



# Full wwPDB X-ray Structure Validation Report ⓘ

May 20, 2026 – 08:08 PM EDT

PDB ID : 9YAS / pdb\_00009yas  
Title : Rana catesbeiana saxiphilin mutant - Y558A:STX-C13-OBz (co-crystal)  
Authors : Chen, Z.; Zakrzewska, S.; Minor, D.L.  
Deposited on : 2025-09-16  
Resolution : 2.45 Å(reported)

This is a Full wwPDB X-ray Structure 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

---

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

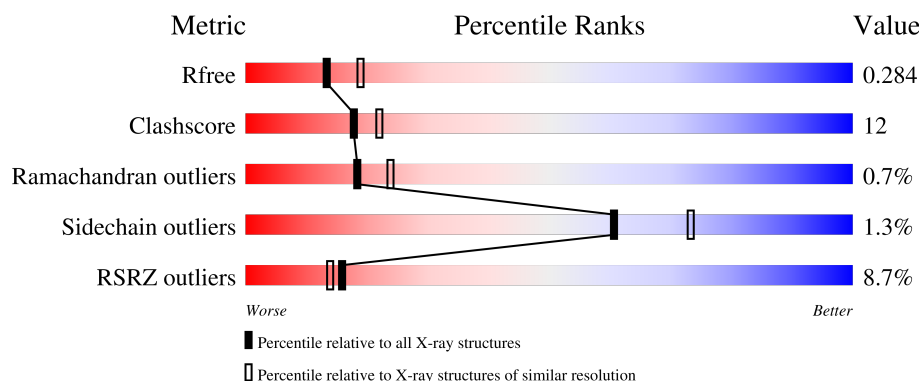
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.45 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))
$R_{free}$	180053	1190 (2.46-2.46)
Clashscore	190562	1229 (2.46-2.46)
Ramachandran outliers	187476	1218 (2.46-2.46)
Sidechain outliers	187428	1218 (2.46-2.46)
RSRZ outliers	180081	1190 (2.46-2.46)

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 of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	853	<div> <div>9%</div> <div>74%</div> <div>22%</div> <div>..</div> </div>
1	B	853	<div> <div>8%</div> <div>69%</div> <div>26%</div> <div>..</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 12881 atoms, of which 0 are hydrogens and 0 are deuteriums.

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.

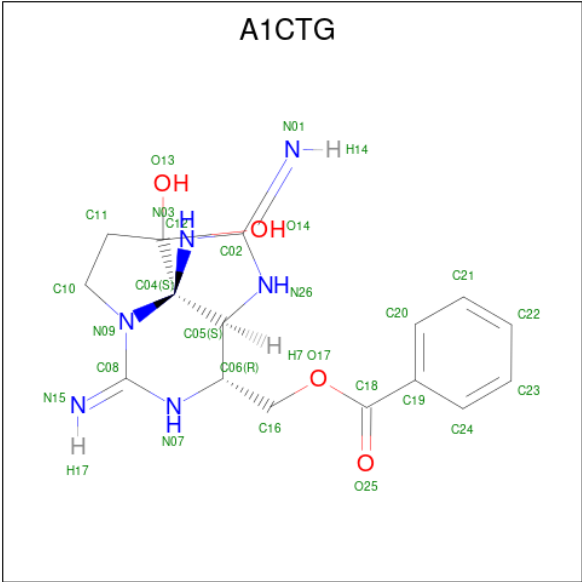
- Molecule 1 is a protein called Saxiphilin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	818	Total	C	N	O	S	0	0	0
			6308	3939	1090	1220	59			
1	B	818	Total	C	N	O	S	0	0	0
			6308	3939	1090	1220	59			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	558	ALA	TYR	engineered mutation	UNP P31226
A	826	SER	-	expression tag	UNP P31226
A	827	ASN	-	expression tag	UNP P31226
A	828	SER	-	expression tag	UNP P31226
A	829	LEU	-	expression tag	UNP P31226
A	830	GLU	-	expression tag	UNP P31226
A	831	VAL	-	expression tag	UNP P31226
A	832	LEU	-	expression tag	UNP P31226
A	833	PHE	-	expression tag	UNP P31226
A	834	GLN	-	expression tag	UNP P31226
B	558	ALA	TYR	engineered mutation	UNP P31226
B	826	SER	-	expression tag	UNP P31226
B	827	ASN	-	expression tag	UNP P31226
B	828	SER	-	expression tag	UNP P31226
B	829	LEU	-	expression tag	UNP P31226
B	830	GLU	-	expression tag	UNP P31226
B	831	VAL	-	expression tag	UNP P31226
B	832	LEU	-	expression tag	UNP P31226
B	833	PHE	-	expression tag	UNP P31226
B	834	GLN	-	expression tag	UNP P31226

- Molecule 2 is [(2Z,3aS,4R,6Z,7R,10aS)-10,10-dihydroxy-2,6-diiminooctahydro-1H,8H-pyrrolo[1,2-c]purin-4-yl]methyl benzoate (CCD ID: A1CTG) (formula: C<sub>16</sub>H<sub>20</sub>N<sub>6</sub>O<sub>4</sub>) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			26	16	6	4		
2	B	1	Total	C	N	O	0	0
			26	16	6	4		

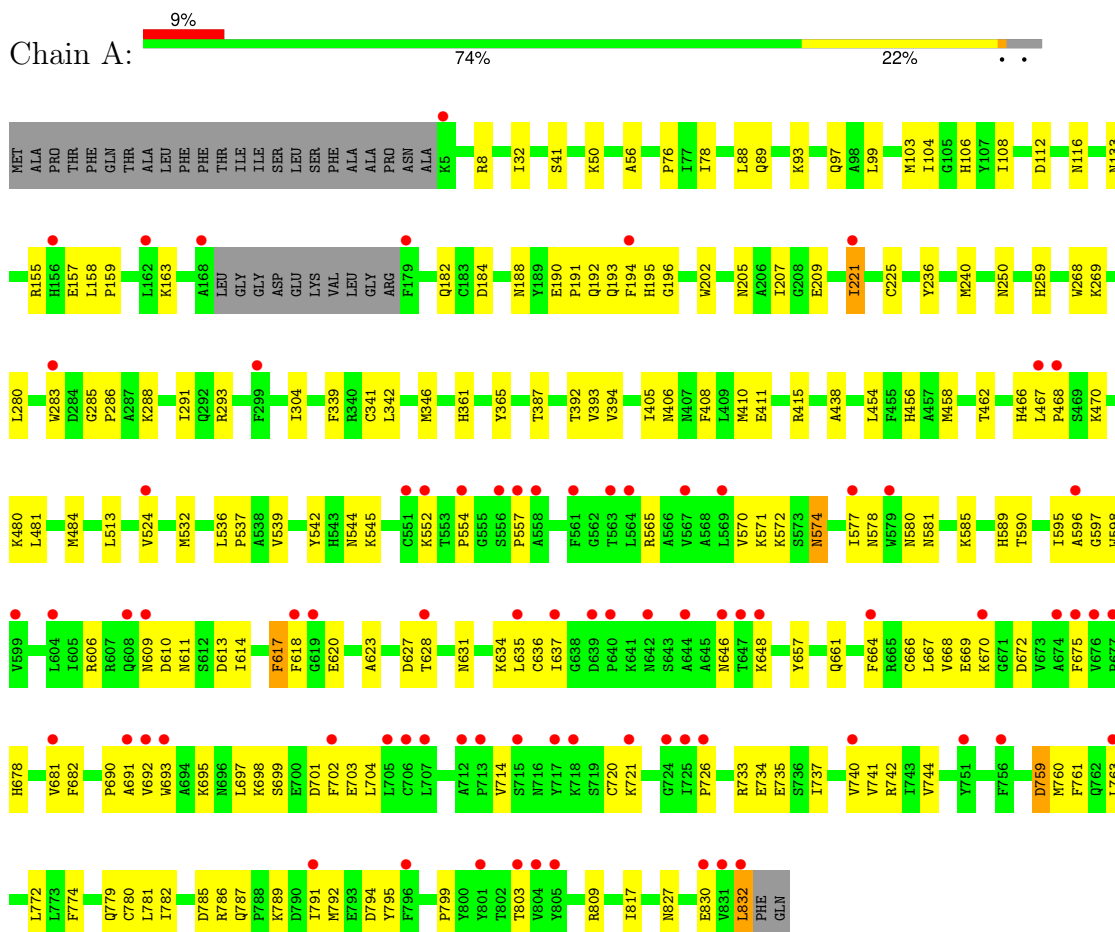
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	112	Total	O	0	0
			112	112		
3	B	101	Total	O	0	0
			101	101		

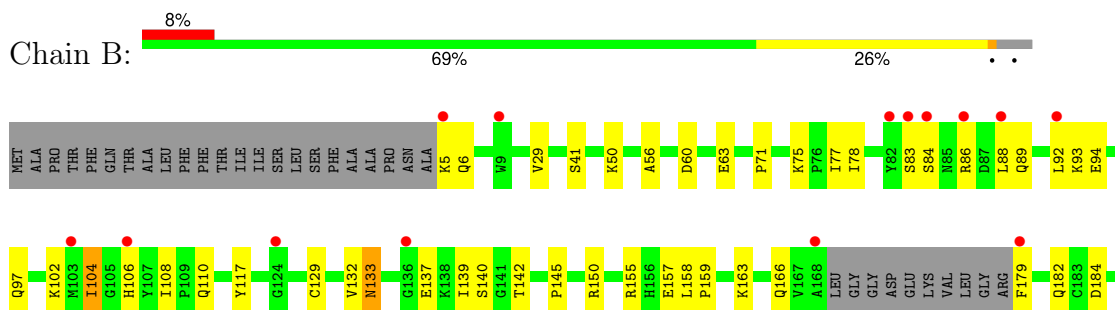
### 3 Residue-property plots [i](#)

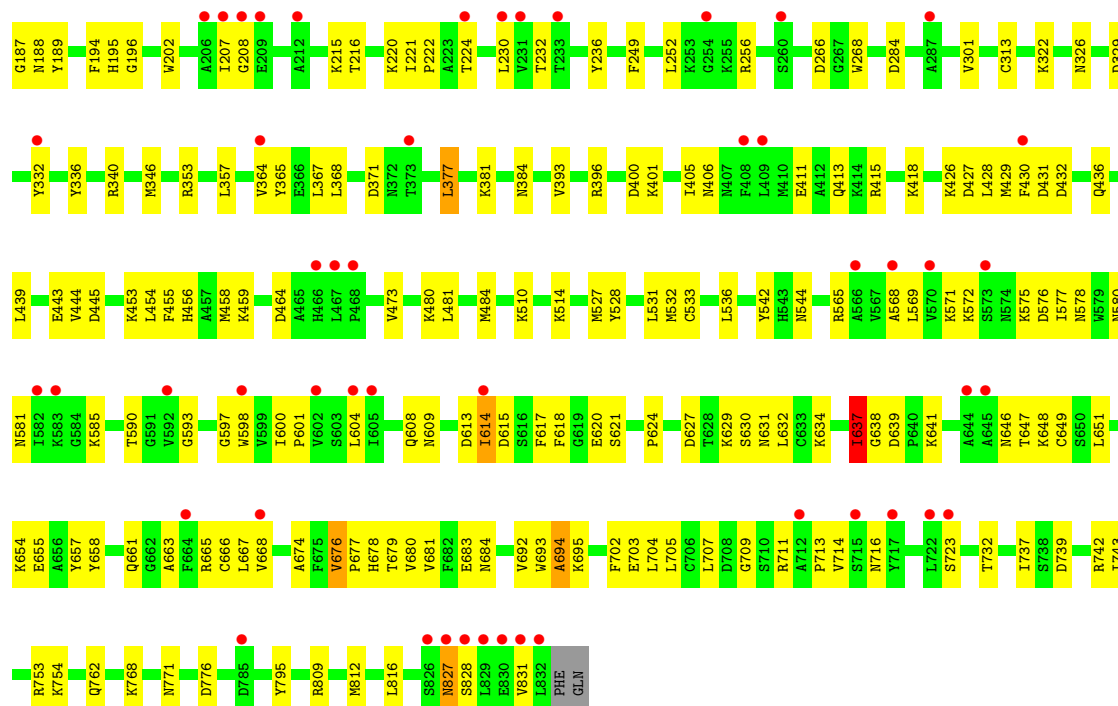
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Saxiphilin



#### • Molecule 1: Saxiphilin





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	96.20Å 109.33Å 254.69Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.10 – 2.45 48.10 – 2.45	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.10-2.45) 99.9 (48.10-2.45)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.53 (at 2.45Å)	Xtriage
Refinement program	PHENIX (1.21.1_5286: ???)	Depositor
R, $R_{free}$	0.249 , 0.284 0.250 , 0.284	Depositor DCC
$R_{free}$ test set	4977 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	63.4	Xtriage
Anisotropy	0.665	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 59.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	12881	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	97.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.89% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: A1CTG

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	0.19	0/6433	0.36	0/8682
1	B	0.21	0/6433	0.39	0/8682
All	All	0.20	0/12866	0.37	0/17364

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts

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 hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	6308	0	6177	139	0
1	B	6308	0	6177	168	0
2	A	26	0	0	3	0
2	B	26	0	0	0	0
3	A	112	0	0	7	0
3	B	101	0	0	13	0
All	All	12881	0	12354	307	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 12.

All (307) close contacts within the same asymmetric unit are listed below, sorted by their clash



magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:341:CYS:SG	3:A:1068:HOH:O	2.19	1.00
1:B:703:GLU:OE1	1:B:711:ARG:NH2	1.98	0.95
1:A:192:GLN:NE2	1:A:225:CYS:SG	2.43	0.91
1:B:84:SER:HB3	1:B:232:THR:H	1.36	0.90
1:A:572:LYS:HD2	1:A:701:ASP:HA	1.59	0.84
1:B:145:PRO:O	1:B:150:ARG:NH2	2.12	0.82
1:A:288:LYS:H	1:A:288:LYS:HD2	1.44	0.81
1:A:240:MET:HE1	1:A:339:PHE:HD1	1.47	0.79
1:A:636:CYS:SG	3:A:1065:HOH:O	2.42	0.78
1:B:577:ILE:HD13	1:B:585:LYS:HG3	1.67	0.75
1:A:589:HIS:HD2	1:A:623:ALA:HB2	1.52	0.75
1:A:346:MET:HB2	3:A:1068:HOH:O	1.85	0.74
1:B:236:TYR:HD2	1:B:384:ASN:HD22	1.35	0.74
1:A:704:LEU:HG	1:A:714:VAL:HA	1.71	0.71
1:A:698:LYS:NZ	1:A:699:SER:OG	2.24	0.70
1:A:293:ARG:NH1	3:A:1004:HOH:O	2.24	0.69
1:B:753:ARG:NH2	3:B:1008:HOH:O	2.24	0.69
1:A:41:SER:OG	3:A:1001:HOH:O	2.09	0.69
1:A:590:THR:HG22	1:A:597:GLY:HA3	1.74	0.69
1:A:240:MET:HE1	1:A:339:PHE:CD1	2.27	0.68
1:B:56:ALA:HB2	1:B:405:ILE:HD13	1.75	0.68
1:B:590:THR:HG22	1:B:597:GLY:HA3	1.74	0.68
1:A:606:ARG:HH12	1:A:614:ILE:N	1.92	0.68
1:B:527:MET:HE3	1:B:531:LEU:HD21	1.76	0.67
1:B:155:ARG:HG2	1:B:157:GLU:HG2	1.75	0.67
1:B:94:GLU:OE1	1:B:117:TYR:OH	2.12	0.67
1:B:78:ILE:HB	1:B:393:VAL:HB	1.78	0.66
1:B:207:ILE:HD11	1:B:230:LEU:HD13	1.77	0.66
1:B:676:VAL:HG22	1:B:677:PRO:HD2	1.77	0.66
1:B:216:THR:HG21	1:B:222:PRO:HA	1.78	0.66
1:B:145:PRO:C	1:B:150:ARG:NH2	2.54	0.65
1:B:129:CYS:H	1:B:150:ARG:NH1	1.95	0.65
1:B:194:PHE:HE1	1:B:196:GLY:HA2	1.61	0.65
1:A:552:LYS:O	1:A:554:PRO:HD3	1.96	0.65
1:B:133:ASN:HD21	1:B:137:GLU:HB2	1.60	0.65
1:B:704:LEU:HD13	1:B:714:VAL:HA	1.78	0.65
1:B:184:ASP:OD2	1:B:188:ASN:N	2.30	0.64
1:B:75:LYS:NZ	1:B:444:VAL:O	2.31	0.64
1:B:140:SER:H	1:B:831:VAL:HG23	1.62	0.64
1:B:158:LEU:HD23	1:B:163:LYS:HG3	1.81	0.63
1:B:473:VAL:HG23	1:B:743:ILE:HD12	1.79	0.63

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:661:GLN:NE2	3:A:1003:HOH:O	2.23	0.63
1:B:455:PHE:O	3:B:1002:HOH:O	2.16	0.63
1:B:615:ASP:O	3:B:1001:HOH:O	2.15	0.62
1:A:184:ASP:OD2	1:A:188:ASN:N	2.26	0.62
1:A:532:MET:HG2	1:A:817:ILE:HG23	1.82	0.62
1:B:381:LYS:HD3	1:B:827:ASN:HB2	1.81	0.62
1:B:637:ILE:HG23	1:B:638:GLY:N	2.14	0.62
1:B:108:ILE:O	1:B:110:GLN:NE2	2.31	0.62
1:A:799:PRO:O	1:A:803:THR:HG23	1.99	0.62
1:A:668:VAL:HG21	1:A:697:LEU:HD11	1.82	0.62
1:B:578:ASN:ND2	1:B:709:GLY:O	2.33	0.61
1:A:737:ILE:O	1:A:741:VAL:HG23	1.99	0.61
1:B:129:CYS:N	1:B:150:ARG:NH1	2.48	0.61
1:B:576:ASP:O	1:B:581:ASN:ND2	2.34	0.61
1:A:735:GLU:N	1:A:735:GLU:OE1	2.34	0.61
1:B:60:ASP:HB3	1:B:63:GLU:HG3	1.81	0.61
1:A:565:ARG:HB2	1:A:678:HIS:HD2	1.66	0.60
1:B:145:PRO:C	1:B:150:ARG:HH22	2.08	0.60
1:A:268:TRP:HZ2	1:A:291:ILE:HG23	1.66	0.60
1:B:620:GLU:HG2	1:B:634:LYS:HD3	1.82	0.60
1:B:5:LYS:NZ	3:B:1010:HOH:O	2.31	0.60
1:B:129:CYS:HB2	1:B:150:ARG:HH11	1.68	0.59
1:B:676:VAL:HG22	1:B:680:VAL:HG21	1.84	0.59
1:B:667:LEU:HB2	1:B:674:ALA:HB2	1.85	0.59
1:A:207:ILE:HD11	1:A:209:GLU:CD	2.28	0.59
1:B:29:VAL:HG13	1:B:411:GLU:HG2	1.85	0.59
1:A:89:GLN:N	1:A:89:GLN:OE1	2.35	0.58
1:B:624:PRO:HA	1:B:649:CYS:HA	1.86	0.58
1:B:665:ARG:HE	1:B:693:TRP:CD1	2.20	0.58
1:A:466:HIS:HE1	1:A:468:PRO:HA	1.67	0.58
1:A:571:LYS:HG3	1:A:667:LEU:HD11	1.83	0.58
1:B:182:GLN:N	1:B:182:GLN:OE1	2.36	0.58
1:A:466:HIS:CE1	1:A:468:PRO:HA	2.38	0.58
1:B:93:LYS:O	1:B:97:GLN:HG3	2.02	0.58
1:B:381:LYS:HB3	1:B:827:ASN:HB2	1.86	0.58
1:B:637:ILE:O	3:B:1003:HOH:O	2.16	0.58
1:A:191:PRO:HG2	1:A:809:ARG:HD2	1.85	0.57
1:A:668:VAL:HG13	1:A:669:GLU:OE2	2.04	0.57
1:B:182:GLN:HE22	1:B:194:PHE:H	1.51	0.57
1:B:357:LEU:HD13	1:B:377:LEU:HG	1.87	0.57
1:A:578:ASN:OD1	1:A:580:ASN:N	2.34	0.57

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:681:VAL:HG13	1:A:682:PHE:CD2	2.39	0.57
1:B:129:CYS:CB	1:B:150:ARG:HH11	2.16	0.57
1:A:158:LEU:HD12	1:A:163:LYS:HE3	1.87	0.57
1:B:189:TYR:HE1	1:B:207:ILE:O	1.88	0.56
1:B:569:LEU:HD22	1:B:681:VAL:HG22	1.87	0.56
1:B:353:ARG:NH1	1:B:431:ASP:OD2	2.38	0.56
1:B:578:ASN:HB3	1:B:580:ASN:OD1	2.06	0.55
1:A:133:ASN:HD22	1:A:830:GLU:HB3	1.70	0.55
1:A:76:PRO:HB3	1:A:392:THR:HG21	1.89	0.55
1:A:269:LYS:HE2	1:A:458:MET:HE1	1.87	0.55
1:A:609:ASN:OD1	1:A:617:PHE:HE1	1.90	0.55
1:B:647:THR:HB	1:B:654:LYS:HB3	1.88	0.55
1:B:565:ARG:HD2	1:B:678:HIS:CD2	2.42	0.55
1:B:86:ARG:HD2	1:B:230:LEU:HB3	1.89	0.54
1:B:639:ASP:OD2	1:B:641:LYS:HB3	2.07	0.54
1:A:574:ASN:HB3	1:A:577:ILE:HD13	1.88	0.54
1:B:182:GLN:NE2	1:B:194:PHE:H	2.06	0.54
1:A:56:ALA:HB2	1:A:405:ILE:HD13	1.90	0.54
1:A:236:TYR:CE1	1:A:387:THR:HG22	2.42	0.54
1:B:754:LYS:O	1:B:754:LYS:HG2	2.07	0.53
1:B:665:ARG:HE	1:B:693:TRP:CG	2.26	0.53
1:A:631:ASN:HA	1:A:634:LYS:HE2	1.89	0.53
1:A:795:TYR:O	2:A:901:A1CTG:N15	2.41	0.53
1:B:568:ALA:HB3	1:B:705:LEU:HB2	1.90	0.53
1:A:182:GLN:HE22	1:A:194:PHE:H	1.55	0.53
1:A:589:HIS:CD2	1:A:623:ALA:HB2	2.38	0.53
1:A:539:VAL:HG11	1:A:781:LEU:HB3	1.90	0.53
1:B:179:PHE:HB2	1:B:195:HIS:CE1	2.44	0.53
1:A:454:LEU:O	1:A:458:MET:HG2	2.08	0.52
1:A:458:MET:O	1:A:462:THR:HG23	2.09	0.52
1:A:761:PHE:HE2	1:A:763:LEU:HD23	1.74	0.52
1:A:585:LYS:O	1:A:618:PHE:HB2	2.10	0.52
1:A:410:MET:HE1	1:A:438:ALA:HA	1.91	0.52
1:B:202:TRP:HB3	1:B:215:LYS:HG2	1.91	0.52
1:A:536:LEU:CD1	1:A:734:GLU:HG3	2.40	0.52
1:A:740:VAL:O	1:A:744:VAL:HG23	2.10	0.52
1:B:624:PRO:HG3	1:B:655:GLU:HG2	1.93	0.51
1:B:159:PRO:O	1:B:163:LYS:HD3	2.11	0.51
1:B:411:GLU:HG3	1:B:415:ARG:HD2	1.91	0.51
1:A:620:GLU:HG2	1:A:634:LYS:HE3	1.93	0.51
1:A:524:VAL:HG11	1:A:726:PRO:HD2	1.93	0.51

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:536:LEU:HD23	1:A:791:ILE:HG12	1.92	0.51
1:B:322:LYS:HD2	1:B:329:ASP:HB3	1.91	0.51
1:B:618:PHE:O	3:B:1001:HOH:O	2.20	0.51
1:B:411:GLU:OE1	1:B:415:ARG:NH1	2.43	0.51
1:A:182:GLN:OE1	1:A:193:GLN:HA	2.10	0.51
1:A:539:VAL:CG1	1:A:781:LEU:HB3	2.41	0.51
1:B:600:ILE:HD12	1:B:723:SER:HB3	1.92	0.50
1:B:661:GLN:HG3	1:B:684:ASN:OD1	2.10	0.50
1:A:669:GLU:O	1:A:670:LYS:HG2	2.10	0.50
1:B:655:GLU:HB3	1:B:658:TYR:HB2	1.92	0.50
1:B:528:TYR:CE2	1:B:812:MET:HB3	2.46	0.50
1:B:189:TYR:CE1	1:B:208:GLY:HA3	2.47	0.50
1:B:569:LEU:HD11	1:B:702:PHE:HB3	1.94	0.50
1:B:83:SER:O	1:B:88:LEU:HD23	2.12	0.50
1:B:159:PRO:HG2	1:B:187:GLY:HA3	1.92	0.50
1:A:606:ARG:HH12	1:A:614:ILE:H	1.58	0.50
1:B:739:ASP:OD1	3:B:1005:HOH:O	2.20	0.49
1:A:103:MET:HE3	1:A:106:HIS:HB3	1.94	0.49
1:B:249:PHE:HD1	1:B:368:LEU:HG	1.77	0.49
1:A:570:VAL:HG13	1:A:703:GLU:HG3	1.95	0.49
1:B:565:ARG:HD2	1:B:678:HIS:HD2	1.76	0.49
1:B:637:ILE:HG13	1:B:638:GLY:H	1.77	0.49
1:A:480:LYS:O	1:A:484:MET:HG3	2.13	0.48
1:A:646:ASN:HA	1:A:648:LYS:NZ	2.28	0.48
1:B:621:SER:HB2	1:B:632:LEU:HA	1.94	0.48
1:A:581:ASN:OD1	1:A:585:LYS:NZ	2.41	0.48
1:B:84:SER:HB3	1:B:232:THR:N	2.17	0.48
1:B:401:LYS:NZ	3:B:1009:HOH:O	2.28	0.48
1:B:655:GLU:HA	3:B:1003:HOH:O	2.13	0.48
1:B:679:THR:O	1:B:683:GLU:HG3	2.12	0.48
1:A:32:ILE:HG21	1:A:408:PHE:HB2	1.95	0.48
1:B:578:ASN:H	1:B:581:ASN:HB3	1.77	0.48
1:A:207:ILE:HA	1:A:832:LEU:HG	1.96	0.48
1:A:627:ASP:CG	1:A:628:THR:H	2.21	0.48
1:B:713:PRO:HG2	1:B:716:ASN:OD1	2.14	0.48
1:B:593:GLY:N	3:B:1019:HOH:O	2.47	0.48
1:B:220:LYS:HD2	1:B:221:ILE:H	1.78	0.47
1:B:481:LEU:HD22	1:B:768:LYS:O	2.14	0.47
1:A:792:MET:HG2	1:A:809:ARG:HH12	1.79	0.47
1:A:690:PRO:HA	1:A:695:LYS:HG2	1.96	0.47
1:A:787:GLN:HB2	1:A:789:LYS:HE3	1.95	0.47

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:481:LEU:HD23	1:B:484:MET:HE3	1.95	0.47
1:B:527:MET:HG3	1:B:795:TYR:CZ	2.49	0.47
1:B:608:GLN:C	1:B:609:ASN:HD22	2.21	0.47
1:A:545:LYS:HD3	1:A:779:GLN:NE2	2.29	0.47
1:B:163:LYS:O	1:B:166:GLN:HG2	2.14	0.47
1:B:367:LEU:HG	1:B:377:LEU:HD12	1.95	0.47
1:A:570:VAL:CG1	1:A:703:GLU:HG3	2.45	0.47
1:B:459:LYS:NZ	1:B:514:LYS:O	2.46	0.47
1:A:259:HIS:CD2	1:A:304:ILE:HG12	2.50	0.47
1:B:651:LEU:HD12	1:B:658:TYR:HE1	1.80	0.47
1:A:691:ALA:O	1:A:693:TRP:N	2.48	0.47
1:B:459:LYS:HG3	1:B:464:ASP:HB2	1.97	0.47
1:B:613:ASP:CG	1:B:614:ILE:H	2.22	0.47
1:A:158:LEU:HD13	1:A:158:LEU:C	2.40	0.47
1:A:542:TYR:CE2	1:A:544:ASN:HB3	2.50	0.47
1:B:453:LYS:HE2	1:B:453:LYS:H	1.80	0.46
1:A:571:LYS:HG3	1:A:667:LEU:CD1	2.46	0.46
1:B:88:LEU:HD12	1:B:89:GLN:H	1.80	0.46
1:B:646:ASN:C	1:B:648:LYS:H	2.22	0.46
1:A:194:PHE:CZ	1:A:196:GLY:HA2	2.50	0.46
1:A:536:LEU:HD11	1:A:734:GLU:HG3	1.98	0.46
1:A:585:LYS:HA	1:A:585:LYS:HD3	1.71	0.46
1:B:224:THR:HG21	1:B:536:LEU:HD21	1.98	0.46
1:B:480:LYS:HE2	1:B:484:MET:HE2	1.98	0.46
1:B:663:ALA:O	1:B:666:CYS:HB3	2.15	0.46
1:A:158:LEU:HD22	1:A:159:PRO:HD2	1.97	0.46
1:A:411:GLU:OE1	1:A:415:ARG:NH1	2.49	0.46
1:B:406:ASN:OD1	1:B:439:LEU:HB2	2.15	0.46
1:A:761:PHE:CE2	1:A:763:LEU:HD23	2.52	0.45
1:B:371:ASP:N	1:B:371:ASP:OD1	2.49	0.45
1:B:453:LYS:H	1:B:453:LYS:CE	2.28	0.45
1:B:676:VAL:CG2	1:B:680:VAL:HG21	2.47	0.45
1:B:207:ILE:CD1	1:B:230:LEU:HD13	2.46	0.45
1:B:581:ASN:OD1	1:B:581:ASN:O	2.34	0.45
1:B:104:ILE:O	1:B:106:HIS:N	2.50	0.45
1:B:657:TYR:CE2	1:B:666:CYS:HA	2.52	0.45
1:B:627:ASP:HB3	1:B:630:SER:HB2	1.99	0.45
1:A:610:ASP:OD1	1:A:610:ASP:C	2.60	0.45
1:A:760:MET:HA	1:A:760:MET:HE2	1.99	0.45
1:A:667:LEU:HD12	1:A:702:PHE:CE2	2.51	0.45
1:B:256:ARG:HG2	1:B:301:VAL:HG22	1.99	0.45

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:445:ASP:N	1:B:445:ASP:OD1	2.49	0.45
1:B:572:LYS:HD3	1:B:703:GLU:HG3	1.99	0.45
1:A:112:ASP:OD1	1:A:116:ASN:N	2.50	0.45
1:A:155:ARG:HD3	1:A:157:GLU:OE1	2.16	0.45
1:A:283:TRP:CD1	1:A:285:GLY:H	2.34	0.45
1:A:794:ASP:OD1	2:A:901:A1CTG:N01	2.50	0.45
1:A:646:ASN:HA	1:A:648:LYS:HZ1	1.82	0.44
1:A:682:PHE:CD2	1:A:682:PHE:N	2.85	0.44
1:A:772:LEU:O	1:A:774:PHE:N	2.45	0.44
1:B:427:ASP:H	1:B:432:ASP:CG	2.25	0.44
1:B:665:ARG:HH21	1:B:693:TRP:HB3	1.82	0.44
1:B:762:GLN:OE1	1:B:762:GLN:HA	2.17	0.44
1:B:77:ILE:HG13	1:B:78:ILE:HG12	1.99	0.44
1:B:313:CYS:SG	3:B:1015:HOH:O	2.62	0.44
1:B:332:TYR:HE1	1:B:340:ARG:HG2	1.81	0.44
1:B:542:TYR:CZ	1:B:544:ASN:HB3	2.52	0.44
1:B:676:VAL:HG13	1:B:677:PRO:O	2.17	0.44
1:B:88:LEU:HD12	1:B:89:GLN:N	2.32	0.44
1:B:428:LEU:O	1:B:430:PHE:N	2.50	0.44
1:A:456:HIS:NE2	1:A:513:LEU:O	2.47	0.44
1:A:733:ARG:O	1:A:737:ILE:HG12	2.17	0.44
1:A:759:ASP:OD2	1:A:760:MET:HG2	2.18	0.44
1:A:786:ARG:NH1	1:A:787:GLN:O	2.51	0.44
1:B:481:LEU:HD23	1:B:484:MET:CE	2.48	0.44
1:B:609:ASN:OD1	1:B:617:PHE:HE1	2.00	0.44
1:B:641:LYS:HE2	1:B:641:LYS:HB2	1.86	0.44
1:A:637:ILE:HG12	1:A:657:TYR:CE2	2.53	0.43
1:A:785:ASP:O	2:A:901:A1CTG:N01	2.51	0.43
1:B:454:LEU:O	1:B:458:MET:HG2	2.17	0.43
1:B:604:LEU:O	1:B:608:GLN:HG3	2.18	0.43
1:A:291:ILE:HD12	1:A:291:ILE:HA	1.81	0.43
1:B:384:ASN:OD1	1:B:384:ASN:C	2.61	0.43
1:B:651:LEU:HD12	1:B:658:TYR:CE1	2.54	0.43
1:A:133:ASN:ND2	1:A:830:GLU:HB3	2.33	0.43
1:B:443:GLU:OE2	1:B:510:LYS:NZ	2.51	0.43
1:A:50:LYS:HB3	1:A:50:LYS:HE3	1.85	0.43
1:A:467:LEU:HD23	1:A:467:LEU:H	1.83	0.43
1:A:570:VAL:HG23	1:A:672:ASP:O	2.19	0.43
1:A:572:LYS:HA	1:A:703:GLU:CD	2.43	0.43
1:A:240:MET:HE2	1:A:342:LEU:HD23	2.01	0.43
1:B:413:GLN:OE1	1:B:436:GLN:HA	2.18	0.43

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:572:LYS:HA	1:B:703:GLU:OE1	2.19	0.43
1:B:575:LYS:HB3	1:B:577:ILE:HG22	2.01	0.43
1:B:629:LYS:HA	1:B:629:LYS:HD2	1.79	0.43
1:B:693:TRP:CE3	1:B:694:ALA:HB3	2.54	0.43
1:A:78:ILE:HB	1:A:393:VAL:HB	2.00	0.43
1:A:88:LEU:HB3	1:A:89:GLN:OE1	2.19	0.42
1:A:589:HIS:CE1	1:A:675:PHE:CD2	3.07	0.42
1:A:611:ASN:OD1	1:A:613:ASP:HB3	2.18	0.42
1:B:615:ASP:OD1	1:B:631:ASN:HB2	2.19	0.42
1:A:470:LYS:O	1:A:742:ARG:NH2	2.53	0.42
1:B:532:MET:SD	1:B:809:ARG:HB3	2.60	0.42
1:B:50:LYS:HE2	1:B:71:PRO:O	2.18	0.42
1:B:129:CYS:N	1:B:150:ARG:HH12	2.18	0.42
1:B:638:GLY:HA2	3:B:1003:HOH:O	2.18	0.42
1:A:250:ASN:HA	1:A:280:LEU:HD21	2.01	0.42
1:A:565:ARG:HB2	1:A:678:HIS:CD2	2.51	0.42
1:A:698:LYS:O	1:A:701:ASP:HB2	2.19	0.42
1:A:8:ARG:HE	1:A:8:ARG:HB2	1.69	0.42
1:B:346:MET:HE3	1:B:346:MET:HA	2.01	0.42
1:B:655:GLU:OE2	1:B:657:TYR:HB2	2.19	0.42
1:A:620:GLU:O	1:A:635:LEU:HD22	2.19	0.42
1:A:620:GLU:HA	1:A:634:LYS:HG3	2.02	0.42
1:B:139:ILE:O	1:B:142:THR:OG1	2.35	0.42
1:A:93:LYS:O	1:A:97:GLN:HG3	2.20	0.42
1:A:221:ILE:H	1:A:221:ILE:HG13	1.64	0.42
1:A:571:LYS:HB2	1:A:574:ASN:CG	2.45	0.42
1:A:617:PHE:CD2	1:A:618:PHE:CD2	3.08	0.42
1:B:742:ARG:NH1	3:B:1005:HOH:O	2.53	0.42
1:A:664:PHE:CD1	1:A:693:TRP:HZ3	2.38	0.41
1:A:695:LYS:HB3	1:A:695:LYS:HE2	1.79	0.41
1:B:6:GLN:O	1:B:401:LYS:HE2	2.20	0.41
1:B:456:HIS:CE1	1:B:514:LYS:HA	2.55	0.41
1:B:533:CYS:SG	1:B:816:LEU:CD1	3.08	0.41
1:A:76:PRO:HA	1:A:394:VAL:HG12	2.02	0.41
1:B:252:LEU:HD12	1:B:252:LEU:HA	1.84	0.41
1:B:692:VAL:HA	1:B:695:LYS:HE2	2.02	0.41
1:B:771:ASN:H	1:B:776:ASP:CG	2.28	0.41
1:A:565:ARG:HG2	1:A:721:LYS:HD3	2.03	0.41
1:A:406:ASN:O	1:A:410:MET:HG2	2.21	0.41
1:A:537:PRO:HD2	1:A:791:ILE:HD11	2.02	0.41
1:A:595:ILE:HG22	1:A:596:ALA:N	2.35	0.41

*Continued on next page...*



Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:205:ASN:HD21	1:A:209:GLU:HB2	1.86	0.41
1:B:41:SER:HB2	1:B:326:ASN:HB2	2.02	0.41
1:B:598:TRP:C	1:B:601:PRO:HD2	2.45	0.41
1:A:190:GLU:OE2	3:A:1002:HOH:O	2.22	0.41
1:A:361:HIS:HB3	1:A:365:TYR:CD2	2.55	0.41
1:B:184:ASP:OD2	1:B:187:GLY:N	2.54	0.41
1:A:99:LEU:HD23	1:A:108:ILE:HD13	2.03	0.41
1:A:195:HIS:HB2	1:A:202:TRP:CH2	2.56	0.41
1:B:249:PHE:N	1:B:368:LEU:HD11	2.35	0.41
1:B:396:ARG:HH21	1:B:400:ASP:CG	2.28	0.41
1:A:481:LEU:HD23	1:A:484:MET:HE2	2.02	0.41
1:B:571:LYS:NZ	1:B:668:VAL:O	2.54	0.40
1:A:286:PRO:HD2	1:A:288:LYS:HZ2	1.86	0.40
1:B:336:TYR:CZ	1:B:426:LYS:HD2	2.56	0.40
1:B:637:ILE:HD12	1:B:637:ILE:HA	1.91	0.40
1:A:780:CYS:SG	1:A:782:ILE:HD11	2.62	0.40
1:B:364:VAL:HG13	1:B:365:TYR:CD2	2.56	0.40
1:B:406:ASN:CG	1:B:439:LEU:HD22	2.46	0.40
1:A:657:TYR:CE1	1:A:666:CYS:HA	2.56	0.40
1:B:732:THR:HG23	1:B:737:ILE:HD13	2.04	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

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	814/853 (95%)	773 (95%)	36 (4%)	5 (1%)	21	27
1	B	814/853 (95%)	760 (93%)	47 (6%)	7 (1%)	14	17
All	All	1628/1706 (95%)	1533 (94%)	83 (5%)	12 (1%)	18	24

All (12) Ramachandran outliers are listed below:



Mol	Chain	Res	Type
1	A	574	ASN
1	A	692	VAL
1	B	133	ASN
1	B	614	ILE
1	B	694	ALA
1	A	598	TRP
1	B	268	TRP
1	A	617	PHE
1	B	429	MET
1	B	828	SER
1	A	557	PRO
1	B	637	ILE

### 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	703/730 (96%)	697 (99%)	6 (1%)	70	81
1	B	703/730 (96%)	691 (98%)	12 (2%)	53	67
All	All	1406/1460 (96%)	1388 (99%)	18 (1%)	61	73

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	104	ILE
1	A	221	ILE
1	A	720	CYS
1	A	759	ASP
1	A	827	ASN
1	A	832	LEU
1	B	92	LEU
1	B	102	LYS
1	B	104	ILE
1	B	132	VAL
1	B	266	ASP
1	B	284	ASP

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	B	377	LEU
1	B	418	LYS
1	B	637	ILE
1	B	676	VAL
1	B	707	LEU
1	B	827	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	148	GLN
1	A	192	GLN
1	A	248	GLN
1	A	466	HIS
1	A	589	HIS
1	A	609	ASN
1	A	631	ASN
1	A	678	HIS
1	A	684	ASN
1	A	689	ASN
1	A	716	ASN
1	A	767	ASN
1	A	779	GLN
1	A	824	HIS
1	B	195	HIS
1	B	309	GLN
1	B	315	GLN
1	B	361	HIS
1	B	466	HIS
1	B	578	ASN
1	B	608	GLN
1	B	609	ASN
1	B	678	HIS
1	B	746	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

## 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

2 ligands are modelled in this entry.

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	$\# Z  > 2$
2	A1CTG	A	901	-	23,29,29	0.63	1 (4%)	22,45,45	1.01	3 (13%)
2	A1CTG	B	901	-	23,29,29	0.52	0	22,45,45	0.89	1 (4%)

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
2	A1CTG	A	901	-	-	2/9/57/57	0/4/4/4
2	A1CTG	B	901	-	-	0/9/57/57	0/4/4/4

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	901	A1CTG	C08-N07	-2.01	1.32	1.34

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	901	A1CTG	N03-C04-N09	-2.66	107.84	112.45

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	901	A1CTG	C10-N09-C08	2.50	126.20	122.11
2	A	901	A1CTG	N03-C02-N26	2.07	111.31	109.21
2	B	901	A1CTG	N03-C02-N26	2.05	111.29	109.21

There are no chirality outliers.

All (2) torsion outliers are listed below:

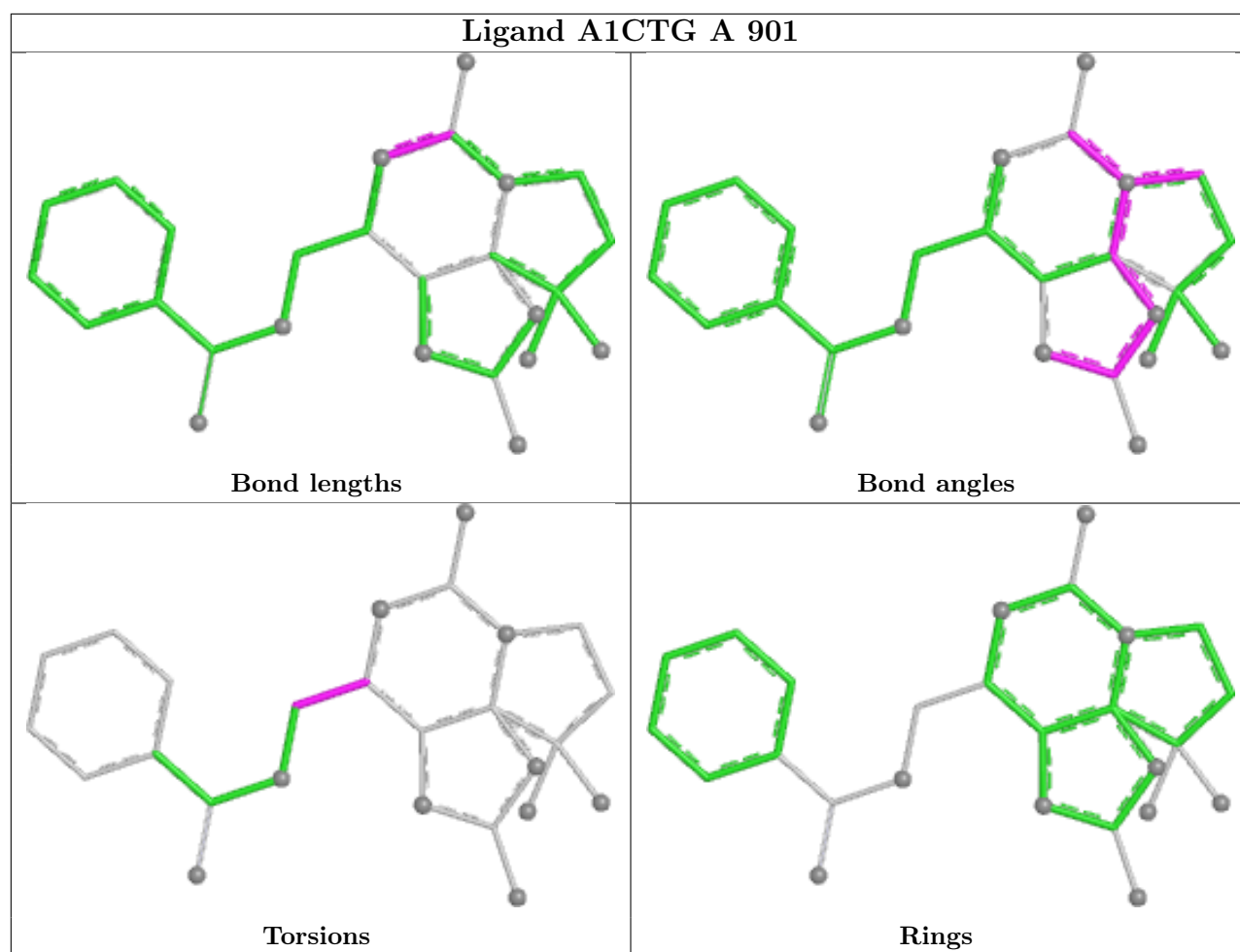
Mol	Chain	Res	Type	Atoms
2	A	901	A1CTG	C05-C06-C16-O17
2	A	901	A1CTG	N07-C06-C16-O17

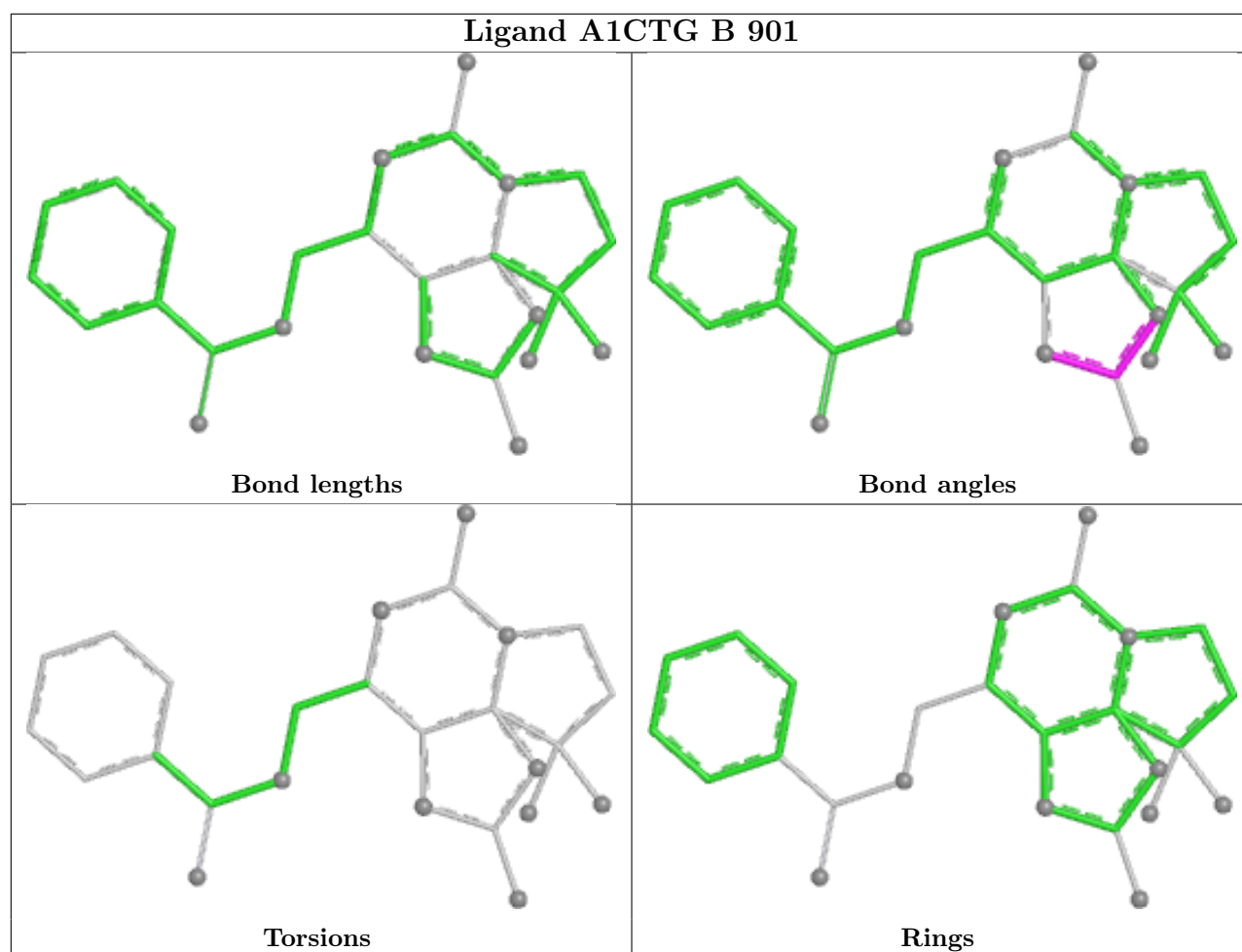
There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	901	A1CTG	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

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	818/853 (95%)	0.69	78 (9%) 14 11	51, 85, 166, 201	0
1	B	818/853 (95%)	0.76	64 (7%) 19 16	56, 90, 151, 196	0
All	All	1636/1706 (95%)	0.73	142 (8%) 16 14	51, 89, 162, 201	0

All (142) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	832	LEU	9.9
1	B	832	LEU	7.8
1	B	829	LEU	6.5
1	B	598	TRP	5.5
1	B	207	ILE	4.9
1	B	82	TYR	4.9
1	B	831	VAL	4.6
1	B	86	ARG	4.4
1	A	664	PHE	4.1
1	A	831	VAL	4.1
1	B	168	ALA	4.0
1	A	644	ALA	3.8
1	A	564	LEU	3.7
1	B	179	PHE	3.7
1	A	640	PRO	3.7
1	A	702	PHE	3.6
1	A	725	ILE	3.6
1	A	726	PRO	3.5
1	B	592	VAL	3.5
1	B	83	SER	3.4
1	B	468	PRO	3.3
1	B	614	ILE	3.3
1	A	467	LEU	3.3
1	B	830	GLU	3.3

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	A	577	ILE	3.3
1	B	664	PHE	3.2
1	B	827	ASN	3.2
1	A	756	PHE	3.2
1	A	468	PRO	3.2
1	A	830	GLU	3.2
1	A	604	LEU	3.1
1	A	221	ILE	3.1
1	A	724	GLY	3.1
1	A	637	ILE	3.1
1	B	84	SER	3.0
1	A	796	PHE	3.0
1	A	804	VAL	3.0
1	B	5	LYS	3.0
1	A	558	ALA	3.0
1	A	283	TRP	2.9
1	B	583	LYS	2.9
1	A	608	GLN	2.9
1	B	644	ALA	2.8
1	A	693	TRP	2.8
1	A	646	ASN	2.8
1	B	224	THR	2.8
1	B	88	LEU	2.8
1	A	5	LYS	2.8
1	B	254	GLY	2.7
1	A	579	TRP	2.7
1	B	9	TRP	2.7
1	B	209	GLU	2.7
1	A	554	PRO	2.7
1	A	563	THR	2.7
1	A	179	PHE	2.7
1	B	408	PHE	2.7
1	A	705	LEU	2.7
1	A	194	PHE	2.6
1	A	674	ALA	2.6
1	B	208	GLY	2.6
1	B	206	ALA	2.6
1	B	467	LEU	2.6
1	A	712	ALA	2.5
1	B	717	TYR	2.5
1	B	785	ASP	2.5
1	A	556	SER	2.5

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	A	707	LEU	2.5
1	B	409	LEU	2.5
1	B	722	LEU	2.5
1	A	691	ALA	2.5
1	A	676	VAL	2.5
1	A	681	VAL	2.5
1	A	168	ALA	2.5
1	B	233	THR	2.5
1	B	230	LEU	2.4
1	A	596	ALA	2.4
1	B	573	SER	2.4
1	B	723	SER	2.4
1	B	828	SER	2.4
1	B	287	ALA	2.4
1	B	712	ALA	2.4
1	A	670	LYS	2.4
1	A	675	PHE	2.4
1	A	609	ASN	2.4
1	A	642	ASN	2.4
1	B	106	HIS	2.3
1	A	718	LYS	2.3
1	B	826	SER	2.3
1	B	605	ILE	2.3
1	A	721	LYS	2.3
1	A	557	PRO	2.3
1	A	647	THR	2.3
1	B	136	GLY	2.3
1	B	231	VAL	2.3
1	A	763	LEU	2.3
1	A	648	LYS	2.3
1	A	751	TYR	2.3
1	A	628	THR	2.2
1	A	552	LYS	2.2
1	A	715	SER	2.2
1	B	715	SER	2.2
1	A	551	CYS	2.2
1	A	801	TYR	2.2
1	B	332	TYR	2.2
1	B	373	THR	2.2
1	A	635	LEU	2.2
1	B	92	LEU	2.2
1	B	582	ILE	2.2

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	A	713	PRO	2.2
1	B	430	PHE	2.2
1	B	212	ALA	2.2
1	B	566	ALA	2.2
1	A	706	CYS	2.2
1	A	677	PRO	2.2
1	B	604	LEU	2.1
1	A	717	TYR	2.1
1	A	524	VAL	2.1
1	A	599	VAL	2.1
1	B	602	VAL	2.1
1	A	567	VAL	2.1
1	A	692	VAL	2.1
1	B	364	VAL	2.1
1	B	570	VAL	2.1
1	B	668	VAL	2.1
1	A	805	TYR	2.1
1	A	639	ASP	2.1
1	A	561	PHE	2.1
1	A	162	LEU	2.1
1	A	619	GLY	2.1
1	B	124	GLY	2.1
1	A	803	THR	2.1
1	A	156	HIS	2.1
1	A	299	PHE	2.0
1	A	618	PHE	2.0
1	A	740	VAL	2.0
1	A	569	LEU	2.0
1	B	103	MET	2.0
1	B	260	SER	2.0
1	B	466	HIS	2.0
1	B	568	ALA	2.0
1	B	645	ALA	2.0
1	A	791	ILE	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 [i](#)

There are no oligosaccharides in this entry.

## 6.4 Ligands [i](#)

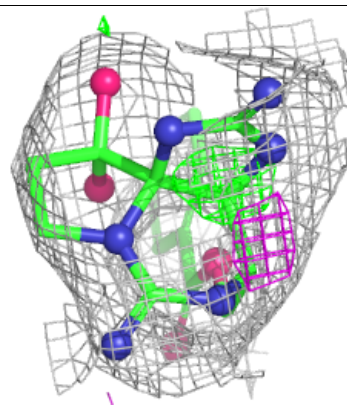
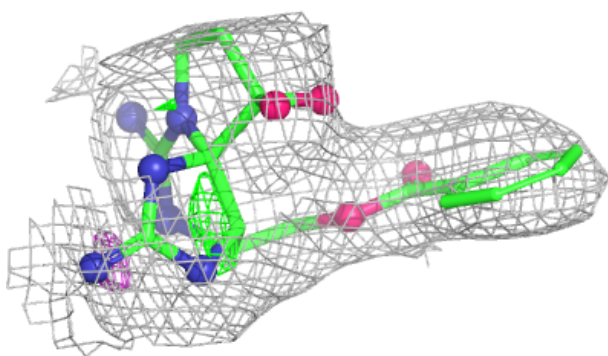
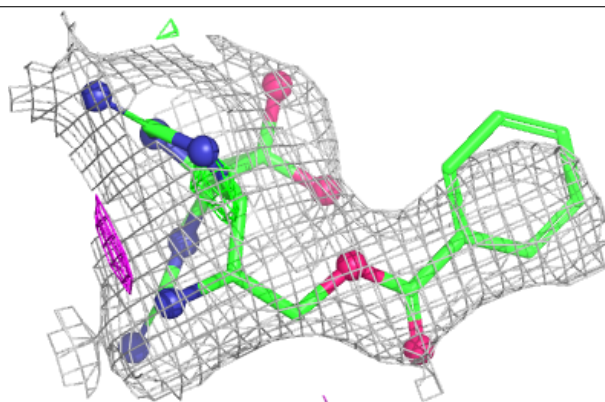
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	A1CTG	A	901	26/26	0.88	0.14	112,123,136,149	0
2	A1CTG	B	901	26/26	0.91	0.12	78,85,95,100	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

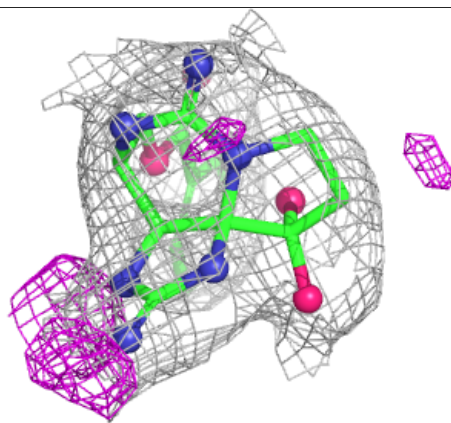
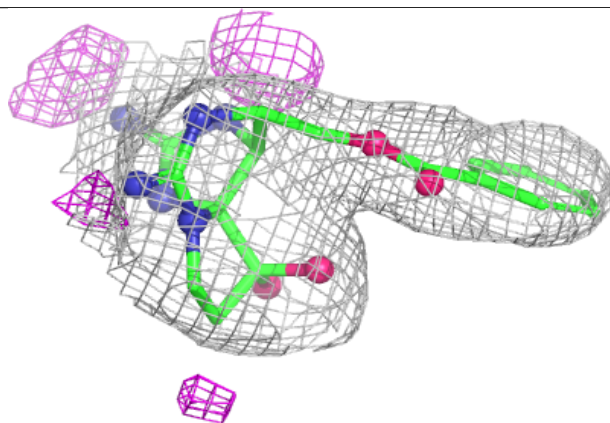
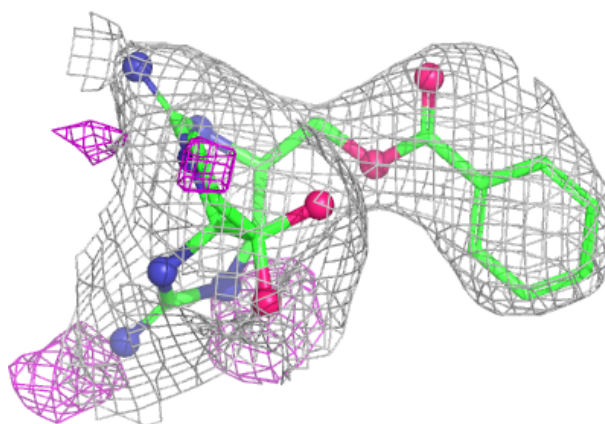
### Electron density around A1CTG A 901:

2mF<sub>o</sub>-DF<sub>c</sub> (at 0.7 rmsd) in gray  
mF<sub>o</sub>-DF<sub>c</sub> (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around A1CTG B 901:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



## 6.5 Other polymers [i](#)

There are no such residues in this entry.