



Full wwPDB X-ray Structure Validation Report ⓘ

Jul 25, 2017 – 07:56 PM EDT

PDB ID : 2PQJ
Title : Crystal structure of active ribosome inactivating protein from maize (b-32), complex with adenine
Authors : Mak, A.N.S.; Au, S.W.N.; Cha, S.S.; Young, J.A.; Wong, K.B.; Shaw, P.C.
Deposited on : unknown
Resolution : 2.80 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20029824
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20029824

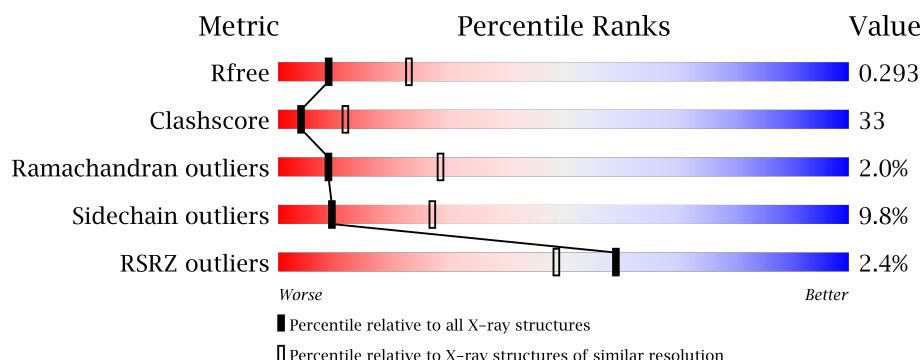
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	100719	2583 (2.80-2.80)
Clashscore	112137	3033 (2.80-2.80)
Ramachandran outliers	110173	2983 (2.80-2.80)
Sidechain outliers	110143	2985 (2.80-2.80)
RSRZ outliers	101464	2610 (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 ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	243	<div> <div>6%</div> <div>47% 42% 6% . .</div> </div>
1	B	243	<div> <div>47% 44% 5% 5%</div> </div>
1	C	243	<div> <div>6%</div> <div>47% 41% 7% 5%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ADE	A	501	-	-	-	X
2	ADE	B	501	-	-	X	X
2	ADE	C	501	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5427 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.

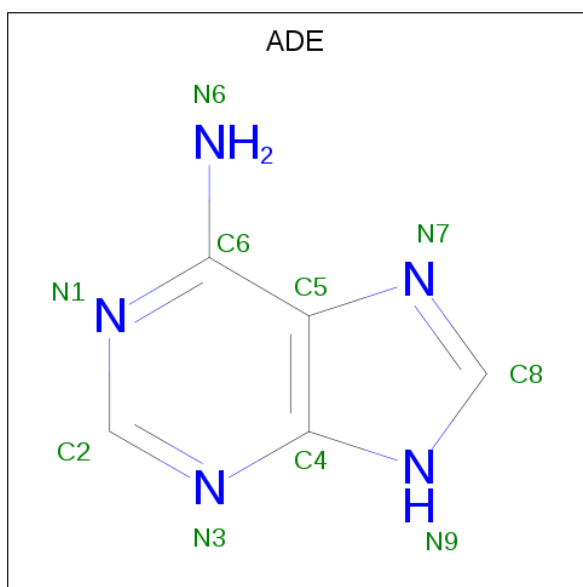
- Molecule 1 is a protein called Ribosome-inactivating protein 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	233	Total	C	N	O	S	0	0	0
			1800	1153	313	326	8			
1	B	232	Total	C	N	O	S	0	0	0
			1786	1144	306	328	8			
1	C	232	Total	C	N	O	S	0	0	0
			1796	1151	312	326	7			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	21	MET	-	EXPRESSION TAG	UNP P25891
A	165	LEU	-	LINKER	UNP P25891
A	166	GLU	-	LINKER	UNP P25891
B	21	MET	-	EXPRESSION TAG	UNP P25891
B	165	LEU	-	LINKER	UNP P25891
B	166	GLU	-	LINKER	UNP P25891
C	21	MET	-	EXPRESSION TAG	UNP P25891
C	165	LEU	-	LINKER	UNP P25891
C	166	GLU	-	LINKER	UNP P25891

- Molecule 2 is ADENINE (three-letter code: ADE) (formula: C₅H₅N₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	N	0	0
			10	5	5		
2	B	1	Total	C	N	0	0
			10	5	5		
2	C	1	Total	C	N	0	0
			10	5	5		

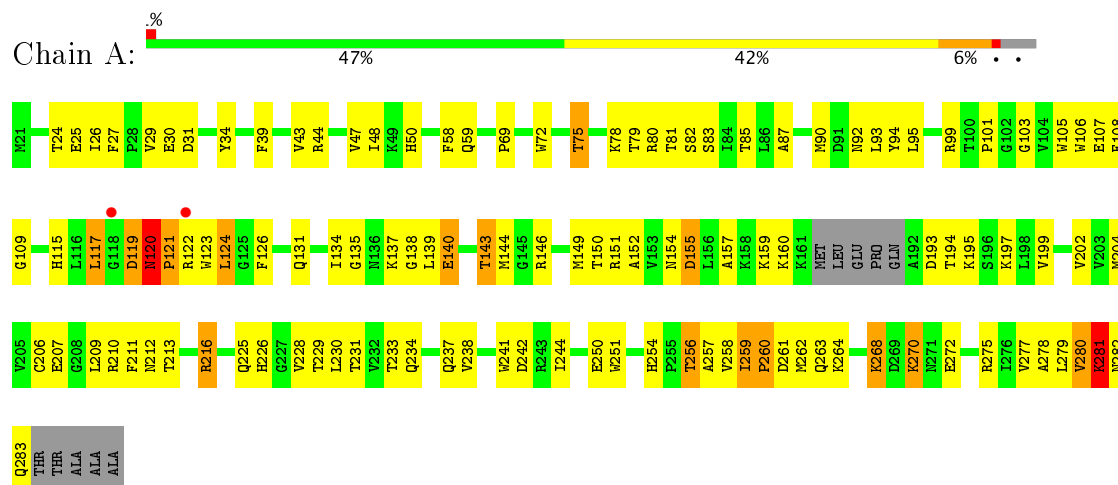
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	6	Total	O	0	0
			6	6		
3	B	6	Total	O	0	0
			6	6		
3	C	3	Total	O	0	0
			3	3		

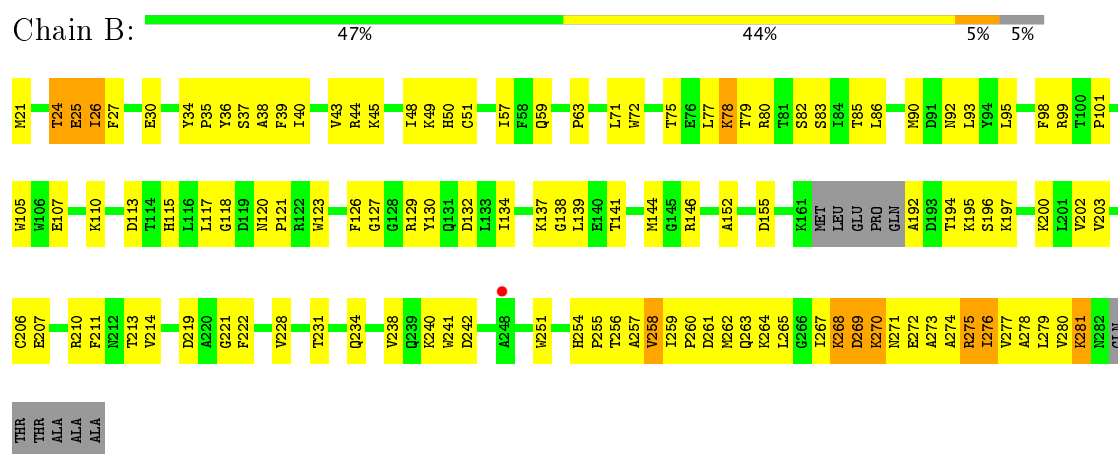
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

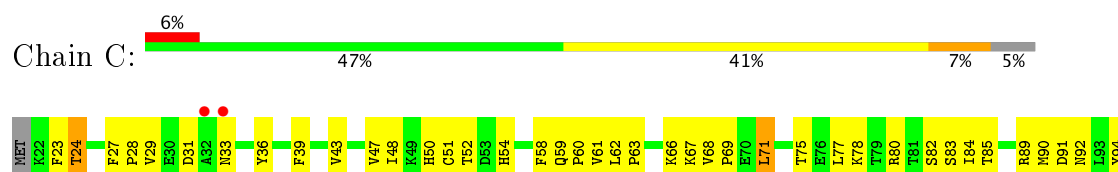
• Molecule 1: Ribosome-inactivating protein 3

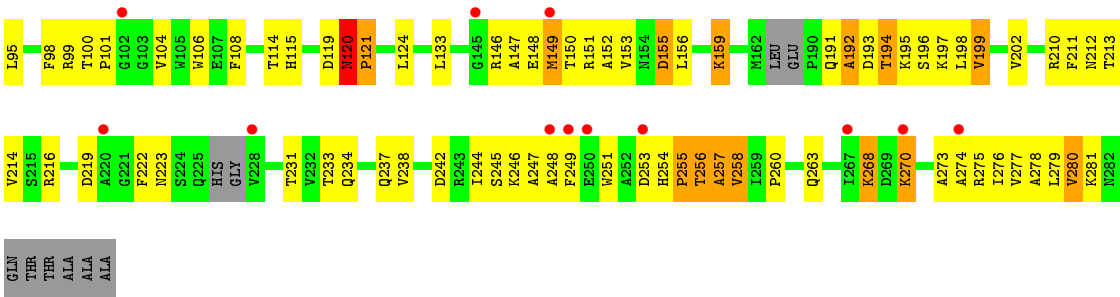


• Molecule 1: Ribosome-inactivating protein 3



• Molecule 1: Ribosome-inactivating protein 3





GLN
THR
THR
ALA
ALA
ALA

4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	115.91Å 115.91Å 44.90Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	40.99 – 2.80 40.99 – 2.80	Depositor EDS
% Data completeness (in resolution range)	98.9 (40.99-2.80) 100.0 (40.99-2.80)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.78 (at 2.81Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.232 , 0.293 0.238 , 0.293	Depositor DCC
R_{free} test set	843 reflections (5.07%)	DCC
Wilson B-factor (Å ²)	63.8	Xtriage
Anisotropy	0.154	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 72.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.028 for -h,-k,l 0.025 for h,-h-k,-l 0.021 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	5427	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.89% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.44	0/1840	0.68	1/2495 (0.0%)
1	B	0.46	0/1826	0.68	0/2480
1	C	0.40	0/1834	0.63	0/2484
All	All	0.43	0/5500	0.66	1/7459 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
1	C	0	2
All	All	0	5

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	120	ASN	C-N-CD	-5.32	108.90	120.60

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	117	LEU	Peptide
1	A	281	LYS	Peptide
1	A	282	ASN	Peptide
1	C	120	ASN	Peptide

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Mol	Chain	Res	Type	Group
1	C	192	ALA	Peptide

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	1800	0	1781	116	0
1	B	1786	0	1753	120	0
1	C	1796	0	1783	127	0
2	A	10	0	4	2	0
2	B	10	0	4	5	0
2	C	10	0	4	2	0
3	A	6	0	0	0	0
3	B	6	0	0	0	0
3	C	3	0	0	0	0
All	All	5427	0	5329	355	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 33.

All (355) 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:59:GLN:HE21	1:A:277:VAL:CG2	1.53	1.21
1:C:152:ALA:HA	1:C:197:LYS:HD2	1.28	1.11
1:C:211:PHE:HB3	1:C:213:THR:HG22	1.37	1.05
1:A:59:GLN:NE2	1:A:277:VAL:CG2	2.24	1.01
1:C:231:THR:H	1:C:234:GLN:HE21	1.12	0.97
1:B:264:LYS:HG3	1:C:254:HIS:CE1	1.99	0.97
1:B:118:GLY:O	1:B:121:PRO:HD3	1.65	0.96
1:A:106:TRP:HZ2	1:A:160:LYS:O	1.49	0.94
1:B:37:SER:HB3	1:B:219:ASP:OD1	1.68	0.94
1:C:212:ASN:O	1:C:216:ARG:HG2	1.69	0.93
1:B:231:THR:H	1:B:234:GLN:HE21	1.10	0.92
1:A:26:ILE:HG21	1:A:78:LYS:HD3	1.53	0.91
1:C:268:LYS:HE2	1:C:268:LYS:N	1.87	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:106:TRP:NE1	1:A:159:LYS:HG2	1.86	0.89
1:A:152:ALA:HA	1:A:197:LYS:HD2	1.57	0.87
1:B:264:LYS:HG3	1:C:254:HIS:NE2	1.88	0.86
1:C:192:ALA:HB1	1:C:195:LYS:HG3	1.55	0.85
1:A:59:GLN:HE21	1:A:277:VAL:HG23	1.37	0.85
1:A:211:PHE:HB3	1:A:213:THR:HG22	1.59	0.85
1:A:106:TRP:CZ2	1:A:160:LYS:O	2.30	0.85
1:B:127:GLY:HA3	1:B:132:ASP:OD2	1.79	0.83
1:B:258:VAL:HG12	1:B:263:GLN:OE1	1.79	0.83
1:B:254:HIS:HD2	1:B:257:ALA:HB2	1.43	0.82
1:C:29:VAL:HG13	1:C:149:MET:HG2	1.62	0.81
1:A:59:GLN:HE21	1:A:277:VAL:HG22	1.43	0.81
1:B:37:SER:CB	1:B:219:ASP:OD1	2.31	0.79
1:C:231:THR:H	1:C:234:GLN:NE2	1.80	0.79
1:C:24:THR:HG22	1:C:75:THR:HA	1.63	0.78
1:B:211:PHE:HB3	1:B:213:THR:HG22	1.64	0.78
1:A:268:LYS:HE2	1:A:268:LYS:N	2.00	0.77
1:A:59:GLN:HB3	1:A:278:ALA:O	1.84	0.77
1:C:268:LYS:HE2	1:C:268:LYS:H	1.48	0.75
1:A:106:TRP:HE1	1:A:159:LYS:HG2	1.51	0.75
1:B:50:HIS:HB2	1:B:90:MET:HE3	1.67	0.75
1:B:24:THR:HB	1:B:75:THR:HG22	1.69	0.75
1:C:52:THR:OG1	1:C:54:HIS:HE1	1.69	0.74
1:A:115:HIS:CE1	1:A:123:TRP:NE1	2.55	0.74
1:C:59:GLN:NE2	1:C:277:VAL:HG23	2.01	0.74
1:A:79:THR:HG21	1:A:154:ASN:HD21	1.50	0.74
1:B:268:LYS:HE2	1:B:268:LYS:N	2.02	0.74
1:B:59:GLN:HB3	1:B:278:ALA:O	1.88	0.74
1:A:59:GLN:HB2	1:A:280:VAL:HG13	1.68	0.73
1:B:192:ALA:HB1	1:B:195:LYS:HG3	1.71	0.73
1:C:51:CYS:SG	1:C:62:LEU:HD23	2.28	0.73
1:A:250:GLU:CD	1:A:259:ILE:HD13	2.09	0.72
1:B:264:LYS:HE3	1:C:254:HIS:CE1	2.24	0.72
1:B:231:THR:H	1:B:234:GLN:NE2	1.86	0.72
1:B:107:GLU:HB2	1:B:117:LEU:HD12	1.72	0.72
1:C:256:THR:O	1:C:257:ALA:O	2.08	0.71
1:B:264:LYS:CG	1:C:254:HIS:CE1	2.71	0.71
1:C:211:PHE:HB2	1:C:214:VAL:HG23	1.70	0.71
1:C:39:PHE:O	1:C:43:VAL:HG23	1.91	0.71
1:A:250:GLU:CD	1:A:259:ILE:CD1	2.59	0.71
1:B:240:LYS:HG3	1:B:265:LEU:HD11	1.73	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:24:THR:HG22	1:B:75:THR:HA	1.72	0.70
1:C:115:HIS:CD2	1:C:121:PRO:HB2	2.27	0.70
1:B:251:TRP:CH2	1:B:262:MET:HG3	2.26	0.70
1:C:62:LEU:HD21	1:C:279:LEU:HD13	1.73	0.70
1:C:194:THR:O	1:C:198:LEU:HG	1.91	0.70
1:B:26:ILE:HG22	1:B:27:PHE:N	2.06	0.70
1:C:133:LEU:HD12	1:C:199:VAL:HG23	1.73	0.70
1:B:24:THR:OG1	1:B:43:VAL:HG22	1.92	0.69
1:B:83:SER:O	1:B:101:PRO:HG3	1.91	0.69
1:A:231:THR:H	1:A:234:GLN:NE2	1.90	0.68
1:B:254:HIS:CD2	1:B:256:THR:OG1	2.46	0.67
1:C:69:PRO:HD3	1:C:89:ARG:NH2	2.09	0.67
1:C:58:PHE:HD2	1:C:274:ALA:CB	2.08	0.67
1:B:259:ILE:O	1:B:262:MET:N	2.27	0.67
1:A:79:THR:HG21	1:A:154:ASN:ND2	2.10	0.66
1:C:254:HIS:O	1:C:256:THR:N	2.28	0.66
1:C:28:PRO:HB2	1:C:31:ASP:HB3	1.79	0.65
1:C:119:ASP:O	1:C:120:ASN:CG	2.34	0.65
1:A:231:THR:H	1:A:234:GLN:HE21	1.44	0.65
1:B:50:HIS:HB2	1:B:90:MET:CE	2.27	0.65
1:B:35:PRO:HG2	1:B:38:ALA:HB2	1.79	0.65
1:C:119:ASP:O	1:C:120:ASN:CB	2.44	0.64
1:A:107:GLU:HG2	1:A:123:TRP:CZ3	2.33	0.64
1:C:59:GLN:HB3	1:C:278:ALA:O	1.98	0.63
1:C:50:HIS:HB2	1:C:90:MET:HE2	1.80	0.63
1:B:130:TYR:CD1	1:B:139:LEU:HD11	2.34	0.63
1:C:71:LEU:HD12	1:C:90:MET:SD	2.38	0.63
1:A:106:TRP:CE3	1:A:122:ARG:O	2.52	0.63
1:B:24:THR:HG21	1:B:75:THR:HB	1.81	0.63
1:A:26:ILE:HG21	1:A:78:LYS:CD	2.27	0.62
1:A:59:GLN:NE2	1:A:277:VAL:HG21	2.12	0.62
1:C:31:ASP:OD1	1:C:33:ASN:HB2	1.99	0.62
1:B:254:HIS:CD2	1:B:257:ALA:HB2	2.32	0.61
1:B:37:SER:N	1:B:219:ASP:OD1	2.33	0.61
1:A:24:THR:HB	1:A:75:THR:HA	1.83	0.60
1:C:52:THR:OG1	1:C:54:HIS:CE1	2.53	0.60
1:B:152:ALA:HA	1:B:197:LYS:HD2	1.83	0.60
1:B:192:ALA:CB	1:B:195:LYS:HG3	2.31	0.60
1:B:24:THR:CB	1:B:75:THR:HG22	2.31	0.60
1:C:59:GLN:HE21	1:C:277:VAL:HG23	1.66	0.60
1:B:272:GLU:O	1:B:275:ARG:HB2	2.01	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:264:LYS:CE	1:C:254:HIS:HE1	2.15	0.59
1:B:207:GLU:HA	1:B:207:GLU:OE1	2.02	0.59
1:C:147:ALA:O	1:C:150:THR:HB	2.02	0.59
1:C:231:THR:N	1:C:234:GLN:HE21	1.92	0.59
1:C:77:LEU:HD22	1:C:149:MET:CE	2.32	0.59
1:A:94:TYR:HA	2:A:501:ADE:C2	2.37	0.59
1:C:58:PHE:CD2	1:C:274:ALA:CB	2.85	0.59
1:A:58:PHE:CD1	1:A:270:LYS:HE2	2.38	0.59
1:C:24:THR:HG21	1:C:75:THR:HB	1.85	0.59
1:C:68:VAL:O	1:C:68:VAL:HG12	2.03	0.59
1:A:115:HIS:CE1	1:A:123:TRP:CE2	2.91	0.58
1:B:130:TYR:CE2	2:B:501:ADE:C6	2.91	0.58
1:B:24:THR:CG2	1:B:75:THR:HA	2.34	0.58
1:A:79:THR:CG2	1:A:154:ASN:HD21	2.15	0.58
1:B:115:HIS:CD2	1:B:121:PRO:HG2	2.39	0.58
1:B:30:GLU:OE2	1:B:79:THR:HB	2.04	0.58
1:B:251:TRP:CE3	1:B:270:LYS:HA	2.39	0.57
1:A:101:PRO:CD	1:A:157:ALA:HB1	2.34	0.57
1:B:254:HIS:O	1:B:256:THR:N	2.37	0.57
1:C:85:THR:O	1:C:98:PHE:HA	2.04	0.57
1:C:260:PRO:HA	1:C:263:GLN:CG	2.34	0.57
1:A:82:SER:OG	1:A:83:SER:N	2.36	0.57
1:B:138:GLY:O	1:B:141:THR:HG23	2.05	0.57
1:C:94:TYR:HA	2:C:501:ADE:C2	2.39	0.57
1:B:264:LYS:HE3	1:C:254:HIS:HE1	1.69	0.56
1:C:24:THR:CG2	1:C:75:THR:HA	2.34	0.56
1:B:35:PRO:HG2	1:B:38:ALA:CB	2.35	0.56
1:B:207:GLU:OE2	2:B:501:ADE:C2	2.58	0.56
1:A:92:ASN:HB3	1:A:281:LYS:HE2	1.87	0.56
1:B:231:THR:N	1:B:234:GLN:HE21	1.91	0.56
1:A:106:TRP:HB3	1:A:124:LEU:HG	1.87	0.56
1:C:50:HIS:HB2	1:C:90:MET:CE	2.36	0.56
1:A:131:GLN:HA	1:A:135:GLY:HA2	1.88	0.55
1:A:260:PRO:O	1:A:264:LYS:HG2	2.06	0.55
1:A:59:GLN:NE2	1:A:277:VAL:HG22	2.11	0.55
1:B:80:ARG:HH11	1:B:80:ARG:HG2	1.71	0.55
1:C:151:ARG:O	1:C:155:ASP:HB2	2.07	0.55
1:B:59:GLN:HB2	1:B:280:VAL:HG23	1.89	0.55
1:C:231:THR:HG23	1:C:234:GLN:NE2	2.22	0.55
1:C:58:PHE:HD2	1:C:274:ALA:HB1	1.72	0.55
1:B:258:VAL:CG1	1:B:263:GLN:OE1	2.53	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:152:ALA:CA	1:C:197:LYS:HD2	2.18	0.55
1:C:69:PRO:HD3	1:C:89:ARG:HH22	1.72	0.54
1:A:149:MET:HB2	1:A:204:MET:CE	2.37	0.54
1:B:26:ILE:CG2	1:B:27:PHE:N	2.68	0.54
1:C:95:LEU:N	2:C:501:ADE:N1	2.52	0.54
1:A:58:PHE:CE1	1:A:270:LYS:HE2	2.43	0.54
1:C:258:VAL:O	1:C:260:PRO:HD3	2.08	0.54
1:B:211:PHE:CZ	1:B:238:VAL:HA	2.43	0.54
1:C:253:ASP:C	1:C:254:HIS:HD1	2.11	0.53
1:C:92:ASN:HB3	1:C:281:LYS:HE2	1.91	0.53
1:C:84:ILE:HG12	1:C:153:VAL:HG13	1.91	0.53
1:C:192:ALA:HB1	1:C:195:LYS:CG	2.36	0.53
1:A:44:ARG:O	1:A:48:ILE:HG13	2.08	0.52
1:B:24:THR:HG21	1:B:75:THR:CB	2.39	0.52
1:A:31:ASP:OD2	1:A:34:TYR:CE2	2.62	0.52
1:A:26:ILE:HG22	1:A:27:PHE:N	2.24	0.52
1:A:241:TRP:CH2	1:A:281:LYS:HG3	2.45	0.52
1:C:149:MET:O	1:C:153:VAL:HG23	2.09	0.52
1:B:110:LYS:O	1:B:113:ASP:HB2	2.10	0.52
1:B:234:GLN:O	1:B:238:VAL:HG23	2.10	0.52
1:B:271:ASN:O	1:B:274:ALA:HB3	2.10	0.52
1:C:90:MET:O	1:C:91:ASP:C	2.48	0.52
1:C:60:PRO:HD2	1:C:279:LEU:HD23	1.92	0.52
1:C:273:ALA:O	1:C:277:VAL:HG22	2.09	0.51
1:C:211:PHE:HB2	1:C:214:VAL:CG2	2.38	0.51
1:C:36:TYR:O	1:C:39:PHE:HB3	2.10	0.51
1:B:231:THR:HG23	1:B:234:GLN:NE2	2.25	0.51
1:A:47:VAL:O	1:A:90:MET:HE1	2.09	0.51
1:C:149:MET:HG3	1:C:149:MET:O	2.10	0.51
1:B:264:LYS:CE	1:C:254:HIS:CE1	2.92	0.51
1:A:250:GLU:OE2	1:A:259:ILE:HD13	2.09	0.51
1:B:78:LYS:HB2	1:B:83:SER:HB3	1.92	0.51
1:C:198:LEU:O	1:C:202:VAL:HG23	2.11	0.51
1:A:211:PHE:CZ	1:A:238:VAL:HA	2.45	0.51
1:C:242:ASP:O	1:C:246:LYS:HG3	2.10	0.50
1:B:144:MET:HB2	1:B:228:VAL:HG13	1.94	0.50
1:A:280:VAL:HG23	1:A:281:LYS:N	2.26	0.50
1:A:39:PHE:O	1:A:43:VAL:HG23	2.11	0.50
1:A:119:ASP:O	1:A:120:ASN:O	2.30	0.50
1:B:130:TYR:CE1	1:B:139:LEU:HD11	2.46	0.50
1:B:59:GLN:HG2	1:B:277:VAL:O	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:99:ARG:CZ	1:A:103:GLY:HA2	2.42	0.50
1:B:254:HIS:C	1:B:256:THR:H	2.15	0.50
1:A:144:MET:O	1:A:228:VAL:HG12	2.12	0.50
1:C:80:ARG:HH11	1:C:80:ARG:HG2	1.76	0.50
1:A:85:THR:HB	1:A:99:ARG:HB3	1.94	0.49
1:C:43:VAL:O	1:C:47:VAL:HG23	2.12	0.49
1:C:95:LEU:HG	1:C:108:PHE:HE1	1.76	0.49
1:B:251:TRP:CG	1:B:270:LYS:HB3	2.47	0.49
1:C:254:HIS:N	1:C:255:PRO:HD3	2.26	0.49
1:B:202:VAL:O	1:B:206:CYS:HB2	2.12	0.49
1:B:24:THR:HG23	1:B:25:GLU:O	2.12	0.49
1:B:129:ARG:O	1:B:132:ASP:N	2.37	0.49
1:B:264:LYS:CG	1:C:254:HIS:HE1	2.24	0.49
1:C:61:VAL:HG11	1:C:249:PHE:HZ	1.77	0.49
1:B:280:VAL:HG12	1:B:281:LYS:N	2.28	0.49
1:B:259:ILE:O	1:B:261:ASP:N	2.46	0.49
1:A:251:TRP:CE3	1:A:270:LYS:HA	2.48	0.48
1:A:48:ILE:C	1:A:50:HIS:N	2.65	0.48
1:B:242:ASP:O	1:B:242:ASP:OD2	2.31	0.48
1:A:154:ASN:HD22	1:A:154:ASN:N	2.10	0.48
1:A:151:ARG:O	1:A:155:ASP:HB2	2.14	0.48
1:A:260:PRO:HA	1:A:263:GLN:HB2	1.95	0.48
1:C:194:THR:O	1:C:198:LEU:CG	2.58	0.48
1:A:30:GLU:OE2	1:A:80:ARG:HG2	2.13	0.48
1:A:83:SER:O	1:A:101:PRO:HG3	2.14	0.48
1:B:48:ILE:HA	1:B:51:CYS:SG	2.54	0.48
1:A:29:VAL:O	1:A:146:ARG:NH2	2.43	0.48
1:C:191:GLN:O	1:C:191:GLN:HG3	2.13	0.48
1:A:225:GLN:OE1	1:A:226:HIS:CE1	2.66	0.48
1:B:36:TYR:O	1:B:39:PHE:HB3	2.13	0.48
1:A:119:ASP:O	1:A:120:ASN:C	2.52	0.48
1:B:127:GLY:HA3	1:B:132:ASP:CG	2.34	0.48
1:C:36:TYR:HD2	1:C:219:ASP:HB2	1.79	0.48
1:C:233:THR:O	1:C:237:GLN:HG3	2.14	0.48
1:C:210:ARG:O	1:C:278:ALA:HB3	2.12	0.48
1:B:211:PHE:HB2	1:B:214:VAL:HG23	1.94	0.47
1:C:156:LEU:O	1:C:159:LYS:HB3	2.14	0.47
1:C:211:PHE:CZ	1:C:238:VAL:HA	2.49	0.47
1:C:245:SER:O	1:C:248:ALA:HB3	2.14	0.47
1:A:126:PHE:CD2	1:A:195:LYS:HE2	2.48	0.47
1:C:66:LYS:O	1:C:68:VAL:N	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:27:PHE:CD1	1:C:149:MET:HE2	2.49	0.47
1:C:85:THR:HB	1:C:99:ARG:HB3	1.96	0.47
1:B:99:ARG:HB2	1:B:105:TRP:CE2	2.50	0.47
1:C:260:PRO:HA	1:C:263:GLN:HG3	1.94	0.47
1:C:279:LEU:C	1:C:280:VAL:CG1	2.82	0.47
1:A:69:PRO:HG2	1:A:72:TRP:CZ2	2.49	0.47
1:A:250:GLU:CD	1:A:259:ILE:HD11	2.36	0.47
1:B:251:TRP:HB3	1:B:270:LYS:HB3	1.96	0.47
1:A:93:LEU:HD11	1:A:279:LEU:HD13	1.96	0.46
1:B:126:PHE:CG	1:B:195:LYS:HE2	2.50	0.46
1:C:237:GLN:HE22	1:C:276:ILE:CD1	2.28	0.46
1:A:122:ARG:O	1:A:123:TRP:C	2.53	0.46
1:B:203:VAL:HA	1:B:207:GLU:HG2	1.96	0.46
1:B:254:HIS:NE2	1:B:256:THR:OG1	2.43	0.46
1:C:237:GLN:NE2	1:C:276:ILE:HD13	2.29	0.46
1:B:210:ARG:O	1:B:278:ALA:N	2.45	0.46
1:B:93:LEU:HD11	1:B:279:LEU:HD12	1.97	0.46
1:B:196:SER:O	1:B:200:LYS:HG3	2.15	0.46
1:A:107:GLU:OE1	1:A:115:HIS:HD2	1.99	0.46
1:C:82:SER:OG	1:C:101:PRO:HG2	2.16	0.46
1:B:36:TYR:HD2	1:B:219:ASP:HB2	1.81	0.46
1:B:262:MET:HB3	1:B:267:ILE:HB	1.98	0.46
1:C:100:THR:OG1	1:C:104:VAL:HB	2.15	0.46
1:A:117:LEU:N	1:A:117:LEU:HD23	2.30	0.46
1:C:242:ASP:OD1	1:C:246:LYS:HE3	2.15	0.46
1:A:24:THR:HG22	1:A:25:GLU:O	2.16	0.46
1:A:30:GLU:HG2	1:A:150:THR:HG21	1.97	0.46
1:A:106:TRP:HE3	1:A:122:ARG:O	1.97	0.46
1:A:126:PHE:CG	1:A:195:LYS:HE2	2.51	0.46
1:A:43:VAL:HB	1:A:209:LEU:HD21	1.98	0.46
1:A:264:LYS:HA	1:A:264:LYS:HD3	1.77	0.45
1:A:95:LEU:HB3	2:A:501:ADE:H61	1.81	0.45
1:A:48:ILE:O	1:A:50:HIS:N	2.50	0.45
1:A:59:GLN:NE2	1:A:277:VAL:HG23	2.13	0.45
1:B:260:PRO:HA	1:B:263:GLN:HG3	1.97	0.45
1:C:24:THR:HG21	1:C:75:THR:CB	2.46	0.45
1:C:48:ILE:C	1:C:50:HIS:N	2.66	0.45
1:C:78:LYS:HB2	1:C:83:SER:HB3	1.97	0.45
1:A:272:GLU:O	1:A:275:ARG:HG2	2.16	0.45
1:A:48:ILE:C	1:A:50:HIS:H	2.19	0.45
1:B:130:TYR:HE2	2:B:501:ADE:N1	2.14	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:138:GLY:CA	1:A:140:GLU:OE1	2.65	0.45
1:B:59:GLN:N	1:B:59:GLN:OE1	2.47	0.45
1:A:24:THR:HG21	1:A:75:THR:HB	1.98	0.45
1:A:259:ILE:O	1:A:262:MET:N	2.49	0.45
1:B:90:MET:HB3	1:B:90:MET:HE2	1.76	0.45
1:A:101:PRO:HD3	1:A:157:ALA:HB1	1.99	0.45
1:A:87:ALA:HB2	1:A:105:TRP:CZ3	2.52	0.45
1:C:29:VAL:CG1	1:C:149:MET:HG2	2.38	0.45
1:C:77:LEU:HD22	1:C:149:MET:HE1	1.99	0.45
1:A:120:ASN:N	1:A:121:PRO:HD3	2.31	0.45
1:A:259:ILE:HA	1:A:260:PRO:HD2	1.71	0.44
1:B:50:HIS:CB	1:B:90:MET:CE	2.94	0.44
1:A:216:ARG:HA	1:A:216:ARG:HD3	1.65	0.44
1:C:119:ASP:O	1:C:120:ASN:HB2	2.15	0.44
1:C:251:TRP:CE3	1:C:270:LYS:HA	2.52	0.44
1:A:256:THR:O	1:A:257:ALA:C	2.56	0.44
1:A:250:GLU:OE1	1:A:259:ILE:HD13	2.17	0.44
1:C:279:LEU:C	1:C:280:VAL:HG12	2.38	0.44
1:A:138:GLY:HA3	1:A:140:GLU:OE1	2.16	0.44
1:A:268:LYS:HD2	1:A:272:GLU:OE1	2.18	0.44
1:A:109:GLY:HA3	1:A:123:TRP:CE3	2.53	0.44
1:A:233:THR:O	1:A:237:GLN:HG3	2.18	0.44
1:C:71:LEU:HB3	1:C:90:MET:HB2	2.00	0.44
1:A:106:TRP:CZ3	1:A:122:ARG:O	2.71	0.44
1:A:244:ILE:HA	1:A:262:MET:SD	2.58	0.44
1:A:108:PHE:O	1:A:123:TRP:HZ3	2.01	0.44
1:A:212:ASN:O	1:A:216:ARG:HG2	2.18	0.44
1:A:93:LEU:HD11	1:A:279:LEU:CD1	2.47	0.44
1:B:280:VAL:CG1	1:B:281:LYS:N	2.81	0.44
1:B:129:ARG:HA	2:B:501:ADE:H62	1.82	0.44
1:C:119:ASP:O	1:C:120:ASN:OD1	2.35	0.44
1:B:37:SER:CA	1:B:219:ASP:OD1	2.66	0.43
1:B:254:HIS:C	1:B:256:THR:N	2.69	0.43
1:B:39:PHE:O	1:B:43:VAL:HG23	2.18	0.43
1:C:254:HIS:C	1:C:256:THR:N	2.72	0.43
1:A:202:VAL:O	1:A:206:CYS:HB2	2.18	0.43
1:B:44:ARG:O	1:B:48:ILE:HG13	2.18	0.43
1:C:66:LYS:O	1:C:67:LYS:C	2.57	0.43
1:B:260:PRO:HA	1:B:263:GLN:CG	2.49	0.43
1:B:50:HIS:CB	1:B:90:MET:HE1	2.48	0.43
1:B:134:ILE:O	1:B:137:LYS:HB2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:244:ILE:O	1:C:247:ALA:HB3	2.19	0.43
1:A:230:LEU:HB3	1:A:234:GLN:HE21	1.82	0.43
1:B:85:THR:O	1:B:98:PHE:HA	2.18	0.43
1:C:29:VAL:HG12	1:C:150:THR:OG1	2.19	0.43
1:C:211:PHE:CB	1:C:214:VAL:HG23	2.42	0.43
1:C:106:TRP:HB3	1:C:124:LEU:HG	2.01	0.43
1:C:133:LEU:CD1	1:C:199:VAL:HG23	2.46	0.42
1:C:146:ARG:HB2	1:C:222:PHE:CD1	2.54	0.42
1:C:254:HIS:N	1:C:255:PRO:CD	2.82	0.42
1:B:130:TYR:OH	1:B:202:VAL:HG12	2.20	0.42
1:B:63:PRO:HG2	1:B:71:LEU:HD12	2.01	0.42
1:B:40:ILE:O	1:B:44:ARG:HG3	2.18	0.42
1:C:106:TRP:NE1	1:C:159:LYS:HG2	2.34	0.42
1:B:45:LYS:O	1:B:49:LYS:HG3	2.19	0.42
1:C:78:LYS:CB	1:C:83:SER:HB3	2.49	0.42
1:A:143:THR:HA	1:A:229:THR:HG22	2.01	0.42
1:A:206:CYS:O	1:A:210:ARG:HG3	2.20	0.42
1:C:146:ARG:HB2	1:C:222:PHE:CE1	2.54	0.42
1:C:237:GLN:HE22	1:C:276:ILE:HD13	1.84	0.42
1:C:77:LEU:HD22	1:C:149:MET:HE2	2.00	0.42
1:C:27:PHE:HE1	1:C:149:MET:SD	2.43	0.42
1:C:58:PHE:CD2	1:C:274:ALA:HB2	2.55	0.42
1:A:138:GLY:O	1:A:140:GLU:N	2.52	0.42
1:B:120:ASN:HD22	1:B:120:ASN:H	1.67	0.42
1:B:251:TRP:CB	1:B:270:LYS:HB3	2.50	0.42
1:C:120:ASN:HA	1:C:121:PRO:HD3	1.38	0.42
1:A:259:ILE:O	1:A:261:ASP:N	2.54	0.41
1:B:77:LEU:HD11	1:B:86:LEU:HD11	2.02	0.41
1:C:242:ASP:OD2	1:C:281:LYS:HD3	2.21	0.41
1:B:21:MET:N	1:B:72:TRP:H	2.18	0.41
1:B:259:ILE:CG2	1:B:262:MET:HG2	2.50	0.41
1:B:120:ASN:HD22	1:B:120:ASN:N	2.17	0.41
1:C:234:GLN:O	1:C:237:GLN:HB2	2.21	0.41
1:A:138:GLY:O	1:A:140:GLU:OE1	2.39	0.41
1:A:258:VAL:HG13	1:A:263:GLN:OE1	2.21	0.41
1:B:273:ALA:C	1:B:275:ARG:N	2.73	0.41
1:B:115:HIS:CE1	1:B:123:TRP:NE1	2.88	0.41
1:A:135:GLY:C	1:A:137:LYS:H	2.24	0.41
1:A:250:GLU:OE2	1:A:259:ILE:CD1	2.68	0.41
1:B:86:LEU:HB3	1:B:95:LEU:HD11	2.03	0.41
1:A:124:LEU:HD22	1:A:126:PHE:CZ	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:29:VAL:O	1:A:146:ARG:HG3	2.21	0.41
1:B:146:ARG:HB2	1:B:222:PHE:CD1	2.56	0.41
1:B:269:ASP:OD1	1:B:271:ASN:HB2	2.21	0.41
1:B:130:TYR:CE2	2:B:501:ADE:N1	2.89	0.41
1:A:107:GLU:HG2	1:A:123:TRP:HZ3	1.85	0.40
1:A:207:GLU:OE1	1:A:210:ARG:HD2	2.22	0.40
1:C:197:LYS:HA	1:C:197:LYS:HD3	1.78	0.40
1:C:254:HIS:C	1:C:256:THR:H	2.25	0.40
1:C:82:SER:O	1:C:83:SER:HB3	2.20	0.40
1:A:30:GLU:CD	1:A:80:ARG:HG2	2.41	0.40
1:B:211:PHE:CZ	1:B:241:TRP:HB2	2.56	0.40
1:A:31:ASP:OD2	1:A:34:TYR:CD2	2.75	0.40
1:A:30:GLU:CG	1:A:150:THR:HG21	2.52	0.40
1:C:192:ALA:HA	1:C:194:THR:HG23	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	229/243 (94%)	198 (86%)	25 (11%)	6 (3%)	6	21
1	B	228/243 (94%)	201 (88%)	24 (10%)	3 (1%)	14	41
1	C	226/243 (93%)	181 (80%)	40 (18%)	5 (2%)	8	26
All	All	683/729 (94%)	580 (85%)	89 (13%)	14 (2%)	9	28

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	119	ASP
1	C	120	ASN
1	C	257	ALA

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Mol	Chain	Res	Type
1	A	139	LEU
1	C	255	PRO
1	A	121	PRO
1	C	121	PRO
1	A	260	PRO
1	B	221	GLY
1	A	120	ASN
1	B	255	PRO
1	A	134	ILE
1	B	276	ILE
1	C	63	PRO

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	187/206 (91%)	168 (90%)	19 (10%)	8	25
1	B	186/206 (90%)	169 (91%)	17 (9%)	11	31
1	C	187/206 (91%)	168 (90%)	19 (10%)	8	25
All	All	560/618 (91%)	505 (90%)	55 (10%)	9	27

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	75	THR
1	A	81	THR
1	A	124	LEU
1	A	140	GLU
1	A	143	THR
1	A	155	ASP
1	A	193	ASP
1	A	194	THR
1	A	199	VAL
1	A	216	ARG
1	A	242	ASP

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Mol	Chain	Res	Type
1	A	254	HIS
1	A	256	THR
1	A	259	ILE
1	A	268	LYS
1	A	270	LYS
1	A	280	VAL
1	A	281	LYS
1	A	283	GLN
1	B	24	THR
1	B	25	GLU
1	B	26	ILE
1	B	34	TYR
1	B	57	ILE
1	B	78	LYS
1	B	82	SER
1	B	92	ASN
1	B	155	ASP
1	B	194	THR
1	B	258	VAL
1	B	268	LYS
1	B	269	ASP
1	B	270	LYS
1	B	275	ARG
1	B	276	ILE
1	B	281	LYS
1	C	23	PHE
1	C	24	THR
1	C	71	LEU
1	C	114	THR
1	C	148	GLU
1	C	149	MET
1	C	155	ASP
1	C	159	LYS
1	C	193	ASP
1	C	194	THR
1	C	196	SER
1	C	199	VAL
1	C	223	ASN
1	C	256	THR
1	C	258	VAL
1	C	268	LYS
1	C	270	LYS

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Mol	Chain	Res	Type
1	C	275	ARG
1	C	280	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	HIS
1	A	115	HIS
1	A	154	ASN
1	A	225	GLN
1	A	226	HIS
1	A	234	GLN
1	B	54	HIS
1	B	115	HIS
1	B	120	ASN
1	B	154	ASN
1	B	234	GLN
1	B	237	GLN
1	B	254	HIS
1	C	50	HIS
1	C	54	HIS
1	C	154	ASN
1	C	234	GLN
1	C	237	GLN

5.3.3 RNA ⓘ

There are no RNA molecules 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

3 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	ADE	A	501	-	9,11,11	0.95	0	7,15,15	3.47	3 (42%)
2	ADE	B	501	-	9,11,11	1.09	0	7,15,15	3.25	3 (42%)
2	ADE	C	501	-	9,11,11	0.98	0	7,15,15	3.29	3 (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
2	ADE	A	501	-	-	0/0/0/0	0/2/2/2
2	ADE	B	501	-	-	0/0/0/0	0/2/2/2
2	ADE	C	501	-	-	0/0/0/0	0/2/2/2

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	ADE	N3-C2-N1	-7.95	121.94	128.86
2	C	501	ADE	N3-C2-N1	-7.43	122.39	128.86
2	B	501	ADE	N3-C2-N1	-7.33	122.48	128.86
2	B	501	ADE	C4-C5-N7	-2.68	106.82	109.41
2	C	501	ADE	C4-C5-N7	-2.61	106.89	109.41
2	A	501	ADE	C4-C5-N7	-2.60	106.90	109.41
2	B	501	ADE	C2-N3-C4	3.27	121.00	113.33
2	C	501	ADE	C2-N3-C4	3.29	121.07	113.33
2	A	501	ADE	C2-N3-C4	3.33	121.16	113.33

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	ADE	2	0
2	B	501	ADE	5	0
2	C	501	ADE	2	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	233/243 (95%)	-0.08	2 (0%) 84 79	20, 55, 79, 101	0
1	B	232/243 (95%)	-0.12	1 (0%) 92 90	28, 55, 77, 90	0
1	C	232/243 (95%)	0.41	14 (6%) 23 14	37, 76, 100, 126	0
All	All	697/729 (95%)	0.07	17 (2%) 59 49	20, 60, 94, 126	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	267	ILE	3.5
1	C	33	ASN	3.1
1	C	220	ALA	3.0
1	C	253	ASP	3.0
1	C	249	PHE	3.0
1	C	145	GLY	2.9
1	C	102	GLY	2.9
1	A	118	GLY	2.8
1	C	250	GLU	2.5
1	C	248	ALA	2.3
1	C	149	MET	2.3
1	C	228	VAL	2.3
1	C	274	ALA	2.3
1	C	270	LYS	2.2
1	C	32	ALA	2.2
1	B	248	ALA	2.2
1	A	122	ARG	2.1

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 carbohydrates 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. 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, 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	LLDF	B-factors(\AA^2)	Q<0.9
2	ADE	B	501	10/10	0.42	0.49	15.74	90,101,103,104	0
2	ADE	A	501	10/10	0.85	0.64	13.92	116,117,117,118	0
2	ADE	C	501	10/10	0.86	0.53	9.14	116,117,118,118	0

6.5 Other polymers [i](#)

There are no such residues in this entry.