



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 06:20 AM UTC

PDB ID : 1TZ3 / pdb_00001tz3
Title : crystal structure of aminoimidazole riboside kinase complexed with aminoimidazole riboside
Authors : Zhang, Y.; Dougherty, M.; Downs, D.M.; Ealick, S.E.
Deposited on : 2004-07-09
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

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
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

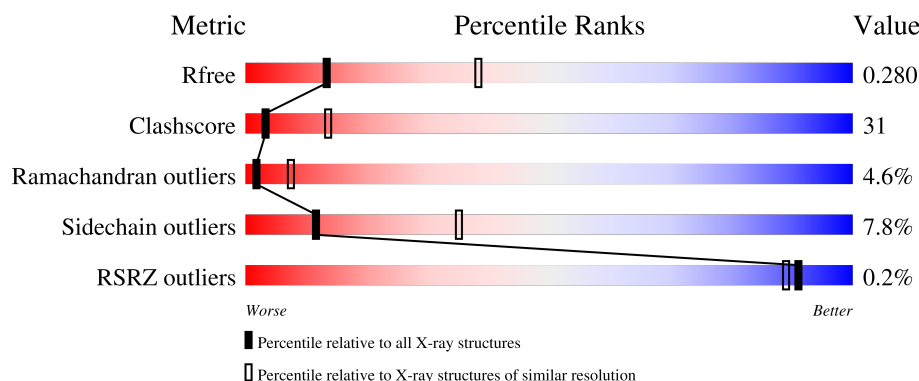
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

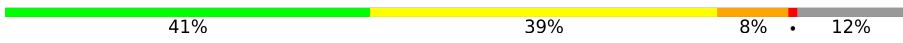
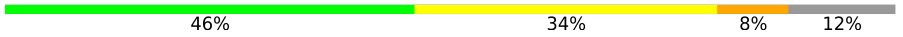
The reported resolution of this entry is 2.90 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 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	339	
1	B	339	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4485 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 putative sugar kinase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	299	Total	C	N	O	S	Se	0	0	0
			2221	1407	377	423	10	4			
1	B	298	Total	C	N	O	S	Se	0	0	0
			2204	1393	376	421	10	4			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	cloning artifact	UNP Q8ZKR2
A	-18	GLY	-	cloning artifact	UNP Q8ZKR2
A	-17	SER	-	cloning artifact	UNP Q8ZKR2
A	-16	SER	-	cloning artifact	UNP Q8ZKR2
A	-15	HIS	-	cloning artifact	UNP Q8ZKR2
A	-14	HIS	-	cloning artifact	UNP Q8ZKR2
A	-13	HIS	-	cloning artifact	UNP Q8ZKR2
A	-12	HIS	-	cloning artifact	UNP Q8ZKR2
A	-11	HIS	-	cloning artifact	UNP Q8ZKR2
A	-10	HIS	-	cloning artifact	UNP Q8ZKR2
A	-9	SER	-	cloning artifact	UNP Q8ZKR2
A	-8	SER	-	cloning artifact	UNP Q8ZKR2
A	-7	GLY	-	cloning artifact	UNP Q8ZKR2
A	-6	LEU	-	cloning artifact	UNP Q8ZKR2
A	-5	VAL	-	cloning artifact	UNP Q8ZKR2
A	-4	PRO	-	cloning artifact	UNP Q8ZKR2
A	-3	ARG	-	cloning artifact	UNP Q8ZKR2
A	-2	GLY	-	cloning artifact	UNP Q8ZKR2
A	-1	SER	-	cloning artifact	UNP Q8ZKR2
A	0	HIS	-	cloning artifact	UNP Q8ZKR2
A	148	MSE	MET	modified residue	UNP Q8ZKR2
A	165	MSE	MET	modified residue	UNP Q8ZKR2
A	286	MSE	MET	modified residue	UNP Q8ZKR2
A	294	MSE	MET	modified residue	UNP Q8ZKR2
B	-19	MET	-	cloning artifact	UNP Q8ZKR2

Continued on next page...

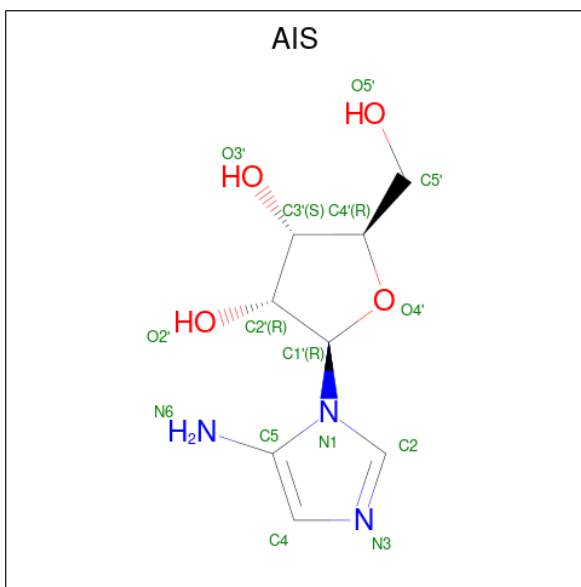
Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	-18	GLY	-	cloning artifact	UNP Q8ZKR2
B	-17	SER	-	cloning artifact	UNP Q8ZKR2
B	-16	SER	-	cloning artifact	UNP Q8ZKR2
B	-15	HIS	-	cloning artifact	UNP Q8ZKR2
B	-14	HIS	-	cloning artifact	UNP Q8ZKR2
B	-13	HIS	-	cloning artifact	UNP Q8ZKR2
B	-12	HIS	-	cloning artifact	UNP Q8ZKR2
B	-11	HIS	-	cloning artifact	UNP Q8ZKR2
B	-10	HIS	-	cloning artifact	UNP Q8ZKR2
B	-9	SER	-	cloning artifact	UNP Q8ZKR2
B	-8	SER	-	cloning artifact	UNP Q8ZKR2
B	-7	GLY	-	cloning artifact	UNP Q8ZKR2
B	-6	LEU	-	cloning artifact	UNP Q8ZKR2
B	-5	VAL	-	cloning artifact	UNP Q8ZKR2
B	-4	PRO	-	cloning artifact	UNP Q8ZKR2
B	-3	ARG	-	cloning artifact	UNP Q8ZKR2
B	-2	GLY	-	cloning artifact	UNP Q8ZKR2
B	-1	SER	-	cloning artifact	UNP Q8ZKR2
B	0	HIS	-	cloning artifact	UNP Q8ZKR2
B	148	MSE	MET	modified residue	UNP Q8ZKR2
B	165	MSE	MET	modified residue	UNP Q8ZKR2
B	286	MSE	MET	modified residue	UNP Q8ZKR2
B	294	MSE	MET	modified residue	UNP Q8ZKR2

- Molecule 2 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	2	Total K 2 2	0	0
2	B	2	Total K 2 2	0	0

- Molecule 3 is 5-AMINOIMIDAZOLE RIBONUCLEOSIDE (CCD ID: AIS) (formula: C₈H₁₃N₃O₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			15	8	3	4		

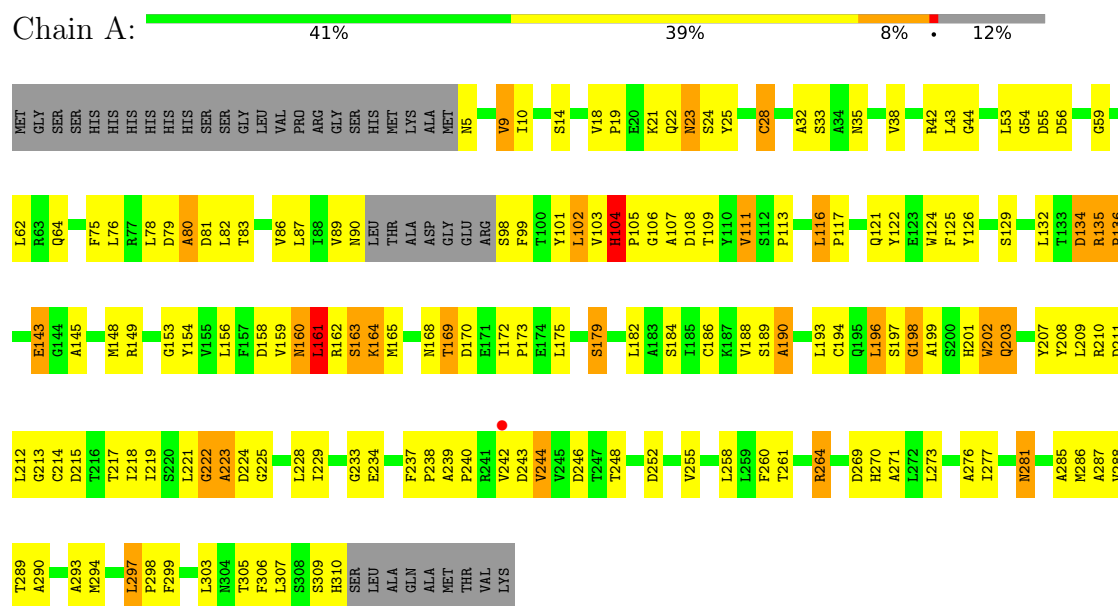
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	20	Total	O	0	0
			20	20		
4	B	21	Total	O	0	0
			21	21		

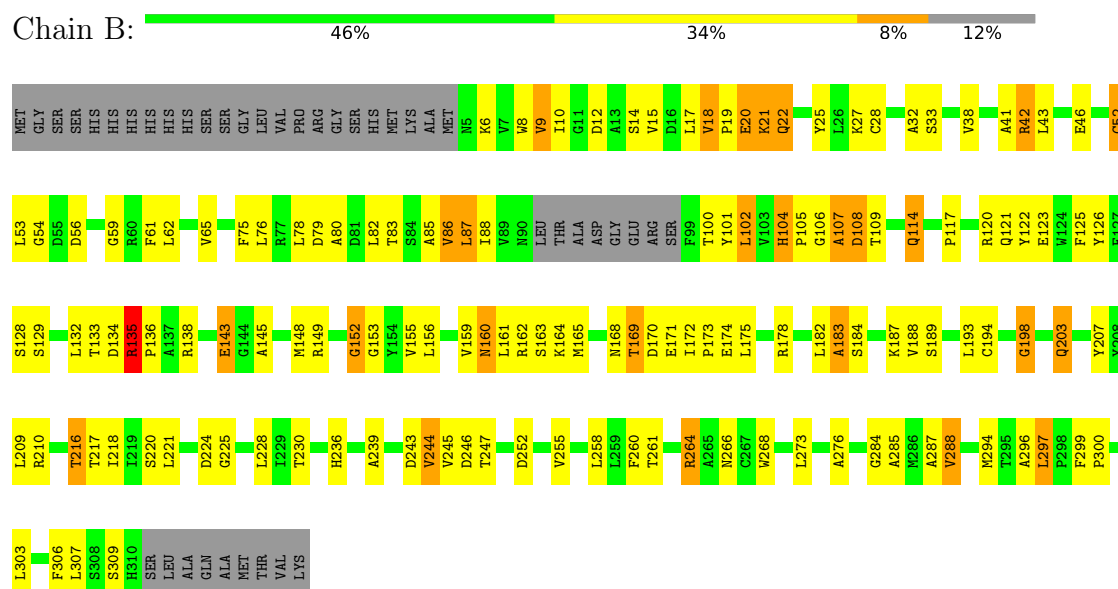
3 Residue-property plots

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: putative sugar kinase



- Molecule 1: putative sugar kinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	135.85Å 54.10Å 89.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.77 – 2.90 28.77 – 2.90	Depositor EDS
% Data completeness (in resolution range)	91.1 (28.77-2.90) 86.9 (28.77-2.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.00 (at 2.69Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.205 , 0.275 0.218 , 0.280	Depositor DCC
R_{free} test set	3490 reflections (9.89%)	wwPDB-VP
Wilson B-factor (Å ²)	59.7	Xtriage
Anisotropy	0.442	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 50.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4485	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.54% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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: K, AIS

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	0.48	0/2264	1.02	20/3077 (0.6%)
1	B	0.49	0/2247	1.01	14/3055 (0.5%)
All	All	0.48	0/4511	1.01	34/6132 (0.6%)

There are no bond length outliers.

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	184	SER	N-CA-C	-8.34	103.63	113.88
1	B	288	VAL	N-CA-C	6.98	116.98	110.42
1	B	135	ARG	CA-C-N	-6.68	112.89	119.64
1	B	135	ARG	C-N-CA	-6.68	112.89	119.64
1	A	28	CYS	CA-C-N	6.61	126.65	119.90
1	A	28	CYS	C-N-CA	6.61	126.65	119.90
1	B	108	ASP	N-CA-C	-6.43	105.42	113.20
1	A	299	PHE	N-CA-C	-6.35	100.84	110.50
1	A	161	LEU	N-CA-C	6.26	119.00	108.23
1	A	143	GLU	N-CA-C	-6.09	104.56	111.14
1	A	202	TRP	N-CA-C	5.98	117.80	111.28
1	A	233	GLY	N-CA-C	5.91	118.43	111.63
1	B	104	HIS	CA-C-N	-5.78	112.61	119.84
1	B	104	HIS	C-N-CA	-5.78	112.61	119.84
1	B	184	SER	N-CA-C	-5.78	105.89	113.12
1	A	18	VAL	N-CA-C	5.71	113.20	107.55
1	A	297	LEU	CA-C-N	5.67	125.57	119.85
1	A	297	LEU	C-N-CA	5.67	125.57	119.85
1	B	152	GLY	N-CA-C	-5.61	105.89	115.08
1	A	179	SER	N-CA-C	-5.57	105.12	111.14
1	A	134	ASP	N-CA-C	5.44	118.85	110.42
1	B	143	GLU	N-CA-C	-5.42	105.28	111.14

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	190	ALA	N-CA-C	-5.36	105.34	111.07
1	A	298	PRO	N-CA-C	5.32	119.83	111.38
1	A	213	GLY	N-CA-C	5.26	120.95	114.69
1	B	12	ASP	N-CA-C	5.20	118.42	110.20
1	A	223	ALA	N-CA-C	-5.19	107.11	112.93
1	A	244	VAL	N-CA-C	5.10	119.95	109.34
1	B	198	GLY	N-CA-C	-5.10	108.12	113.58
1	A	201	HIS	N-CA-C	-5.07	100.64	108.90
1	A	305	THR	N-CA-C	-5.07	106.20	112.38
1	B	297	LEU	CA-C-N	5.05	125.32	119.92
1	B	297	LEU	C-N-CA	5.05	125.32	119.92
1	B	183	ALA	N-CA-C	5.05	117.04	110.43

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	2221	0	2131	129	0
1	B	2204	0	2098	141	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	15	0	13	3	0
4	A	20	0	0	3	0
4	B	21	0	0	3	0
All	All	4485	0	4242	265	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 31.

All (265) 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:203:GLN:NE2	1:A:203:GLN:H	1.43	1.14

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:35:ASN:HD22	1:A:294:MSE:HE2	1.15	1.04
1:A:104:HIS:HB2	1:A:105:PRO:HD3	1.36	1.04
1:A:194:CYS:HA	1:A:199:ALA:HB3	1.45	0.97
1:A:203:GLN:H	1:A:203:GLN:HE21	1.05	0.94
1:B:21:LYS:HD3	1:B:22:GLN:N	1.81	0.94
1:A:160:ASN:ND2	1:A:162:ARG:HH12	1.66	0.91
1:A:156:LEU:HG	4:A:618:HOH:O	1.70	0.91
1:B:294:MSE:HA	1:B:297:LEU:HD13	1.53	0.90
1:A:203:GLN:HE21	1:A:203:GLN:N	1.71	0.89
1:B:203:GLN:HE21	1:B:203:GLN:H	0.88	0.88
1:B:156:LEU:HG	4:B:622:HOH:O	1.74	0.88
1:B:203:GLN:HE21	1:B:203:GLN:N	1.72	0.87
1:A:104:HIS:HB2	1:A:105:PRO:CD	2.08	0.82
1:A:303:LEU:O	1:A:307:LEU:HG	1.79	0.82
1:B:203:GLN:H	1:B:203:GLN:NE2	1.74	0.80
1:B:102:LEU:N	1:B:102:LEU:HD12	1.97	0.79
1:A:35:ASN:HD22	1:A:294:MSE:CE	1.95	0.78
1:A:104:HIS:HD1	1:A:105:PRO:HD2	1.48	0.78
1:B:264:ARG:HE	1:B:264:ARG:HA	1.48	0.78
1:A:35:ASN:ND2	1:A:294:MSE:HE2	1.96	0.78
1:A:203:GLN:NE2	1:A:203:GLN:N	2.26	0.74
1:A:285:ALA:O	1:A:288:VAL:HG22	1.88	0.74
1:A:35:ASN:OD1	1:A:293:ALA:HB1	1.89	0.72
1:B:207:TYR:HB2	4:B:611:HOH:O	1.90	0.71
1:B:83:THR:O	1:B:107:ALA:HB2	1.92	0.70
1:B:162:ARG:HG2	1:B:164:LYS:HZ1	1.57	0.69
1:A:264:ARG:HE	1:A:264:ARG:HA	1.57	0.69
1:A:90:ASN:HB2	1:A:98:SER:OG	1.93	0.69
1:A:175:LEU:O	1:A:179:SER:HB2	1.93	0.69
1:B:244:VAL:HA	1:B:288:VAL:CG1	2.24	0.68
1:B:21:LYS:HD3	1:B:22:GLN:H	1.57	0.68
1:B:244:VAL:HA	1:B:288:VAL:HG11	1.75	0.68
1:A:21:LYS:O	1:A:23:ASN:N	2.28	0.67
1:B:104:HIS:C	1:B:106:GLY:H	2.03	0.67
1:B:120:ARG:O	1:B:123:GLU:HB2	1.95	0.66
1:B:164:LYS:H	1:B:164:LYS:HD3	1.60	0.66
1:B:228:LEU:HD11	1:B:230:THR:HG23	1.77	0.66
1:B:145:ALA:HB3	1:B:182:LEU:HD13	1.76	0.66
1:A:161:LEU:HD13	1:A:163:SER:H	1.59	0.66
1:A:229:ILE:HG12	1:A:234:GLU:HG3	1.77	0.65
1:B:216:THR:HB	1:B:230:THR:HG22	1.77	0.65

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:109:THR:O	1:B:136:PRO:HB2	1.96	0.64
1:A:286:MSE:O	1:A:289:THR:HG23	1.99	0.63
1:B:294:MSE:HG3	1:B:297:LEU:HD22	1.80	0.63
1:B:216:THR:HG21	1:B:268:TRP:HH2	1.64	0.63
1:B:247:THR:HA	1:B:288:VAL:HG22	1.81	0.63
1:B:104:HIS:C	1:B:106:GLY:N	2.56	0.63
1:A:104:HIS:CB	1:A:105:PRO:HD3	2.21	0.62
1:A:42:ARG:NH1	1:A:297:LEU:HB2	2.14	0.62
1:B:172:ILE:HB	1:B:173:PRO:HD3	1.82	0.62
1:A:261:THR:HG21	1:A:276:ALA:HA	1.82	0.61
1:A:32:ALA:HA	1:A:252:ASP:OD2	2.00	0.61
1:A:169:THR:O	1:A:169:THR:HG22	1.99	0.61
1:B:132:LEU:HD22	1:B:138:ARG:HG3	1.83	0.61
1:B:268:TRP:H	1:B:268:TRP:CD1	2.19	0.61
1:A:104:HIS:CB	1:A:105:PRO:CD	2.78	0.60
1:B:14:SER:C	1:B:62:LEU:HD11	2.26	0.60
1:B:209:LEU:H	1:B:209:LEU:HD22	1.68	0.59
1:B:65:VAL:HG11	1:B:294:MSE:HE1	1.83	0.59
1:A:135:ARG:CB	1:A:136:PRO:HD3	2.32	0.59
1:A:246:ASP:HB3	1:A:290:ALA:O	2.03	0.59
1:B:224:ASP:O	1:B:236:HIS:HE1	1.84	0.59
1:B:75:PHE:CZ	1:B:117:PRO:HB3	2.38	0.58
1:B:299:PHE:HB3	1:B:300:PRO:HD2	1.85	0.58
1:A:158:ASP:HB2	4:A:618:HOH:O	2.03	0.58
1:A:104:HIS:HD1	1:A:105:PRO:CD	2.16	0.58
1:A:199:ALA:HB2	4:A:619:HOH:O	2.03	0.58
1:A:196:LEU:HB3	1:A:208:TYR:CE2	2.38	0.58
1:A:162:ARG:HE	3:A:402:AIS:H2	1.68	0.58
1:B:32:ALA:HA	1:B:252:ASP:OD2	2.03	0.58
1:B:42:ARG:O	1:B:43:LEU:HD12	2.03	0.58
1:B:87:LEU:HB3	1:B:100:THR:O	2.05	0.57
1:B:53:LEU:HD12	1:B:76:LEU:HD11	1.86	0.57
1:B:161:LEU:O	1:B:162:ARG:HG3	2.04	0.57
1:A:35:ASN:HA	1:A:294:MSE:HE2	1.87	0.57
1:B:187:LYS:HE2	1:B:220:SER:HB2	1.86	0.57
1:A:225:GLY:HA3	1:A:237:PHE:O	2.04	0.56
1:A:210:ARG:C	1:A:212:LEU:H	2.12	0.56
1:A:148:MSE:SE	1:A:153:GLY:HA3	2.55	0.56
1:B:102:LEU:N	1:B:102:LEU:CD1	2.68	0.56
1:B:123:GLU:HB3	1:B:148:MSE:HE1	1.86	0.56
1:A:75:PHE:CE2	1:A:117:PRO:HB3	2.41	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:14:SER:HB2	1:B:85:ALA:O	2.05	0.56
1:A:240:PRO:HB3	1:A:310:HIS:CD2	2.41	0.56
1:A:218:ILE:HD12	1:A:258:LEU:HD21	1.88	0.55
1:B:17:LEU:HB3	1:B:25:TYR:HB3	1.88	0.55
1:B:108:ASP:OD1	1:B:108:ASP:N	2.38	0.55
1:B:41:ALA:C	1:B:43:LEU:H	2.14	0.55
1:A:55:ASP:HB3	1:A:80:ALA:HB2	1.87	0.55
1:B:75:PHE:CE2	1:B:117:PRO:HB3	2.41	0.55
1:B:264:ARG:HA	1:B:264:ARG:NE	2.19	0.55
1:A:101:TYR:HE1	3:A:402:AIS:N3	2.04	0.55
1:A:161:LEU:CD1	1:A:163:SER:H	2.20	0.55
1:B:261:THR:HG21	1:B:276:ALA:HA	1.89	0.55
1:A:228:LEU:HD11	1:A:273:LEU:HD13	1.89	0.54
1:B:169:THR:HG22	1:B:169:THR:O	2.07	0.54
1:B:123:GLU:HB3	1:B:148:MSE:CE	2.38	0.54
1:B:169:THR:HG23	1:B:172:ILE:HD12	1.90	0.54
1:B:65:VAL:CG1	1:B:294:MSE:HE1	2.38	0.54
1:B:285:ALA:O	1:B:288:VAL:N	2.41	0.54
1:A:224:ASP:O	1:A:238:PRO:HA	2.08	0.53
1:B:188:VAL:HG11	1:B:193:LEU:HD21	1.90	0.53
1:B:10:ILE:HG21	1:B:125:PHE:CZ	2.43	0.53
1:B:38:VAL:HG21	1:B:294:MSE:SE	2.59	0.53
1:A:54:GLY:C	1:A:56:ASP:H	2.15	0.53
1:A:145:ALA:HB3	1:A:182:LEU:HD13	1.90	0.53
1:A:197:SER:O	1:A:198:GLY:C	2.52	0.53
1:B:168:ASN:C	1:B:170:ASP:H	2.17	0.53
1:A:64:GLN:HA	1:A:64:GLN:NE2	2.25	0.52
1:B:101:TYR:C	1:B:102:LEU:HD12	2.34	0.52
1:A:126:TYR:CD1	1:A:126:TYR:C	2.88	0.52
1:A:264:ARG:HA	1:A:264:ARG:NE	2.24	0.52
1:B:46:GLU:HG2	4:B:635:HOH:O	2.09	0.52
1:A:35:ASN:ND2	1:A:294:MSE:CE	2.65	0.52
1:A:38:VAL:HG21	1:A:294:MSE:SE	2.60	0.52
1:A:261:THR:HG21	1:A:276:ALA:CA	2.39	0.52
1:B:42:ARG:O	1:B:42:ARG:HG2	2.10	0.52
1:A:168:ASN:C	1:A:170:ASP:H	2.17	0.52
1:B:18:VAL:HG23	1:B:28:CYS:SG	2.50	0.52
1:B:148:MSE:SE	1:B:153:GLY:HA3	2.59	0.52
1:A:42:ARG:HH11	1:A:297:LEU:HB2	1.75	0.52
1:B:126:TYR:CD1	1:B:126:TYR:C	2.87	0.52
1:B:194:CYS:O	1:B:198:GLY:N	2.43	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:79:ASP:OD1	1:B:80:ALA:N	2.43	0.52
1:B:155:VAL:HB	1:B:183:ALA:HA	1.91	0.52
1:A:9:VAL:HG11	1:A:33:SER:O	2.10	0.51
1:A:193:LEU:O	1:A:196:LEU:HB2	2.10	0.51
1:B:15:VAL:HG22	1:B:62:LEU:HG	1.91	0.51
1:B:18:VAL:CG2	1:B:28:CYS:SG	2.99	0.51
1:B:135:ARG:HH11	1:B:138:ARG:HH11	1.56	0.51
1:B:162:ARG:CG	1:B:164:LYS:HZ1	2.23	0.51
1:A:10:ILE:HG22	1:A:126:TYR:O	2.11	0.51
1:B:86:VAL:C	1:B:87:LEU:HD23	2.35	0.51
1:B:125:PHE:HB2	1:B:148:MSE:HG2	1.92	0.51
1:B:216:THR:HG21	1:B:268:TRP:CH2	2.46	0.51
1:A:89:VAL:HG22	1:A:99:PHE:HE1	1.76	0.51
1:A:208:TYR:O	1:A:212:LEU:HD12	2.10	0.51
1:B:210:ARG:HG3	1:B:210:ARG:HH11	1.75	0.51
1:A:24:SER:N	1:B:27:LYS:HD3	2.26	0.50
1:A:79:ASP:C	1:A:81:ASP:H	2.18	0.50
1:A:53:LEU:O	1:A:78:LEU:HA	2.11	0.50
1:B:21:LYS:HD3	1:B:21:LYS:C	2.32	0.50
1:A:160:ASN:ND2	1:A:162:ARG:NH1	2.47	0.50
1:A:121:GLN:HG2	1:A:122:TYR:CD2	2.47	0.50
1:A:221:LEU:O	1:A:222:GLY:C	2.55	0.50
1:A:162:ARG:O	1:A:163:SER:C	2.54	0.49
1:A:190:ALA:HB2	1:A:221:LEU:HD23	1.94	0.49
1:B:88:ILE:HG13	1:B:88:ILE:O	2.12	0.49
1:B:228:LEU:HD11	1:B:230:THR:CG2	2.42	0.49
1:B:126:TYR:CE2	1:B:255:VAL:HG11	2.47	0.49
1:A:108:ASP:O	1:A:111:VAL:HG23	2.13	0.49
1:B:162:ARG:HG2	1:B:164:LYS:NZ	2.27	0.49
1:B:221:LEU:HB2	1:B:225:GLY:O	2.12	0.49
1:A:225:GLY:HA2	1:A:281:ASN:ND2	2.28	0.49
1:A:87:LEU:HD21	3:A:402:AIS:O2'	2.13	0.49
1:A:269:ASP:OD1	1:A:271:ALA:HB3	2.13	0.49
1:B:210:ARG:HG3	1:B:210:ARG:NH1	2.28	0.49
1:B:284:GLY:O	1:B:287:ALA:HB3	2.12	0.48
1:A:125:PHE:HB2	1:A:148:MSE:HG2	1.95	0.48
1:A:113:PRO:O	1:A:116:LEU:HB2	2.13	0.48
1:B:225:GLY:HA2	1:B:239:ALA:HB2	1.94	0.48
1:B:287:ALA:HA	1:B:296:ALA:HB1	1.94	0.48
1:B:121:GLN:HG2	1:B:122:TYR:CD1	2.49	0.48
1:A:104:HIS:CD2	1:A:104:HIS:H	2.24	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:9:VAL:HG13	1:B:33:SER:HB3	1.95	0.48
1:B:135:ARG:N	1:B:136:PRO:HD3	2.29	0.48
1:B:135:ARG:NH1	1:B:138:ARG:HH11	2.12	0.48
1:B:10:ILE:HG22	1:B:126:TYR:O	2.14	0.47
1:B:86:VAL:O	1:B:102:LEU:HD13	2.14	0.47
1:B:210:ARG:HD2	1:B:210:ARG:C	2.38	0.47
1:A:135:ARG:O	1:A:136:PRO:C	2.57	0.47
1:B:52:CYS:O	1:B:108:ASP:HA	2.14	0.47
1:A:143:GLU:OE1	1:A:143:GLU:HA	2.15	0.47
1:A:175:LEU:HD22	1:A:175:LEU:H	1.80	0.47
1:A:210:ARG:HA	1:A:214:CYS:O	2.15	0.46
1:A:87:LEU:C	1:A:87:LEU:HD12	2.41	0.46
1:A:188:VAL:HG22	1:A:189:SER:N	2.30	0.46
1:A:188:VAL:O	1:A:219:ILE:HA	2.16	0.46
1:B:6:LYS:HD3	1:B:8:TRP:CZ2	2.50	0.46
1:A:121:GLN:O	1:A:122:TYR:HB2	2.14	0.45
1:A:225:GLY:HA2	1:A:281:ASN:HD21	1.81	0.45
1:B:307:LEU:C	1:B:309:SER:H	2.24	0.45
1:A:135:ARG:CB	1:A:136:PRO:CD	2.93	0.45
1:B:56:ASP:OD1	1:B:56:ASP:N	2.44	0.45
1:B:261:THR:HG21	1:B:276:ALA:CA	2.46	0.45
1:A:124:TRP:CD2	1:A:154:TYR:HB2	2.52	0.45
1:B:104:HIS:O	1:B:106:GLY:N	2.50	0.45
1:B:209:LEU:HD22	1:B:209:LEU:N	2.30	0.45
1:B:168:ASN:ND2	1:B:170:ASP:HB2	2.31	0.45
1:A:306:PHE:CD1	1:A:306:PHE:C	2.95	0.45
1:B:128:SER:HB2	1:B:160:ASN:HB2	1.99	0.45
1:B:133:THR:HG22	1:B:134:ASP:N	2.32	0.45
1:A:14:SER:O	1:A:62:LEU:HD21	2.16	0.44
1:B:54:GLY:O	1:B:59:GLY:HA3	2.18	0.44
1:B:121:GLN:HG3	1:B:152:GLY:HA3	1.99	0.44
1:B:134:ASP:C	1:B:136:PRO:CD	2.90	0.44
1:A:55:ASP:HB3	1:A:80:ALA:CB	2.46	0.44
1:A:172:ILE:HB	1:A:173:PRO:HD3	2.00	0.44
1:A:210:ARG:C	1:A:212:LEU:N	2.75	0.44
1:A:242:VAL:HG12	1:A:285:ALA:HB1	1.98	0.44
1:B:114:GLN:OE1	1:B:114:GLN:N	2.51	0.44
1:A:103:VAL:HG22	1:B:102:LEU:HB3	1.99	0.44
1:B:120:ARG:CB	1:B:123:GLU:OE2	2.66	0.44
1:B:228:LEU:HD22	1:B:273:LEU:HD22	1.99	0.44
1:B:243:ASP:O	1:B:244:VAL:CB	2.66	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:132:LEU:HD13	1:A:175:LEU:HB3	1.99	0.44
1:A:286:MSE:HG2	1:A:306:PHE:CD2	2.53	0.44
1:B:168:ASN:C	1:B:170:ASP:N	2.75	0.43
1:B:303:LEU:O	1:B:307:LEU:HG	2.18	0.43
1:A:82:LEU:HB3	1:A:107:ALA:HB2	2.00	0.43
1:B:178:ARG:HD3	1:B:178:ARG:HA	1.81	0.43
1:B:122:TYR:CD2	1:B:266:ASN:HB3	2.53	0.43
1:B:135:ARG:HH11	1:B:138:ARG:NH1	2.17	0.43
1:A:102:LEU:HD13	1:B:86:VAL:HG11	2.01	0.43
1:B:82:LEU:HD12	1:B:82:LEU:N	2.34	0.43
1:A:53:LEU:HD12	1:A:76:LEU:HD11	1.99	0.43
1:A:109:THR:HB	1:A:134:ASP:HB2	2.01	0.43
1:A:190:ALA:HB1	1:A:202:TRP:CD1	2.53	0.43
1:B:174:GLU:OE1	1:B:174:GLU:HA	2.19	0.43
1:B:171:GLU:HB3	1:B:175:LEU:HD21	2.01	0.43
1:B:20:GLU:O	1:B:21:LYS:C	2.62	0.42
1:B:61:PHE:O	1:B:65:VAL:HG23	2.19	0.42
1:A:223:ALA:O	1:A:239:ALA:HB3	2.18	0.42
1:A:5:ASN:OD1	1:A:122:TYR:HB3	2.19	0.42
1:B:19:PRO:HG3	1:B:25:TYR:HE1	1.85	0.42
1:A:168:ASN:HD21	1:A:170:ASP:CG	2.27	0.42
1:A:207:TYR:HB3	1:A:211:ASP:OD2	2.20	0.42
1:A:162:ARG:O	1:A:165:MSE:HG2	2.20	0.42
1:A:286:MSE:HG2	1:A:306:PHE:CE2	2.54	0.42
1:A:75:PHE:CZ	1:A:117:PRO:HB3	2.55	0.42
1:A:55:ASP:HB3	1:A:80:ALA:CA	2.49	0.41
1:B:135:ARG:N	1:B:136:PRO:CD	2.82	0.41
1:A:156:LEU:HD21	1:A:255:VAL:HG13	2.02	0.41
1:A:306:PHE:O	1:A:307:LEU:C	2.63	0.41
1:A:294:MSE:HA	1:A:297:LEU:HG	2.01	0.41
1:A:43:LEU:O	1:A:260:PHE:CE1	2.73	0.41
1:B:261:THR:O	1:B:264:ARG:HB2	2.20	0.41
1:A:42:ARG:C	1:A:44:GLY:H	2.29	0.41
1:A:86:VAL:HB	1:B:102:LEU:HD23	2.02	0.41
1:B:9:VAL:HG11	1:B:33:SER:O	2.21	0.41
1:A:19:PRO:HG3	1:A:25:TYR:HE1	1.84	0.41
1:B:101:TYR:HB2	1:B:104:HIS:CE1	2.55	0.41
1:A:168:ASN:C	1:A:168:ASN:ND2	2.79	0.41
1:A:186:CYS:HB3	1:A:209:LEU:HD13	2.03	0.41
1:B:101:TYR:CD1	1:B:101:TYR:N	2.88	0.41
1:B:162:ARG:HA	1:B:164:LYS:HZ3	1.85	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:306:PHE:O	1:B:309:SER:HB3	2.19	0.41
1:B:188:VAL:HG22	1:B:189:SER:N	2.35	0.41
1:B:246:ASP:O	1:B:288:VAL:HA	2.21	0.41
1:A:54:GLY:O	1:A:59:GLY:HA3	2.20	0.41
1:A:104:HIS:CD2	1:B:105:PRO:HD3	2.56	0.41
1:A:104:HIS:ND1	1:A:105:PRO:HD2	2.24	0.41
1:A:163:SER:O	1:A:164:LYS:C	2.63	0.41
1:A:273:LEU:O	1:A:277:ILE:HG13	2.20	0.41
1:B:143:GLU:OE1	1:B:143:GLU:HA	2.20	0.41
1:B:260:PHE:O	1:B:264:ARG:HD2	2.21	0.41
1:A:168:ASN:C	1:A:170:ASP:N	2.78	0.41
1:B:20:GLU:HG3	1:B:21:LYS:H	1.86	0.41
1:A:168:ASN:O	1:A:170:ASP:N	2.54	0.40
1:A:269:ASP:O	1:A:270:HIS:C	2.65	0.40
1:B:161:LEU:O	1:B:162:ARG:CG	2.68	0.40
1:B:100:THR:HG22	1:B:101:TYR:N	2.36	0.40
1:B:207:TYR:CD1	1:B:210:ARG:NH2	2.89	0.40
1:B:218:ILE:HD12	1:B:258:LEU:HD21	2.02	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	295/339 (87%)	241 (82%)	36 (12%)	18 (6%)	1	4
1	B	294/339 (87%)	260 (88%)	25 (8%)	9 (3%)	3	14
All	All	589/678 (87%)	501 (85%)	61 (10%)	27 (5%)	2	7

All (27) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	22	GLN
1	A	23	ASN
1	A	104	HIS
1	A	135	ARG
1	A	196	LEU
1	A	198	GLY
1	B	20	GLU
1	B	22	GLN
1	B	245	VAL
1	A	169	THR
1	A	222	GLY
1	B	244	VAL
1	A	244	VAL
1	A	248	THR
1	A	287	ALA
1	B	42	ARG
1	B	107	ALA
1	A	102	LEU
1	B	135	ARG
1	B	169	THR
1	A	80	ALA
1	A	163	SER
1	A	164	LYS
1	B	163	SER
1	A	136	PRO
1	A	309	SER
1	A	106	GLY

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	226/263 (86%)	209 (92%)	17 (8%)	12	37
1	B	222/263 (84%)	204 (92%)	18 (8%)	11	33
All	All	448/526 (85%)	413 (92%)	35 (8%)	11	35

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	VAL
1	A	28	CYS
1	A	83	THR
1	A	104	HIS
1	A	111	VAL
1	A	116	LEU
1	A	129	SER
1	A	149	ARG
1	A	159	VAL
1	A	160	ASN
1	A	161	LEU
1	A	203	GLN
1	A	215	ASP
1	A	217	THR
1	A	243	ASP
1	A	264	ARG
1	A	281	ASN
1	B	9	VAL
1	B	18	VAL
1	B	21	LYS
1	B	52	CYS
1	B	78	LEU
1	B	86	VAL
1	B	87	LEU
1	B	102	LEU
1	B	114	GLN
1	B	129	SER
1	B	149	ARG
1	B	159	VAL
1	B	160	ASN
1	B	165	MSE
1	B	203	GLN
1	B	216	THR
1	B	217	THR
1	B	264	ARG

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	5	ASN
1	A	35	ASN
1	A	64	GLN
1	A	67	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	90	ASN
1	A	160	ASN
1	A	168	ASN
1	A	195	GLN
1	A	203	GLN
1	A	281	ASN
1	A	310	HIS
1	B	5	ASN
1	B	35	ASN
1	B	64	GLN
1	B	67	GLN
1	B	104	HIS
1	B	121	GLN
1	B	160	ASN
1	B	168	ASN
1	B	195	GLN
1	B	203	GLN
1	B	236	HIS
1	B	266	ASN
1	B	281	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 4 are monoatomic - leaving 1 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond

length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	AIS	A	402	-	15,16,16	1.05	2 (13%)	21,23,23	2.12	9 (42%)

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
3	AIS	A	402	-	-	0/6/22/22	0/2/2/2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	402	AIS	C4-C5	-2.38	1.33	1.36
3	A	402	AIS	C4-N3	-2.36	1.33	1.37

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	402	AIS	C4-C5-N6	-5.23	124.60	131.28
3	A	402	AIS	C4'-O4'-C1'	-3.46	101.82	109.47
3	A	402	AIS	N1-C2-N3	-3.28	108.57	111.68
3	A	402	AIS	C4-C5-N1	3.22	109.99	107.48
3	A	402	AIS	C5-C4-N3	-2.80	106.49	109.66
3	A	402	AIS	O4'-C1'-C2'	-2.64	100.97	106.62
3	A	402	AIS	C4-N3-C2	2.53	108.55	105.24
3	A	402	AIS	O4'-C1'-N1	2.46	113.93	108.36
3	A	402	AIS	C5'-C4'-C3'	-2.04	110.29	115.10

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	402	AIS	3	0

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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, 95th 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 ²)	Q<0.9
1	A	295/339 (87%)	-0.11	1 (0%) 90 87	30, 52, 75, 99	0
1	B	294/339 (86%)	-0.08	0 100 100	32, 53, 75, 91	0
All	All	589/678 (86%)	-0.10	1 (0%) 91 89	30, 52, 75, 99	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	242	VAL	2.8

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

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, 95th 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(Å ²)	Q<0.9
3	AIS	A	402	15/15	0.83	0.14	75,75,75,75	0
2	K	B	406	1/1	0.86	0.07	64,64,64,64	0
2	K	A	405	1/1	0.86	0.07	69,69,69,69	0
2	K	B	407	1/1	0.92	0.05	52,52,52,52	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	K	A	404	1/1	0.99	0.06	52,52,52,52	0

6.5 Other polymers [i](#)

There are no such residues in this entry.