



# Full wwPDB X-ray Structure Validation Report

May 22, 2020 – 04:25 am BST

PDB ID : 2G27  
Title : Ketopiperazine-Based Renin Inhibitors: Optimization of the "C" Ring  
Authors : Holsworth, D.D.; Jalaiea, M.; Zhanga, E.; Mcconnella, P.  
Deposited on : 2006-02-15  
Resolution : 2.90 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.

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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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

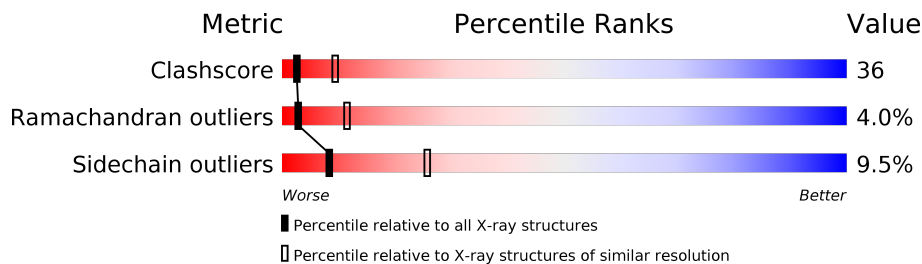
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.90 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))
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)

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 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  $\leq 5\%$ .

Mol	Chain	Length	Quality of chain
1	A	333	
1	B	333	

## 2 Entry composition [i](#)

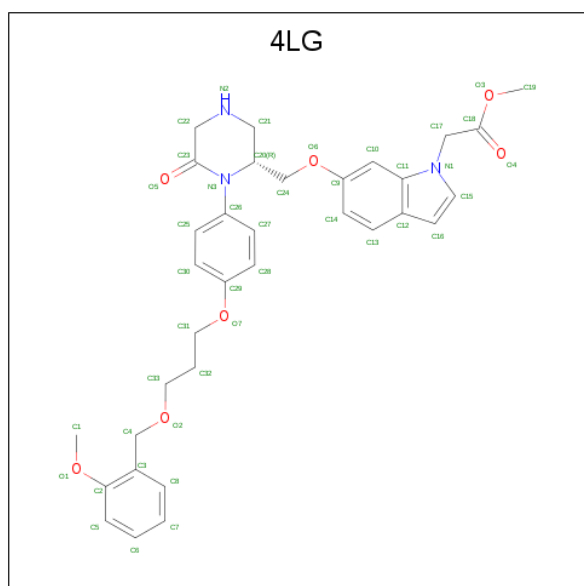
There are 2 unique types of molecules in this entry. The entry contains 5154 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Renin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	328	Total 2534	C 1622	N 409	O 489	S 14	0	0	0
1	B	328	Total 2534	C 1622	N 409	O 489	S 14	0	0	0

- Molecule 2 is METHYL (6-{{(2R)-1-(4-{{3-{{(2-METHOXYBENZYL)OXY}PROPOXY}PHENYL)-6-OXOPIPERAZIN-2-YL}METHOXY}-1H-INDOL-1-YL)ACETATE (three-letter code: 4LG) (formula: C<sub>33</sub>H<sub>37</sub>N<sub>3</sub>O<sub>7</sub>).

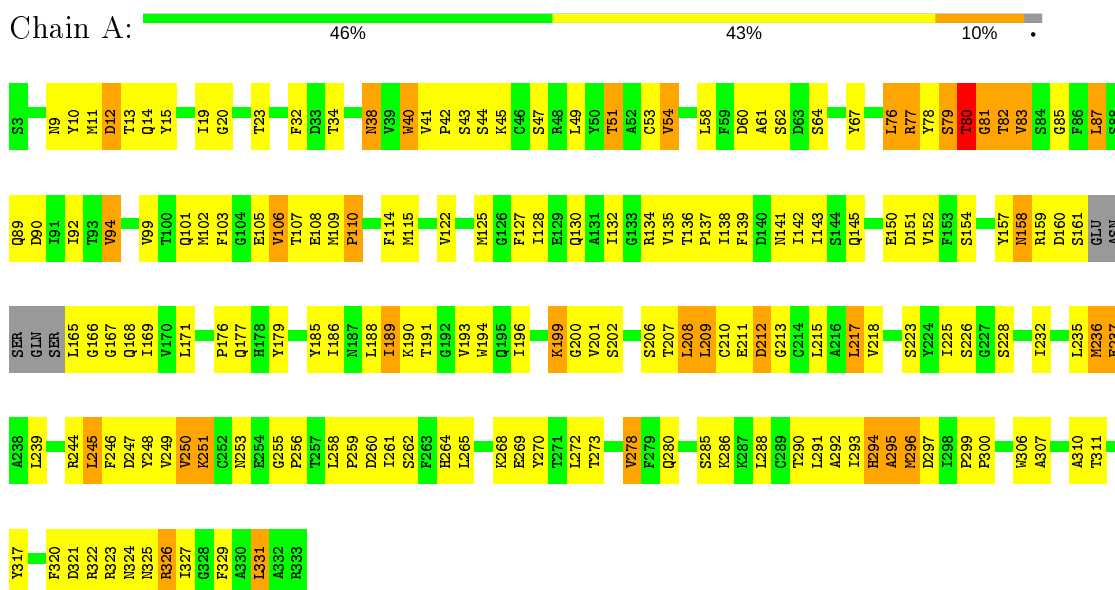


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	Total 43	C 33	N 3	O 7	0	0
2	B	1	Total 43	C 33	N 3	O 7	0	0

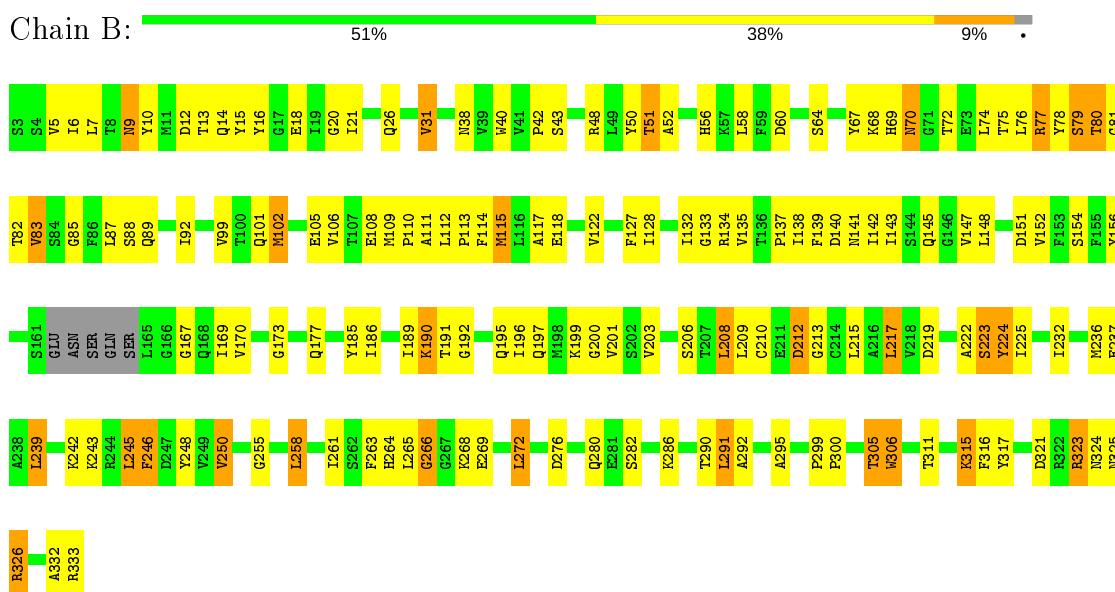
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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. 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. 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: Renin



- Molecule 1: Renin



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	141.41Å 141.41Å 141.41Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.90 47.14 – 2.90	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-2.90) 91.1 (47.14-2.90)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.95 (at 2.91Å)	Xtrriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.230 , 0.310 0.246 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	51.4	Xtrriage
Anisotropy	0.000	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 29.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	0.044 for l,-k,h	Xtrriage
$F_o, F_c$ correlation	0.80	EDS
Total number of atoms	5154	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	45.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.24% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 4LG

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	0.39	0/2593	0.72	1/3514 (0.0%)
1	B	0.44	0/2593	0.75	1/3514 (0.0%)
All	All	0.42	0/5186	0.73	2/7028 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	306	TRP	N-CA-C	-6.36	93.82	111.00
1	A	80	THR	N-CA-C	-5.96	94.91	111.00

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-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 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	2534	0	2470	190	0
1	B	2534	0	2470	178	0
2	A	43	0	37	9	0
2	B	43	0	37	12	0
All	All	5154	0	5014	371	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 36.

All (371) 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:82:THR:HG23	1:A:83:VAL:H	1.02	1.10
1:B:152:VAL:HG12	1:B:321:ASP:HA	1.30	1.09
1:A:43:SER:HB2	1:A:105:GLU:HB3	1.35	1.06
1:A:158:ASN:ND2	1:A:159:ARG:H	1.53	1.05
1:A:158:ASN:HD22	1:A:159:ARG:N	1.56	1.04
1:A:77:ARG:H	1:A:77:ARG:HD2	1.20	1.04
1:B:323:ARG:HH11	1:B:323:ARG:HG3	1.34	0.93
1:A:82:THR:HG23	1:A:83:VAL:N	1.84	0.93
1:B:324:ASN:HB2	1:B:326:ARG:HD2	1.54	0.90
1:A:82:THR:CG2	1:A:83:VAL:H	1.86	0.89
1:B:43:SER:HB2	1:B:105:GLU:HB3	1.54	0.88
1:A:128:ILE:HD13	1:A:137:PRO:HD3	1.56	0.87
1:B:245:LEU:HD23	1:B:246:PHE:N	1.90	0.87
1:B:14:GLN:HA	2:B:803:4LG:O4	1.74	0.87
1:A:152:VAL:HG12	1:A:321:ASP:HA	1.58	0.85
1:B:315:LYS:HG2	1:B:316:PHE:CE1	2.13	0.83
1:A:130:GLN:NE2	1:A:193:VAL:HG22	1.93	0.83
1:B:76:LEU:HD11	1:B:132:ILE:HG13	1.61	0.82
1:B:14:GLN:HG2	2:B:803:4LG:H15	1.61	0.81
1:A:78:TYR:H	1:A:81:GLY:HA2	1.45	0.81
1:A:51:THR:HB	1:A:115:MET:HG3	1.63	0.80
1:B:217:LEU:HD22	1:B:219:ASP:HB2	1.63	0.80
1:A:158:ASN:HD22	1:A:159:ARG:H	0.83	0.80
1:A:19:ILE:HG22	1:A:94:VAL:HG13	1.62	0.80
1:B:245:LEU:HD23	1:B:246:PHE:H	1.46	0.79
1:A:77:ARG:N	1:A:77:ARG:HD2	1.97	0.79
1:A:82:THR:HG23	1:A:110:PRO:HD3	1.66	0.78
1:A:249:VAL:HG21	1:A:288:LEU:HD13	1.66	0.78
1:B:101:GLN:HA	1:B:101:GLN:HE21	1.51	0.76
1:B:72:THR:HB	1:B:87:LEU:HD12	1.67	0.75
1:B:203:VAL:HB	1:B:208:LEU:HD11	1.67	0.75
1:A:158:ASN:ND2	1:A:159:ARG:N	2.24	0.75
1:B:258:LEU:N	1:B:258:LEU:HD23	2.01	0.74
1:B:128:ILE:HG13	1:B:137:PRO:HD3	1.68	0.73
1:A:130:GLN:HE22	1:A:193:VAL:HG22	1.53	0.73
1:A:11:MET:O	1:A:13:THR:HG23	1.87	0.73
1:A:132:ILE:O	1:A:135:VAL:HG23	1.90	0.72
1:B:87:LEU:HD21	1:B:135:VAL:HG21	1.71	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:9:ASN:HD21	1:B:12:ASP:H	1.38	0.71
1:A:268:LYS:NZ	1:A:269:GLU:H	1.89	0.71
1:A:9:ASN:OD1	1:A:160:ASP:HB2	1.90	0.70
1:A:40:TRP:HB3	2:A:804:4LG:H331	1.73	0.70
1:A:211:GLU:O	1:A:212:ASP:HB2	1.92	0.70
1:A:87:LEU:HD13	1:A:102:MET:HE2	1.73	0.70
1:A:235:LEU:O	1:A:239:LEU:HG	1.92	0.69
1:B:321:ASP:HB3	1:B:326:ARG:HG2	1.75	0.69
1:B:79:SER:O	1:B:80:THR:HG22	1.93	0.69
1:B:268:LYS:HG3	1:B:269:GLU:N	2.07	0.69
1:A:47:SER:OG	1:A:49:LEU:HD23	1.94	0.68
1:A:208:LEU:O	1:A:209:LEU:HG	1.94	0.68
1:A:45:LYS:HD3	1:A:45:LYS:O	1.94	0.67
1:A:150:GLU:HB2	1:A:152:VAL:HG22	1.74	0.67
1:A:87:LEU:HD13	1:A:102:MET:CE	2.25	0.67
1:B:51:THR:HB	1:B:115:MET:SD	2.35	0.67
1:B:170:VAL:HB	1:B:173:GLY:O	1.95	0.67
1:A:199:LYS:HB3	1:A:264:HIS:HD2	1.58	0.66
1:B:139:PHE:O	1:B:143:ILE:HG12	1.95	0.66
1:A:76:LEU:HD13	1:A:76:LEU:O	1.96	0.66
1:B:82:THR:HG23	1:B:110:PRO:HG3	1.77	0.66
1:B:89:GLN:HB2	1:B:102:MET:HE1	1.78	0.66
1:B:201:VAL:HG23	1:B:209:LEU:HB2	1.78	0.66
1:A:76:LEU:HB3	1:A:132:ILE:HD11	1.77	0.65
1:B:245:LEU:HD23	1:B:246:PHE:CB	2.26	0.65
1:A:249:VAL:HG12	1:A:290:THR:HA	1.78	0.65
1:B:208:LEU:HD13	1:B:208:LEU:H	1.61	0.65
1:B:50:TYR:CZ	1:B:108:GLU:HB2	2.32	0.65
1:B:127:PHE:CB	1:B:192:GLY:HA2	2.28	0.64
1:B:224:TYR:HB3	1:B:292:ALA:O	1.97	0.64
1:B:74:LEU:HD21	1:B:87:LEU:HG	1.78	0.64
1:B:154:SER:HB3	1:B:317:TYR:CE1	2.33	0.64
1:A:82:THR:CG2	1:A:110:PRO:HD3	2.28	0.64
1:B:83:VAL:HA	1:B:108:GLU:O	1.98	0.64
1:A:101:GLN:HG3	1:A:102:MET:N	2.14	0.63
1:A:43:SER:O	1:A:58:LEU:HD13	1.96	0.63
1:A:13:THR:O	2:A:804:4LG:O4	2.16	0.63
1:A:87:LEU:HD23	1:A:103:PHE:O	1.99	0.63
1:B:76:LEU:CD1	1:B:132:ILE:HG13	2.27	0.63
1:A:78:TYR:O	1:A:81:GLY:N	2.32	0.63
1:B:323:ARG:NH1	1:B:323:ARG:HG3	2.09	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:102:MET:HB3	1:B:135:VAL:CG1	2.29	0.62
1:A:199:LYS:HB3	1:A:264:HIS:CD2	2.35	0.62
1:A:268:LYS:HZ2	1:A:269:GLU:H	1.46	0.62
1:A:128:ILE:CD1	1:A:137:PRO:HD3	2.29	0.62
1:A:185:TYR:C	1:A:186:ILE:HD12	2.20	0.62
1:B:127:PHE:HB2	1:B:192:GLY:HA2	1.82	0.62
1:B:26:GLN:HE22	1:B:60:ASP:H	1.47	0.62
1:A:223:SER:O	1:A:310:ALA:HB3	2.00	0.61
1:B:111:ALA:O	1:B:115:MET:HB2	2.00	0.61
1:A:106:VAL:HG11	1:A:109:MET:HG3	1.81	0.61
1:A:41:VAL:O	1:A:106:VAL:HG23	2.00	0.61
1:B:199:LYS:NZ	1:B:266:GLY:HA2	2.15	0.61
1:A:43:SER:HB2	1:A:105:GLU:CB	2.23	0.61
1:B:315:LYS:HG2	1:B:316:PHE:CD1	2.34	0.61
1:A:99:VAL:HG11	1:A:142:ILE:HG12	1.81	0.61
1:A:38:ASN:HD21	1:A:132:ILE:H	1.48	0.61
1:B:152:VAL:CG1	1:B:321:ASP:HA	2.19	0.61
1:A:20:GLY:O	1:A:92:ILE:HA	2.01	0.61
1:B:236:MET:HG3	1:B:248:TYR:CD2	2.36	0.61
1:B:210:CYS:SG	1:B:210:CYS:O	2.58	0.60
1:A:253:ASN:CG	1:A:285:SER:HA	2.22	0.60
1:B:89:GLN:HB2	1:B:102:MET:CE	2.31	0.60
1:B:99:VAL:HG21	1:B:141:ASN:HB3	1.84	0.60
1:A:14:GLN:HA	2:A:804:4LG:O4	2.01	0.60
1:A:14:GLN:HG2	2:A:804:4LG:H15	1.84	0.60
1:B:9:ASN:HD21	1:B:12:ASP:N	2.00	0.60
1:A:245:LEU:HD22	1:A:246:PHE:CE1	2.37	0.59
1:A:154:SER:HB3	1:A:317:TYR:CE1	2.37	0.59
1:A:77:ARG:CD	1:A:77:ARG:H	2.02	0.59
1:B:189:ILE:HD11	1:B:197:GLN:HB2	1.85	0.59
1:A:196:ILE:CD1	1:A:265:LEU:HD22	2.33	0.58
1:A:211:GLU:O	1:A:212:ASP:CB	2.50	0.58
1:A:209:LEU:HD11	1:A:235:LEU:HB2	1.83	0.58
1:B:199:LYS:O	1:B:264:HIS:HB3	2.04	0.58
1:A:76:LEU:O	1:A:83:VAL:HG13	2.03	0.58
1:B:215:LEU:CB	1:B:305:THR:HG22	2.34	0.58
1:A:206:SER:HB2	1:A:208:LEU:HG	1.86	0.58
1:A:138:ILE:O	1:A:142:ILE:HG13	2.04	0.58
1:A:82:THR:O	1:A:83:VAL:O	2.23	0.57
1:A:253:ASN:OD1	1:A:285:SER:HA	2.04	0.57
1:A:226:SER:HB2	1:A:296:MET:HB3	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:190:LYS:HG3	1:B:191:THR:N	2.20	0.57
1:A:81:GLY:O	1:A:82:THR:HG22	2.04	0.57
1:B:239:LEU:HD21	1:B:261:ILE:HD11	1.85	0.57
1:A:150:GLU:CB	1:A:152:VAL:HG22	2.35	0.56
1:A:76:LEU:HD13	1:A:83:VAL:HG13	1.87	0.56
1:B:323:ARG:HH11	1:B:323:ARG:CG	2.11	0.56
1:A:122:VAL:HG23	2:A:804:4LG:H11	1.87	0.56
1:B:143:ILE:CD1	1:B:148:LEU:HD12	2.35	0.56
1:B:78:TYR:O	1:B:79:SER:C	2.44	0.56
1:A:134:ARG:HH21	1:A:134:ARG:HB2	1.70	0.56
1:A:89:GLN:O	1:A:90:ASP:HB2	2.04	0.56
1:B:106:VAL:HG21	2:B:803:4LG:H331	1.86	0.56
1:B:38:ASN:HB2	1:B:40:TRP:CZ3	2.41	0.55
1:B:191:THR:HG22	1:B:192:GLY:N	2.22	0.55
1:B:10:TYR:HB3	1:B:14:GLN:HB2	1.89	0.55
1:B:13:THR:O	2:B:803:4LG:O4	2.25	0.55
1:A:210:CYS:SG	1:A:210:CYS:O	2.65	0.55
1:A:101:GLN:NE2	1:A:138:ILE:HA	2.22	0.54
1:B:154:SER:HB3	1:B:317:TYR:HE1	1.73	0.54
1:A:165:LEU:HD11	1:A:167:GLY:O	2.07	0.54
1:B:272:LEU:HD23	1:B:276:ASP:HB3	1.90	0.54
1:A:78:TYR:H	1:A:81:GLY:CA	2.18	0.54
1:B:250:VAL:HG21	1:B:258:LEU:HD11	1.88	0.54
1:A:186:ILE:HD12	1:A:186:ILE:N	2.22	0.54
1:B:232:ILE:HB	1:B:295:ALA:HB2	1.90	0.54
1:B:177:GLN:HA	1:B:333:ARG:HH21	1.73	0.54
1:B:101:GLN:HA	1:B:101:GLN:NE2	2.21	0.54
1:B:74:LEU:N	1:B:85:GLY:O	2.39	0.53
1:B:195:GLN:OE1	1:B:215:LEU:HD22	2.08	0.53
1:A:152:VAL:HG12	1:A:321:ASP:CA	2.35	0.53
1:A:235:LEU:C	1:A:235:LEU:HD23	2.29	0.53
1:A:76:LEU:HB3	1:A:132:ILE:CD1	2.38	0.53
1:B:245:LEU:HD23	1:B:246:PHE:CA	2.38	0.53
1:A:106:VAL:CG1	1:A:109:MET:HG3	2.39	0.53
1:B:324:ASN:CB	1:B:326:ARG:HD2	2.33	0.53
1:B:5:VAL:CG1	1:B:169:ILE:HB	2.39	0.53
1:B:245:LEU:CD2	1:B:246:PHE:N	2.67	0.53
1:B:79:SER:O	1:B:80:THR:CG2	2.57	0.53
1:B:79:SER:O	1:B:80:THR:CB	2.57	0.53
1:A:83:VAL:HA	1:A:108:GLU:O	2.09	0.53
1:A:165:LEU:HD12	1:A:166:GLY:N	2.23	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:176:PRO:HA	1:A:179:TYR:CE1	2.44	0.52
1:A:190:LYS:HG3	1:A:190:LYS:O	2.08	0.52
1:A:159:ARG:O	1:A:160:ASP:HB3	2.09	0.52
1:A:165:LEU:HD12	1:A:166:GLY:H	1.73	0.52
1:A:201:VAL:H	1:A:210:CYS:HB3	1.74	0.52
1:B:190:LYS:HG3	1:B:191:THR:O	2.10	0.52
1:B:48:ARG:NH1	1:B:58:LEU:HD13	2.25	0.52
1:B:75:THR:O	1:B:76:LEU:HD13	2.09	0.52
1:A:78:TYR:O	1:A:80:THR:HG22	2.09	0.51
1:B:102:MET:HB3	1:B:135:VAL:HG13	1.91	0.51
1:B:7:LEU:HB2	1:B:167:GLY:C	2.30	0.51
1:A:78:TYR:O	1:A:80:THR:N	2.44	0.51
1:B:291:LEU:N	1:B:291:LEU:HD23	2.25	0.51
1:B:232:ILE:HG21	1:B:295:ALA:HB2	1.93	0.51
1:A:101:GLN:HE22	1:A:138:ILE:HA	1.75	0.51
1:A:82:THR:CG2	1:A:83:VAL:N	2.56	0.51
1:A:291:LEU:C	1:A:293:ILE:H	2.14	0.51
1:B:224:TYR:HA	1:B:311:THR:OG1	2.11	0.51
1:B:9:ASN:ND2	1:B:12:ASP:H	2.04	0.51
1:B:208:LEU:CD1	1:B:208:LEU:H	2.24	0.51
1:A:136:THR:HG21	1:A:141:ASN:ND2	2.26	0.50
1:A:260:ASP:C	1:A:261:ILE:HD13	2.32	0.50
1:A:10:TYR:HB3	1:A:14:GLN:HB2	1.93	0.50
1:B:42:PRO:HB2	1:B:58:LEU:HD23	1.92	0.50
1:B:10:TYR:CD2	1:B:118:GLU:HG3	2.46	0.50
1:A:9:ASN:HD22	1:A:10:TYR:N	2.10	0.50
1:A:278:VAL:HG12	1:A:291:LEU:HD22	1.94	0.50
1:B:20:GLY:O	1:B:92:ILE:HA	2.12	0.50
1:A:82:THR:HG22	1:A:110:PRO:HB3	1.94	0.50
1:A:232:ILE:HG13	1:A:295:ALA:HA	1.93	0.50
1:B:14:GLN:CG	2:B:803:4LG:H15	2.38	0.50
1:B:88:SER:O	1:B:102:MET:HE2	2.12	0.49
1:A:83:VAL:HB	1:A:110:PRO:HD2	1.93	0.49
1:B:203:VAL:HB	1:B:208:LEU:CD1	2.39	0.49
1:A:188:LEU:HG	1:A:325:ASN:O	2.12	0.49
1:B:208:LEU:CD2	1:B:209:LEU:HG	2.42	0.49
1:A:151:ASP:OD2	1:A:322:ARG:HB2	2.13	0.49
1:B:245:LEU:CD2	1:B:246:PHE:H	2.20	0.49
1:B:48:ARG:HH12	1:B:58:LEU:HD13	1.77	0.49
1:A:9:ASN:HD21	1:A:12:ASP:N	2.11	0.49
1:B:21:ILE:HG12	1:B:92:ILE:HG12	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:239:LEU:HD22	1:A:259:PRO:HG2	1.94	0.49
1:B:232:ILE:CB	1:B:295:ALA:HB2	2.42	0.49
1:B:75:THR:C	1:B:76:LEU:HD22	2.33	0.49
1:B:201:VAL:CG2	1:B:209:LEU:HB2	2.43	0.48
2:B:803:4LG:H42	2:B:803:4LG:H13A	1.94	0.48
1:A:128:ILE:HD13	1:A:136:THR:HA	1.96	0.48
1:A:262:SER:HA	1:A:270:TYR:O	2.14	0.48
1:A:237:GLU:C	1:A:239:LEU:H	2.16	0.48
1:A:42:PRO:HA	1:A:106:VAL:HG23	1.96	0.48
1:B:56:HIS:CE1	2:B:803:4LG:H6	2.48	0.48
1:B:76:LEU:HD22	1:B:76:LEU:N	2.28	0.48
1:B:78:TYR:H	1:B:81:GLY:HA3	1.79	0.48
1:B:101:GLN:HE21	1:B:101:GLN:CA	2.16	0.48
1:A:15:TYR:CZ	1:A:157:TYR:HB3	2.48	0.48
1:A:202:SER:HB2	1:A:207:THR:HG23	1.95	0.48
1:A:78:TYR:O	1:A:79:SER:C	2.52	0.48
1:A:81:GLY:O	1:A:82:THR:CB	2.62	0.48
1:B:89:GLN:HE21	1:B:102:MET:CE	2.26	0.48
1:B:83:VAL:H	1:B:110:PRO:HD3	1.76	0.48
1:A:251:LYS:HD2	1:A:286:LYS:O	2.14	0.47
1:B:232:ILE:CG2	1:B:295:ALA:HB2	2.44	0.47
1:A:244:ARG:HB2	1:A:247:ASP:O	2.15	0.47
1:A:196:ILE:HD11	1:A:265:LEU:HD22	1.96	0.47
1:A:299:PRO:HA	1:A:300:PRO:HD3	1.71	0.47
1:B:263:PHE:N	1:B:263:PHE:CD1	2.82	0.47
1:A:225:ILE:HG13	1:A:311:THR:HB	1.97	0.47
1:A:60:ASP:OD1	1:A:62:SER:HB2	2.14	0.47
1:A:23:THR:O	1:A:64:SER:HA	2.14	0.47
1:B:40:TRP:CZ2	1:B:76:LEU:HD21	2.49	0.47
1:B:82:THR:OG1	1:B:83:VAL:N	2.48	0.47
1:B:323:ARG:NH1	1:B:323:ARG:CG	2.70	0.47
1:B:9:ASN:HD22	1:B:10:TYR:N	2.13	0.47
2:B:803:4LG:H10	2:B:803:4LG:H241	1.76	0.46
1:B:128:ILE:CG1	1:B:137:PRO:HD3	2.43	0.46
1:A:151:ASP:HB3	1:A:323:ARG:HB2	1.97	0.46
1:B:15:TYR:O	1:B:31:VAL:HG22	2.16	0.46
1:A:294:HIS:HD2	1:A:294:HIS:O	1.98	0.46
1:B:185:TYR:CD2	1:B:326:ARG:HG3	2.50	0.46
1:A:189:ILE:CD1	1:A:215:LEU:HG	2.46	0.46
1:A:44:SER:HA	1:A:58:LEU:HB3	1.97	0.46
1:B:89:GLN:HE21	1:B:102:MET:HE3	1.81	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:76:LEU:O	1:B:83:VAL:HG13	2.16	0.46
1:A:226:SER:OG	1:A:307:ALA:HB3	2.15	0.45
1:A:150:GLU:C	1:A:152:VAL:H	2.20	0.45
1:A:154:SER:HB2	1:A:317:TYR:OH	2.17	0.45
1:B:142:ILE:O	1:B:145:GLN:HG2	2.16	0.45
1:B:133:GLY:O	1:B:134:ARG:HB3	2.16	0.45
1:B:200:GLY:HA2	1:B:210:CYS:O	2.16	0.45
1:B:272:LEU:HD23	1:B:276:ASP:CB	2.46	0.45
1:A:255:GLY:O	1:A:258:LEU:HG	2.17	0.45
1:B:87:LEU:CD2	1:B:135:VAL:HG21	2.45	0.45
1:B:112:LEU:HA	1:B:113:PRO:HA	1.67	0.45
1:B:50:TYR:O	1:B:51:THR:C	2.54	0.45
1:A:9:ASN:HD21	1:A:12:ASP:HA	1.82	0.44
1:A:44:SER:HA	1:A:58:LEU:HD13	2.00	0.44
1:B:64:SER:HB3	1:B:67:TYR:HB2	2.00	0.44
1:B:78:TYR:N	1:B:81:GLY:HA3	2.32	0.44
1:A:327:ILE:HG22	1:A:329:PHE:CE1	2.52	0.44
1:A:77:ARG:HA	1:A:82:THR:H	1.82	0.44
1:B:78:TYR:HE2	1:B:113:PRO:HD3	1.83	0.44
1:A:196:ILE:HD12	1:A:265:LEU:HD22	1.98	0.44
1:A:268:LYS:NZ	1:A:269:GLU:HB2	2.33	0.44
1:B:215:LEU:HB2	1:B:305:THR:HG22	1.97	0.44
1:B:5:VAL:HG13	1:B:5:VAL:O	2.17	0.44
1:A:106:VAL:HG11	1:A:109:MET:CG	2.48	0.44
1:B:6:ILE:N	1:B:6:ILE:HD12	2.32	0.44
1:A:218:VAL:HG21	1:A:327:ILE:HD13	2.00	0.44
1:A:250:VAL:O	1:A:251:LYS:C	2.54	0.44
1:A:32:PHE:HB2	1:A:169:ILE:HD11	1.99	0.44
1:A:134:ARG:HB2	1:A:134:ARG:NH2	2.32	0.44
1:B:50:TYR:CZ	1:B:108:GLU:CB	3.01	0.44
1:B:255:GLY:HA3	1:B:280:GLN:HE22	1.82	0.44
1:B:156:TYR:HB2	1:B:317:TYR:CE2	2.53	0.44
1:A:171:LEU:N	1:A:171:LEU:HD22	2.32	0.44
1:A:225:ILE:HD11	1:A:272:LEU:HD11	1.99	0.44
1:A:138:ILE:HG23	1:A:139:PHE:N	2.34	0.43
1:A:250:VAL:HG13	1:A:251:LYS:O	2.19	0.43
1:B:232:ILE:O	1:B:236:MET:HG2	2.18	0.43
1:B:69:HIS:ND1	1:B:70:ASN:N	2.66	0.43
1:B:245:LEU:HD23	1:B:246:PHE:HB3	2.01	0.43
1:B:222:ALA:O	2:B:803:4LG:H191	2.17	0.43
1:B:109:MET:HA	1:B:110:PRO:HD3	1.78	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:68:LYS:HB3	1:B:89:GLN:HB3	2.00	0.43
1:A:12:ASP:OD2	1:A:160:ASP:HB2	2.19	0.43
1:A:80:THR:O	1:A:81:GLY:O	2.36	0.43
1:A:81:GLY:O	1:A:82:THR:HB	2.18	0.43
1:B:128:ILE:HA	1:B:137:PRO:HG2	2.00	0.43
1:B:145:GLN:O	1:B:147:VAL:HG13	2.18	0.43
1:A:160:ASP:OD2	1:A:161:SER:N	2.50	0.43
1:B:10:TYR:CG	1:B:118:GLU:HG3	2.54	0.43
1:B:151:ASP:HB3	1:B:323:ARG:HB2	2.00	0.43
1:B:217:LEU:CD2	1:B:219:ASP:HB2	2.43	0.43
1:B:40:TRP:CB	2:B:803:4LG:H321	2.49	0.43
1:A:236:MET:HG3	1:A:248:TYR:CZ	2.53	0.43
1:A:85:GLY:HA3	1:A:105:GLU:O	2.18	0.43
1:B:232:ILE:HG22	1:B:232:ILE:O	2.19	0.43
1:B:290:THR:CG2	1:B:291:LEU:N	2.82	0.43
1:B:70:ASN:O	1:B:87:LEU:O	2.37	0.43
1:A:85:GLY:HA2	1:A:107:THR:OG1	2.19	0.42
1:A:110:PRO:HG2	1:A:114:PHE:HE1	1.84	0.42
1:A:223:SER:HA	2:A:804:4LG:H192	2.00	0.42
1:B:50:TYR:C	1:B:52:ALA:N	2.70	0.42
1:B:79:SER:O	1:B:80:THR:HB	2.19	0.42
1:A:194:TRP:CH2	1:A:320:PHE:HB3	2.55	0.42
1:A:236:MET:SD	1:A:248:TYR:CG	3.12	0.42
1:A:290:THR:CG2	1:A:291:LEU:N	2.82	0.42
1:A:290:THR:HG22	1:A:291:LEU:N	2.34	0.42
1:B:122:VAL:HG23	2:B:803:4LG:C1	2.49	0.42
1:A:239:LEU:CD2	1:A:259:PRO:HG2	2.48	0.42
1:A:40:TRP:CZ2	1:A:132:ILE:HD12	2.54	0.42
1:A:324:ASN:O	1:A:326:ARG:HG3	2.20	0.42
1:B:305:THR:HB	1:B:306:TRP:H	1.19	0.42
1:B:225:ILE:HG13	1:B:311:THR:HB	2.02	0.42
1:A:61:ALA:HB1	1:A:67:TYR:CD1	2.55	0.42
1:B:232:ILE:HG13	1:B:295:ALA:HA	2.00	0.42
1:B:13:THR:O	1:B:223:SER:OG	2.38	0.42
1:A:38:ASN:ND2	1:A:132:ILE:H	2.13	0.42
1:A:268:LYS:HA	1:A:268:LYS:HD2	1.70	0.42
1:B:191:THR:CG2	1:B:192:GLY:N	2.83	0.42
1:B:245:LEU:CD2	1:B:246:PHE:CB	2.96	0.42
1:A:177:GLN:HA	1:A:177:GLN:NE2	2.35	0.42
1:A:294:HIS:C	1:A:294:HIS:CD2	2.93	0.42
1:A:83:VAL:HG21	2:A:804:4LG:H322	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:102:MET:HB3	1:B:135:VAL:HG11	2.01	0.42
1:A:150:GLU:C	1:A:152:VAL:N	2.73	0.41
1:B:199:LYS:HZ2	1:B:266:GLY:HA2	1.81	0.41
1:A:99:VAL:HG12	1:A:145:GLN:OE1	2.20	0.41
1:A:228:SER:O	1:A:232:ILE:HG12	2.19	0.41
1:A:44:SER:HA	1:A:58:LEU:CD1	2.50	0.41
1:A:53:CYS:O	1:A:54:VAL:HG23	2.20	0.41
1:B:134:ARG:O	1:B:134:ARG:HG2	2.20	0.41
1:A:217:LEU:HD23	1:A:217:LEU:HA	1.83	0.41
1:A:331:LEU:HA	1:A:331:LEU:HD23	1.81	0.41
1:A:76:LEU:CD1	1:A:83:VAL:HG22	2.50	0.41
1:B:201:VAL:CG2	1:B:209:LEU:CB	2.98	0.41
1:B:16:TYR:HB3	1:B:31:VAL:HG23	2.02	0.41
1:B:7:LEU:HB2	1:B:167:GLY:CA	2.50	0.41
1:B:9:ASN:ND2	1:B:9:ASN:C	2.73	0.41
1:A:42:PRO:HB2	1:A:58:LEU:CD2	2.50	0.41
1:B:299:PRO:HA	1:B:300:PRO:HD3	1.86	0.41
1:A:139:PHE:CE1	1:A:143:ILE:HD11	2.55	0.41
1:A:81:GLY:O	1:A:82:THR:CG2	2.67	0.41
1:B:48:ARG:C	1:B:50:TYR:H	2.23	0.41
1:A:236:MET:HG3	1:A:248:TYR:CE1	2.55	0.41
1:B:117:ALA:HB3	2:B:803:4LG:H5	2.01	0.41
1:B:212:ASP:N	1:B:212:ASP:OD2	2.54	0.41
1:B:245:LEU:CG	1:B:246:PHE:N	2.83	0.41
1:B:77:ARG:N	1:B:77:ARG:HD2	2.35	0.41
1:A:103:PHE:N	1:A:103:PHE:CD1	2.89	0.41
2:A:804:4LG:H241	2:A:804:4LG:H10	1.84	0.41
1:A:168:GLN:HG2	1:A:169:ILE:N	2.36	0.41
1:B:236:MET:HG3	1:B:248:TYR:CG	2.56	0.41
1:B:50:TYR:O	1:B:52:ALA:N	2.54	0.41
1:A:109:MET:HE2	2:A:804:4LG:H8	2.03	0.41
1:A:201:VAL:HG21	1:A:306:TRP:CZ2	2.56	0.41
1:B:242:LYS:O	1:B:248:TYR:HA	2.21	0.41
1:A:190:LYS:O	1:A:191:THR:C	2.59	0.41
1:A:207:THR:HG22	1:A:207:THR:O	2.21	0.41
1:A:9:ASN:HD21	1:A:12:ASP:CA	2.35	0.40
1:A:225:ILE:HA	1:A:307:ALA:O	2.21	0.40
1:B:128:ILE:HG12	1:B:128:ILE:O	2.21	0.40
1:B:264:HIS:O	1:B:265:LEU:HD23	2.21	0.40
1:A:127:PHE:CD2	1:A:322:ARG:NH1	2.89	0.40
1:B:40:TRP:HZ2	1:B:76:LEU:HD21	1.86	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:200:GLY:N	1:A:210:CYS:SG	2.87	0.40
1:B:186:ILE:HG21	1:B:196:ILE:HD12	2.03	0.40
1:A:34:THR:HA	1:A:125:MET:HB2	2.04	0.40
1:B:143:ILE:HD13	1:B:148:LEU:HD12	2.02	0.40
1:B:317:TYR:N	1:B:332:ALA:HB2	2.36	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	324/333 (97%)	270 (83%)	37 (11%)	17 (5%)	2	6
1	B	324/333 (97%)	287 (89%)	28 (9%)	9 (3%)	5	19
All	All	648/666 (97%)	557 (86%)	65 (10%)	26 (4%)	3	11

All (26) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	79	SER
1	A	80	THR
1	A	81	GLY
1	A	82	THR
1	A	83	VAL
1	A	212	ASP
1	A	295	ALA
1	B	83	VAL
1	A	209	LEU
1	B	70	ASN
1	B	79	SER
1	B	213	GLY
1	A	296	MET

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Mol	Chain	Res	Type
1	B	80	THR
1	B	114	PHE
1	B	206	SER
1	B	266	GLY
1	A	12	ASP
1	A	54	VAL
1	A	189	ILE
1	A	236	MET
1	A	292	ALA
1	A	110	PRO
1	A	256	PRO
1	A	213	GLY
1	B	138	ILE

### 5.3.2 Protein sidechains [i](#)

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	279/284 (98%)	256 (92%)	23 (8%)	11	32
1	B	279/284 (98%)	249 (89%)	30 (11%)	6	20
All	All	558/568 (98%)	505 (90%)	53 (10%)	8	26

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	ASN
1	A	40	TRP
1	A	51	THR
1	A	76	LEU
1	A	77	ARG
1	A	87	LEU
1	A	94	VAL
1	A	106	VAL
1	A	158	ASN
1	A	199	LYS

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	A	208	LEU
1	A	217	LEU
1	A	237	GLU
1	A	245	LEU
1	A	250	VAL
1	A	251	LYS
1	A	273	THR
1	A	278	VAL
1	A	280	GLN
1	A	294	HIS
1	A	297	ASP
1	A	326	ARG
1	A	331	LEU
1	B	9	ASN
1	B	18	GLU
1	B	31	VAL
1	B	51	THR
1	B	77	ARG
1	B	102	MET
1	B	115	MET
1	B	140	ASP
1	B	190	LYS
1	B	208	LEU
1	B	212	ASP
1	B	217	LEU
1	B	223	SER
1	B	224	TYR
1	B	237	GLU
1	B	239	LEU
1	B	243	LYS
1	B	245	LEU
1	B	246	PHE
1	B	250	VAL
1	B	258	LEU
1	B	272	LEU
1	B	282	SER
1	B	286	LYS
1	B	291	LEU
1	B	305	THR
1	B	315	LYS
1	B	323	ARG
1	B	325	ASN

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Mol	Chain	Res	Type
1	B	326	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	ASN
1	A	38	ASN
1	A	101	GLN
1	A	130	GLN
1	A	141	ASN
1	A	145	GLN
1	A	158	ASN
1	A	168	GLN
1	A	177	GLN
1	A	264	HIS
1	A	294	HIS
1	B	9	ASN
1	B	26	GLN
1	B	89	GLN
1	B	101	GLN
1	B	168	GLN
1	B	280	GLN

### 5.3.3 RNA [i](#)

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 carbohydrates 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	4LG	A	804	-	47,47,47	1.69	13 (27%)	52,63,63	2.10	11 (21%)
2	4LG	B	803	-	47,47,47	1.88	14 (29%)	52,63,63	2.22	13 (25%)

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	4LG	A	804	-	-	9/26/40/40	0/5/5/5
2	4LG	B	803	-	-	10/26/40/40	0/5/5/5

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	803	4LG	C11-N1	-4.91	1.33	1.39
2	B	803	4LG	C13-C14	4.23	1.45	1.36
2	A	804	4LG	C13-C14	3.96	1.44	1.36
2	A	804	4LG	C11-N1	-3.47	1.35	1.39
2	B	803	4LG	C10-C11	-3.25	1.34	1.40
2	B	803	4LG	O1-C2	3.25	1.42	1.37
2	B	803	4LG	C23-N3	3.10	1.42	1.36
2	A	804	4LG	C22-C23	3.09	1.57	1.50
2	A	804	4LG	C14-C9	3.01	1.44	1.38
2	B	803	4LG	C14-C9	2.88	1.44	1.38
2	A	804	4LG	C10-C11	-2.78	1.35	1.40
2	B	803	4LG	C22-C23	2.76	1.56	1.50
2	B	803	4LG	C17-C18	-2.68	1.46	1.52
2	A	804	4LG	C17-C18	-2.65	1.46	1.52
2	A	804	4LG	C21-C20	2.59	1.55	1.52
2	B	803	4LG	C21-C20	2.57	1.55	1.52
2	A	804	4LG	C21-N2	-2.45	1.43	1.46
2	A	804	4LG	C27-C26	2.45	1.44	1.39
2	A	804	4LG	C23-N3	2.38	1.41	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	803	4LG	C28-C29	2.24	1.43	1.38
2	B	803	4LG	C27-C26	2.21	1.43	1.39
2	B	803	4LG	C21-N2	-2.14	1.43	1.46
2	B	803	4LG	C28-C27	2.11	1.42	1.38
2	A	804	4LG	C28-C29	2.11	1.42	1.38
2	A	804	4LG	C2-C3	2.10	1.43	1.40
2	A	804	4LG	C28-C27	2.03	1.42	1.38
2	B	803	4LG	C20-N3	2.00	1.51	1.48

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	803	4LG	C17-N1-C15	8.80	137.71	124.73
2	A	804	4LG	C17-N1-C15	8.10	136.68	124.73
2	A	804	4LG	O1-C2-C3	5.58	123.75	115.97
2	B	803	4LG	C10-C11-N1	-5.35	126.23	132.50
2	B	803	4LG	O1-C2-C3	4.93	122.83	115.97
2	A	804	4LG	C10-C11-N1	-4.65	127.05	132.50
2	A	804	4LG	O1-C2-C5	-4.49	116.67	124.37
2	B	803	4LG	O1-C2-C5	-3.85	117.77	124.37
2	B	803	4LG	C24-C20-N3	3.49	121.50	111.17
2	A	804	4LG	C24-C20-N3	3.28	120.86	111.17
2	B	803	4LG	O2-C4-C3	3.21	119.88	109.83
2	B	803	4LG	C9-C10-C11	2.89	123.95	118.10
2	A	804	4LG	C4-O2-C33	2.84	121.75	112.86
2	B	803	4LG	O5-C23-C22	-2.75	115.10	120.14
2	B	803	4LG	C14-C9-C10	-2.71	117.16	120.81
2	A	804	4LG	C9-C10-C11	2.59	123.35	118.10
2	B	803	4LG	C27-C26-N3	2.56	122.95	120.13
2	A	804	4LG	C26-N3-C23	2.47	121.03	118.60
2	A	804	4LG	C14-C9-C10	-2.39	117.59	120.81
2	B	803	4LG	C16-C12-C11	-2.25	104.28	106.20
2	A	804	4LG	O5-C23-C22	-2.20	116.11	120.14
2	A	804	4LG	C16-C12-C11	-2.20	104.32	106.20
2	B	803	4LG	C4-C3-C2	-2.02	117.20	120.49
2	B	803	4LG	C26-N3-C20	2.02	120.58	117.45

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	804	4LG	C18-C17-N1-C15

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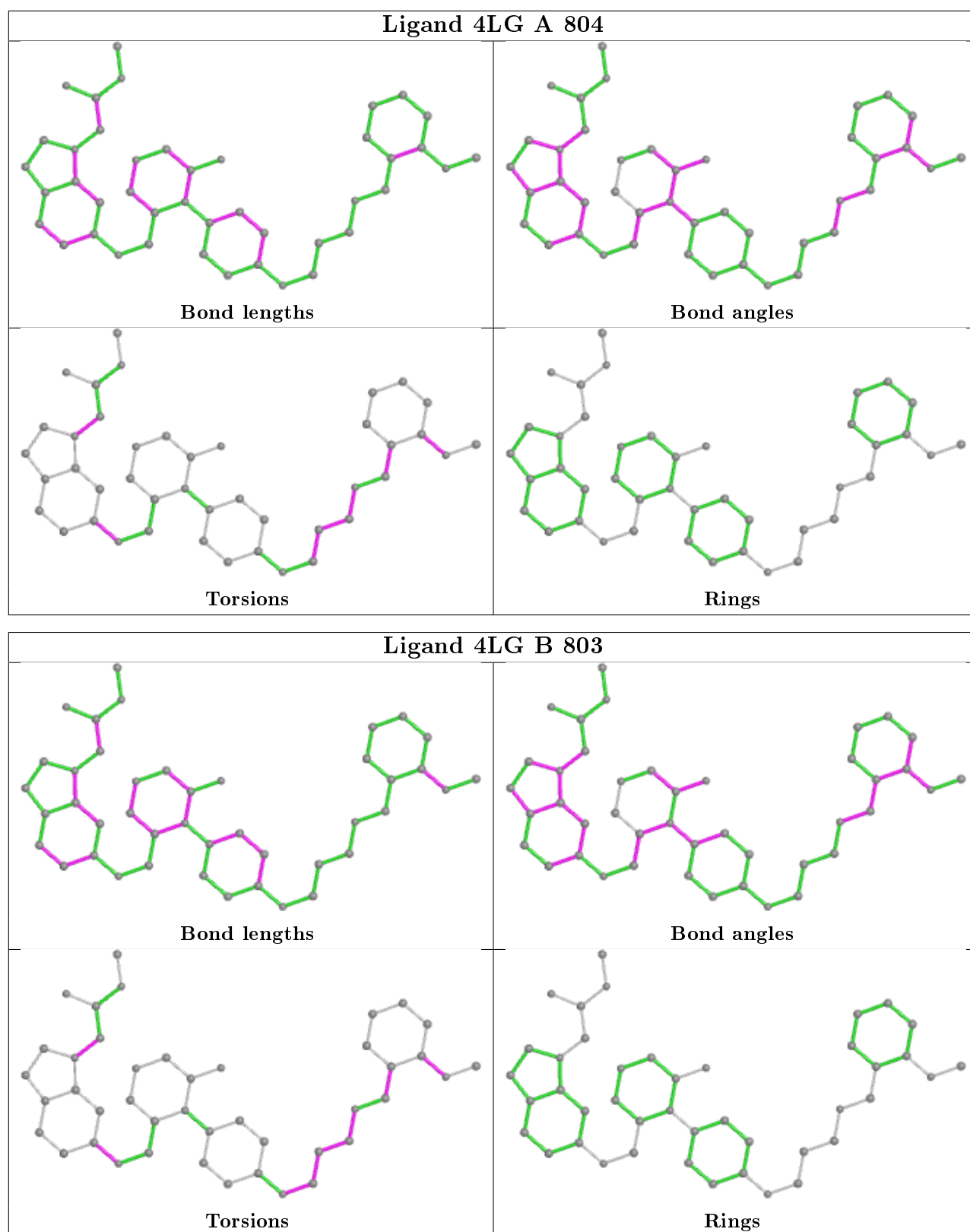
Mol	Chain	Res	Type	Atoms
2	B	803	4LG	C18-C17-N1-C15
2	B	803	4LG	C3-C2-O1-C1
2	A	804	4LG	C3-C2-O1-C1
2	B	803	4LG	C32-C33-O2-C4
2	A	804	4LG	C5-C2-O1-C1
2	A	804	4LG	O7-C31-C32-C33
2	B	803	4LG	C5-C2-O1-C1
2	A	804	4LG	C31-C32-C33-O2
2	A	804	4LG	C14-C9-O6-C24
2	B	803	4LG	C14-C9-O6-C24
2	A	804	4LG	C10-C9-O6-C24
2	B	803	4LG	C10-C9-O6-C24
2	B	803	4LG	C31-C32-C33-O2
2	B	803	4LG	C32-C31-O7-C29
2	A	804	4LG	C32-C33-O2-C4
2	B	803	4LG	O7-C31-C32-C33
2	B	803	4LG	C2-C3-C4-O2
2	A	804	4LG	C2-C3-C4-O2

There are no ring outliers.

2 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	804	4LG	9	0
2	B	803	4LG	12	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

There are no chain breaks in this entry.



## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.3 Carbohydrates

Unable to reproduce the depositors R factor - this section is therefore empty.

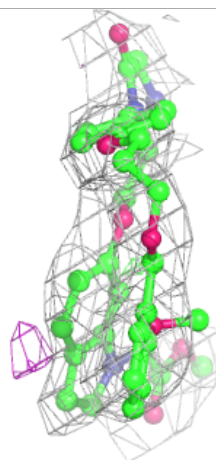
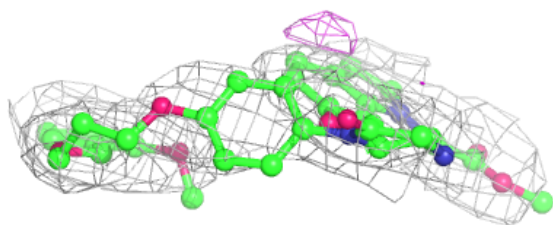
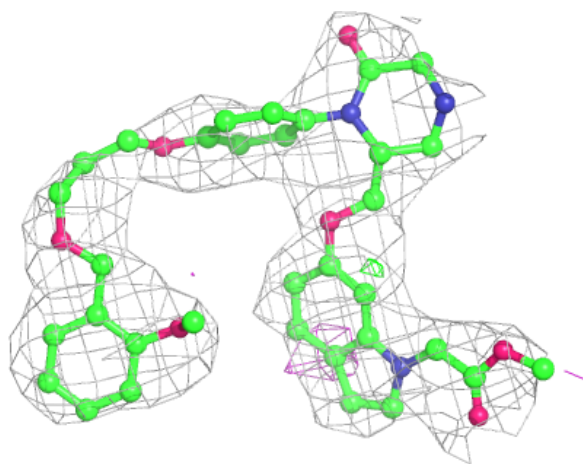
### 6.4 Ligands

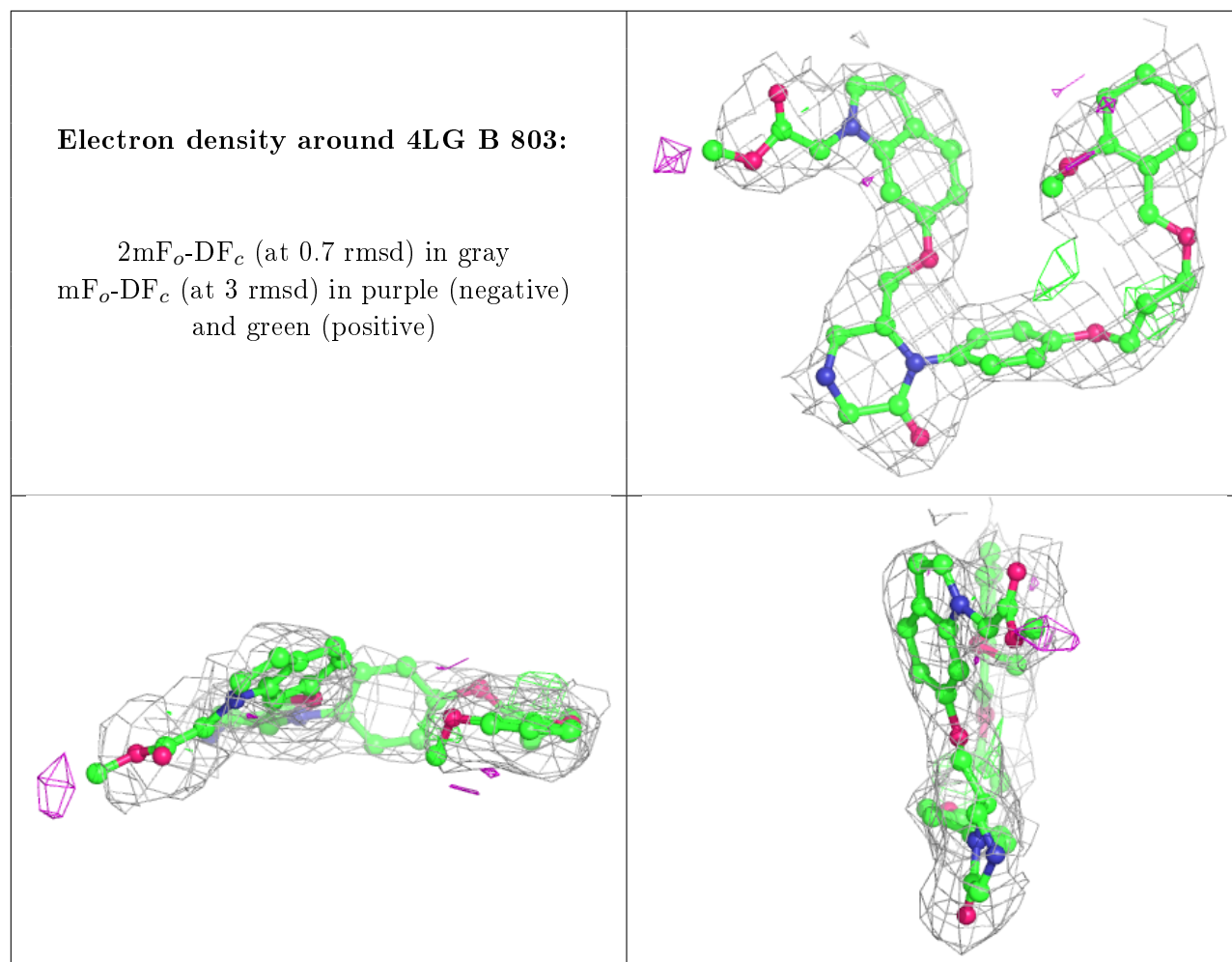
Unable to reproduce the depositors R factor - this section is therefore empty.

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 4LG A 804:**

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





## 6.5 Other polymers [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.