-N2	-2.22	106.45	111.76
3	B	486	615	C12-N1-C14	2.18	119.72	118.05
5	B	3	CIT	O6-C6-O5	-2.12	117.10	123.76
5	B	3	CIT	O7-C3-C6	2.09	111.97	108.95
4	A	4[B]	INN	CA-C0-C	2.04	116.85	112.23
4	A	4[B]	INN	O2-C10-C5	-2.04	117.92	121.05

All (5) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	B	486	615	C22
4	A	4[B]	INN	C5
4	A	4[B]	INN	CA
4	A	4[B]	INN	C11
3	A	486[A]	615	C22

There are no torsion outliers.

There are no ring outliers.

## 5.7 Other polymers ⓘ

There are no such residues in this entry.



## 5.8 Polymer linkage issues

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	251/271 (92%)	-0.25	2 (0%) 83 85	16, 25, 47, 70	0
1	B	254/271 (93%)	0.08	8 (3%) 47 48	19, 32, 58, 71	0
All	All	505/542 (93%)	-0.09	10 (1%) 62 63	16, 29, 55, 71	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	293	PRO	5.4
1	B	294	GLY	4.2
1	B	466	ALA	3.1
1	B	292	LYS	2.8
1	B	270	ALA	2.6
1	B	220	PRO	2.4
1	A	445	GLU	2.2
1	A	221	MET	2.1
1	B	291	VAL	2.1
1	B	327	GLU	2.0

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	615	A	486[A]	35/35	0.28	6.38	3,24,54,59	35
4	INN	A	4[B]	29/29	0.23	6.13	15,32,89,99	29
5	CIT	B	3	13/13	0.17	3.18	36,37,40,42	13
3	615	B	486	35/35	0.10	0.33	18,24,30,36	0
2	ZN	A	1	1/1	0.07	-0.96	24,24,24,24	0
2	ZN	B	2	1/1	0.05	-2.71	25,25,25,25	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.