



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 11:16 AM GMT

PDB ID : 2FPB  
Title : Structure of Strictosidine Synthase, the Biosynthetic Entry to the Monoterpenoid Indole Alkaloid Family  
Authors : Panjikar, S.  
Deposited on : 2006-01-16  
Resolution : 2.80 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>

---

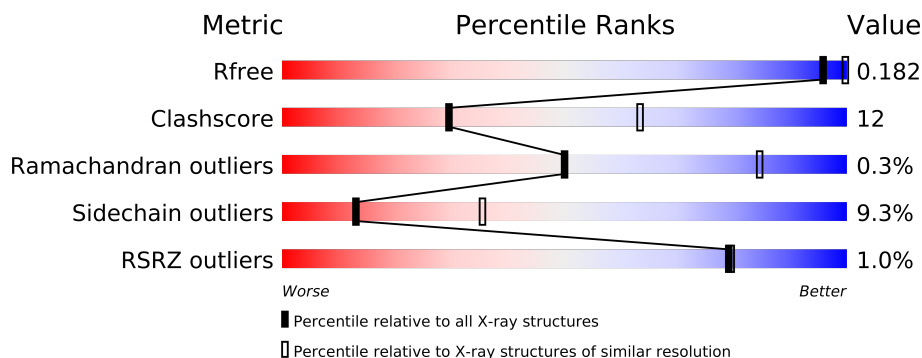
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.80 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	1799 (2.80-2.80)
Clashscore	79885	2295 (2.80-2.80)
Ramachandran outliers	78287	2252 (2.80-2.80)
Sidechain outliers	78261	2254 (2.80-2.80)
RSRZ outliers	66119	1802 (2.80-2.80)

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	322	
1	B	322	

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

Mol	Type	Chain	Res	Geometry	Electron density
2	TSS	A	1001	-	X
2	TSS	B	2001	-	X

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5007 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.

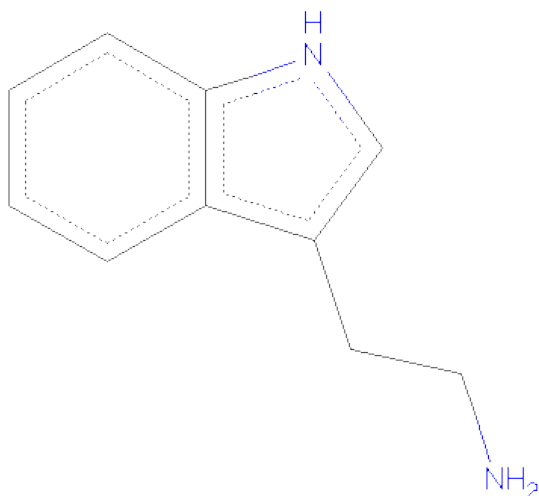
- Molecule 1 is a protein called Strictosidine Synthase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	305	Total	C	N	O	S	Se	0	0	0
			2405	1542	391	463	3	6			
1	B	305	Total	C	N	O	S	Se	0	0	0
			2405	1542	391	463	3	6			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	65	MSE	ILE	ENGINEERED	UNP P68175
A	116	MSE	LEU	ENGINEERED	UNP P68175
A	180	MSE	MET	MODIFIED RESIDUE	UNP P68175
A	190	MSE	ILE	ENGINEERED	UNP P68175
A	203	MSE	LEU	ENGINEERED	UNP P68175
A	276	MSE	MET	MODIFIED RESIDUE	UNP P68175
B	65	MSE	ILE	ENGINEERED	UNP P68175
B	116	MSE	LEU	ENGINEERED	UNP P68175
B	180	MSE	MET	MODIFIED RESIDUE	UNP P68175
B	190	MSE	ILE	ENGINEERED	UNP P68175
B	203	MSE	LEU	ENGINEERED	UNP P68175
B	276	MSE	MET	MODIFIED RESIDUE	UNP P68175

- Molecule 2 is 2-(1H-INDOL-3-YL)ETHANAMINE (three-letter code: TSS) (formula: C<sub>10</sub>H<sub>12</sub>N<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	N	0	0
			12	10	2		
2	B	1	Total	C	N	0	0
			12	10	2		

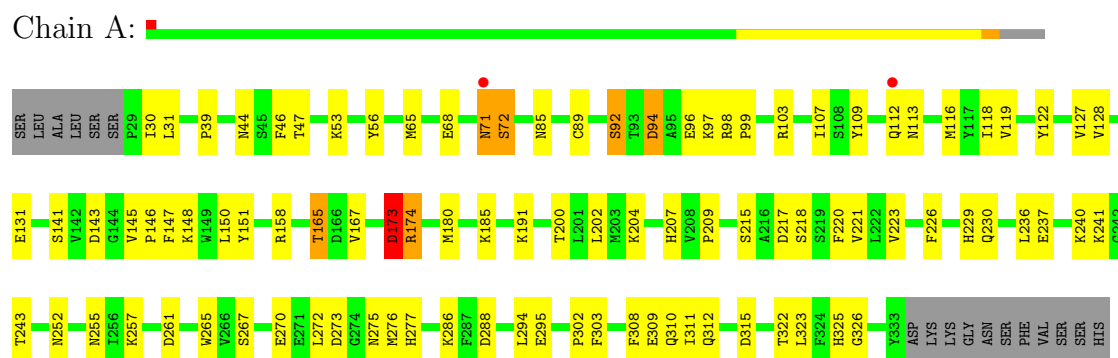
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	93	Total	O	0	0
			93	93		
3	B	80	Total	O	0	0
			80	80		

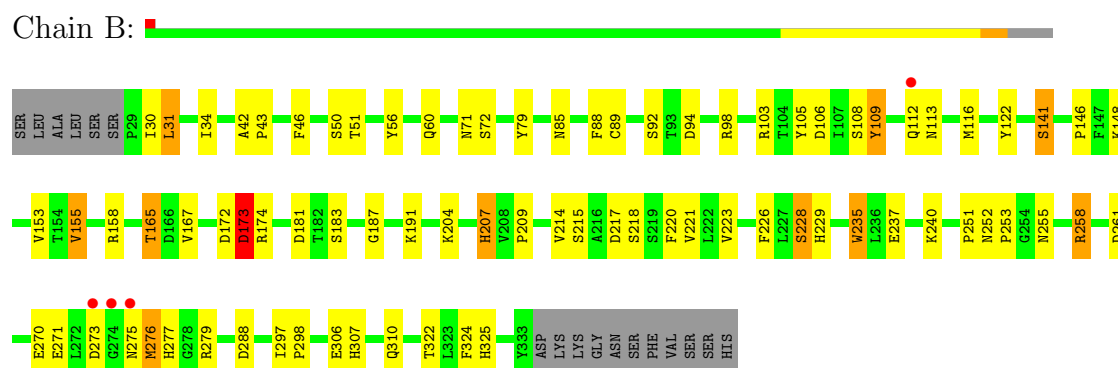
### 3 Residue-property plots

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: Strictosidine Synthase



#### • Molecule 1: Strictosidine Synthase



## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	148.94Å 148.94Å 121.34Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.80 19.65 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.9 (20.00-2.80) 99.9 (19.65-2.80)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.01 (at 2.79Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, $R_{free}$	0.164 , 0.217 0.182 , 0.182	Depositor DCC
$R_{free}$ test set	1234 reflections (5.27%)	DCC
Wilson B-factor (Å <sup>2</sup> )	49.7	Xtriage
Anisotropy	0.742	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 27.0	EDS
Estimated twinning fraction	0.028 for h,-h-k,-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 24636 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	5007	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.12% of the height of the origin peak. No significant pseudotranslation is detected.*

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

## 5 Model quality

### 5.1 Standard geometry

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

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	0.86	0/2465	0.95	5/3340 (0.1%)
1	B	0.77	0/2465	0.91	6/3340 (0.2%)
All	All	0.82	0/4930	0.93	11/6680 (0.2%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	143	ASP	CB-CG-OD2	7.49	125.04	118.30
1	A	273	ASP	CB-CG-OD2	7.12	124.71	118.30
1	B	94	ASP	CB-CG-OD2	6.50	124.15	118.30
1	B	273	ASP	CB-CG-OD2	5.80	123.52	118.30
1	A	261	ASP	CB-CG-OD2	5.66	123.39	118.30
1	B	288	ASP	CB-CG-OD2	5.48	123.24	118.30
1	B	261	ASP	CB-CG-OD2	5.46	123.21	118.30
1	B	172	ASP	CB-CG-OD2	5.41	123.17	118.30
1	A	315	ASP	CB-CG-OD2	5.37	123.13	118.30
1	B	181	ASP	CB-CG-OD2	5.09	122.88	118.30
1	A	173	ASP	CB-CA-C	-5.03	100.34	110.40

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 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	2405	0	2309	68	0
1	B	2405	0	2309	41	0
2	A	12	0	12	6	0
2	B	12	0	12	5	0
3	A	93	0	0	2	0
3	B	80	0	0	2	0
All	All	5007	0	4642	116	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 12.

All (116) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:276:MSE:CE	1:B:276:MSE:SE	2.15	1.44
1:A:180:MSE:CE	1:A:180:MSE:SE	2.18	1.42
1:A:276:MSE:SE	1:A:276:MSE:CE	2.18	1.38
2:B:2001:TSS:HN11	2:B:2001:TSS:HD1	1.09	1.08
1:A:255:ASN:HD22	1:A:310:GLN:HB2	1.17	1.05
2:A:1001:TSS:N1	2:A:1001:TSS:HD1	1.71	1.03
2:A:1001:TSS:HN11	2:A:1001:TSS:CD1	1.70	0.96
2:A:1001:TSS:HD1	2:A:1001:TSS:HN11	1.28	0.95
1:B:229:HIS:HD2	1:B:252:ASN:H	1.00	0.93
1:A:229:HIS:HD2	1:A:252:ASN:H	1.03	0.91
2:A:1001:TSS:N1	2:A:1001:TSS:CD1	2.31	0.90
1:B:229:HIS:CD2	1:B:252:ASN:H	1.90	0.89
2:B:2001:TSS:N1	2:B:2001:TSS:HD1	1.87	0.87
1:A:68:GLU:HB2	1:A:72:SER:HB2	1.56	0.86
2:B:2001:TSS:HN11	2:B:2001:TSS:CD1	1.84	0.86
1:A:158:ARG:HB2	1:A:158:ARG:NH1	1.91	0.85
1:A:96:GLU:HB3	3:A:1046:HOH:O	1.78	0.83
1:A:39:PRO:HG3	1:A:65:MSE:HE2	1.64	0.80
1:A:229:HIS:CD2	1:A:252:ASN:H	1.95	0.79
1:A:158:ARG:HH11	1:A:158:ARG:CB	1.98	0.77
1:B:255:ASN:HD22	1:B:310:GLN:HB2	1.48	0.76
1:A:158:ARG:CB	1:A:158:ARG:NH1	2.48	0.76
1:A:221:VAL:CG2	1:A:236:LEU:HD11	2.18	0.73
1:B:167:VAL:HA	1:B:209:PRO:HD2	1.70	0.73
1:A:221:VAL:HG23	1:A:236:LEU:HD11	1.71	0.73
1:A:322:THR:OG1	1:A:325:HIS:HD2	1.73	0.72
1:B:60:GLN:HE22	1:B:98:ARG:HH21	1.35	0.71

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:275:ASN:ND2	1:A:277:HIS:H	1.90	0.69
1:A:71:ASN:HD22	1:A:72:SER:H	1.40	0.69
1:B:165:THR:HG21	1:B:209:PRO:O	1.93	0.68
1:B:306:GLU:HG2	1:B:324:PHE:CZ	2.29	0.68
1:B:141:SER:HB2	1:B:146:PRO:HA	1.75	0.67
1:A:158:ARG:HH11	1:A:158:ARG:HB3	1.61	0.65
1:A:255:ASN:ND2	1:A:310:GLN:HB2	2.01	0.64
1:B:106:ASP:HB3	1:B:153:VAL:HG12	1.78	0.64
1:A:94:ASP:N	1:A:94:ASP:OD1	2.32	0.61
1:A:122:TYR:OH	1:A:173:ASP:HB2	2.02	0.60
1:B:207:HIS:HD2	1:B:228:SER:OG	1.85	0.59
1:A:98:ARG:HD2	1:A:173:ASP:OD1	2.03	0.58
1:B:217:ASP:O	1:B:218:SER:HB2	2.02	0.58
1:A:229:HIS:HD2	1:A:252:ASN:N	1.88	0.58
1:B:103:ARG:HD2	1:B:105:TYR:CZ	2.37	0.58
1:A:180:MSE:HE3	3:A:1052:HOH:O	2.04	0.57
1:A:275:ASN:HD22	1:A:277:HIS:H	1.52	0.57
1:B:215:SER:HB3	1:B:220:PHE:CZ	2.40	0.56
1:A:191:LYS:HD3	1:A:202:LEU:HD21	1.87	0.56
1:A:267:SER:HB2	1:A:308:PHE:CZ	2.41	0.55
1:B:324:PHE:C	1:B:324:PHE:CD2	2.79	0.55
1:B:235:TRP:CZ3	1:B:240:LYS:HD2	2.42	0.55
1:A:226:PHE:CE1	2:A:1001:TSS:HB1	2.43	0.54
1:B:103:ARG:HD2	1:B:105:TYR:CE2	2.43	0.54
1:B:322:THR:HG21	1:B:325:HIS:HB3	1.89	0.52
1:B:252:ASN:ND2	1:B:271:GLU:H	2.07	0.52
1:B:324:PHE:O	1:B:324:PHE:CD2	2.63	0.51
1:B:226:PHE:HA	1:B:253:PRO:HD2	1.92	0.51
1:B:271:GLU:OE1	1:B:307:HIS:HE1	1.94	0.51
1:A:141:SER:CB	1:A:146:PRO:HA	2.41	0.50
1:B:306:GLU:HG2	1:B:324:PHE:CE1	2.46	0.50
1:B:214:VAL:O	1:B:258:ARG:NH1	2.45	0.50
1:B:85:ASN:HD22	1:B:88:PHE:H	1.58	0.49
1:A:309:GLU:OE2	2:A:1001:TSS:N1	2.45	0.49
1:A:158:ARG:CZ	1:A:158:ARG:HB2	2.43	0.49
1:A:174:ARG:HG3	1:A:174:ARG:HH11	1.78	0.49
1:A:46:PHE:HA	1:A:56:TYR:O	2.14	0.48
1:B:60:GLN:NE2	1:B:98:ARG:HH21	2.08	0.48
1:A:257:LYS:CB	1:A:312:GLN:HG3	2.44	0.48
1:A:151:TYR:HB2	1:A:167:VAL:HG13	1.96	0.48
1:A:119:VAL:O	1:A:119:VAL:HG23	2.15	0.47
1:A:302:PRO:HD2	1:A:303:PHE:CE2	2.49	0.47

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:165:THR:HG21	1:A:209:PRO:O	2.15	0.47
1:B:42:ALA:N	1:B:43:PRO:HD3	2.30	0.47
1:A:85:ASN:O	1:A:89:CYS:HB2	2.15	0.46
2:B:2001:TSS:HA1	3:B:2064:HOH:O	2.16	0.46
2:B:2001:TSS:CD1	2:B:2001:TSS:N1	2.54	0.46
1:A:215:SER:HB3	1:A:220:PHE:CZ	2.51	0.46
1:A:217:ASP:O	1:A:218:SER:HB2	2.17	0.45
1:B:122:TYR:OH	1:B:173:ASP:HB2	2.17	0.45
1:A:148:LYS:HA	1:A:148:LYS:HD2	1.90	0.44
1:A:98:ARG:NH1	1:A:173:ASP:HB3	2.32	0.44
1:A:96:GLU:H	1:A:96:GLU:CD	2.20	0.44
1:A:322:THR:HG21	1:A:325:HIS:HB2	1.99	0.44
1:A:323:LEU:HD23	1:A:323:LEU:HA	1.72	0.44
1:A:286:LYS:HB3	1:A:295:GLU:HB3	2.00	0.44
1:A:98:ARG:N	1:A:99:PRO:CD	2.80	0.44
1:A:288:ASP:HB3	1:A:294:LEU:HD11	1.99	0.44
1:A:167:VAL:HA	1:A:209:PRO:HD2	2.00	0.43
1:A:191:LYS:HG2	1:A:200:THR:HB	2.00	0.43
1:A:236:LEU:O	1:A:241:LYS:HG3	2.18	0.43
1:A:44:ASN:OD1	1:A:103:ARG:NH1	2.52	0.43
1:A:141:SER:HB3	1:A:146:PRO:HA	2.00	0.43
1:B:229:HIS:HD2	1:B:252:ASN:N	1.86	0.43
1:A:221:VAL:HG22	1:A:236:LEU:HD11	1.95	0.43
1:A:92:SER:HB2	1:A:97:LYS:HD2	2.01	0.43
1:A:240:LYS:O	1:A:243:THR:HB	2.18	0.43
1:B:297:ILE:HA	1:B:298:PRO:HD2	1.91	0.43
1:A:147:PHE:CG	1:A:150:LEU:HD21	2.55	0.42
1:A:145:VAL:HA	1:A:146:PRO:HD3	1.78	0.42
1:B:85:ASN:O	1:B:89:CYS:HB2	2.20	0.42
1:A:265:TRP:HB3	1:A:311:ILE:HG23	2.02	0.42
1:A:116:MSE:HE3	1:A:118:ILE:HG12	2.02	0.42
1:A:71:ASN:HD22	1:A:72:SER:N	2.11	0.41
1:B:240:LYS:HB2	3:B:2043:HOH:O	2.20	0.41
1:A:85:ASN:HA	1:A:85:ASN:HD22	1.74	0.41
1:B:46:PHE:HA	1:B:56:TYR:O	2.21	0.41
1:A:326:GLY:H	1:B:71:ASN:HA	1.85	0.41
1:A:47:THR:HG22	1:A:107:ILE:HD12	2.03	0.41
1:B:108:SER:OG	1:B:155:VAL:HG23	2.21	0.41
1:A:270:GLU:HG2	1:A:272:LEU:HD23	2.03	0.41
1:B:275:ASN:ND2	1:B:277:HIS:H	2.18	0.41
1:B:187:GLY:HA3	1:B:209:PRO:HD3	2.02	0.41
1:B:71:ASN:N	1:B:71:ASN:ND2	2.69	0.41

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:71:ASN:ND2	1:A:71:ASN:N	2.68	0.40
1:A:31:LEU:HB3	1:B:31:LEU:HD23	2.02	0.40
1:A:127:VAL:HG22	1:A:128:VAL:N	2.37	0.40
1:B:109:TYR:CE1	1:B:116:MSE:CE	3.05	0.40
1:B:251:PRO:HB2	1:B:270:GLU:HG3	2.02	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	303/322 (94%)	289 (95%)	13 (4%)	1 (0%)	50	84
1	B	303/322 (94%)	289 (95%)	13 (4%)	1 (0%)	50	84
All	All	606/644 (94%)	578 (95%)	26 (4%)	2 (0%)	50	84

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	173	ASP
1	B	173	ASP

### 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	263/272 (97%)	244 (93%)	19 (7%)	21	50
1	B	263/272 (97%)	233 (89%)	30 (11%)	8	24
All	All	526/544 (97%)	477 (91%)	49 (9%)	13	35

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	ILE
1	A	53	LYS
1	A	71	ASN
1	A	72	SER
1	A	92	SER
1	A	94	ASP
1	A	109	TYR
1	A	112	GLN
1	A	113	ASN
1	A	131	GLU
1	A	165	THR
1	A	173	ASP
1	A	174	ARG
1	A	185	LYS
1	A	204	LYS
1	A	207	HIS
1	A	223	VAL
1	A	230	GLN
1	A	237	GLU
1	B	30	ILE
1	B	31	LEU
1	B	34	ILE
1	B	50	SER
1	B	51	THR
1	B	72	SER
1	B	79	TYR
1	B	92	SER
1	B	109	TYR
1	B	112	GLN
1	B	113	ASN
1	B	141	SER
1	B	148	LYS
1	B	155	VAL
1	B	158	ARG
1	B	165	THR
1	B	173	ASP
1	B	174	ARG
1	B	183	SER
1	B	191	LYS
1	B	204	LYS
1	B	207	HIS
1	B	221	VAL

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	B	223	VAL
1	B	228	SER
1	B	235	TRP
1	B	237	GLU
1	B	258	ARG
1	B	276	MSE
1	B	279	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	71	ASN
1	A	85	ASN
1	A	113	ASN
1	A	115	GLN
1	A	207	HIS
1	A	229	HIS
1	A	230	GLN
1	A	252	ASN
1	A	255	ASN
1	A	275	ASN
1	A	292	ASN
1	A	310	GLN
1	A	312	GLN
1	A	325	HIS
1	B	60	GLN
1	B	71	ASN
1	B	85	ASN
1	B	91	ASN
1	B	113	ASN
1	B	207	HIS
1	B	229	HIS
1	B	230	GLN
1	B	252	ASN
1	B	255	ASN
1	B	275	ASN
1	B	307	HIS
1	B	310	GLN
1	B	312	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 ⓘ

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	TSS	A	1001	-	13,13,13	1.51	2 (15%)	17,17,17	0.73	0
2	TSS	B	2001	-	13,13,13	1.17	1 (7%)	17,17,17	0.99	1 (5%)

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	TSS	A	1001	-	-	0/3/3/3	0/0/2/2
2	TSS	B	2001	-	-	0/3/3/3	0/0/2/2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1001	TSS	CD1-CG	-3.18	1.32	1.37

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1001	TSS	CD1-NE1	-2.33	1.33	1.37
2	B	2001	TSS	CB-CG	2.01	1.56	1.51

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2001	TSS	CZ3-CE3-CD2	-2.22	117.64	120.88

There are no chirality outliers.

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	305/322 (94%)	-0.62	2 (0%) 84 85	15, 21, 28, 66	0
1	B	305/322 (94%)	-0.53	4 (1%) 74 75	15, 21, 28, 67	0
All	All	610/644 (94%)	-0.58	6 (0%) 79 79	15, 21, 28, 67	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	274	GLY	3.8
1	B	112	GLN	2.8
1	A	112	GLN	2.7
1	A	71	ASN	2.6
1	B	275	ASN	2.1
1	B	273	ASP	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( $\text{\AA}^2$ )	Q < 0.9
2	TSS	B	2001	12/12	0.24	3.38	55,70,74,75	0
2	TSS	A	1001	12/12	0.22	2.68	55,65,74,74	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.