



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 07:31 PM UTC

PDB ID : 2Y1V / pdb_00002y1v
Title : Full length structure of RrgB Pilus protein from *Streptococcus pneumoniae*
Authors : El-Mortaji, L.; Contreras-Martel, C.; Manzano, C.; Vernet, T.; Dessen, A.; DiGuilmi, A.M.
Deposited on : 2010-12-10
Resolution : 2.39 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

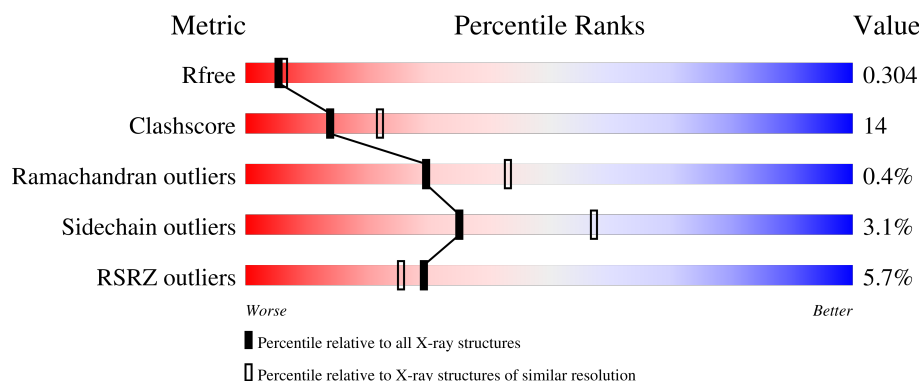
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.39 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	605	5% 76% 22% .
1	B	605	6% 76% 22% .
1	C	605	7% 66% 32% .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 14167 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 CELL WALL SURFACE ANCHOR FAMILY PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	604	Total	C	N	O	Se	0	0	0
			4576	2866	763	942	5			
1	B	604	Total	C	N	O	Se	0	0	0
			4576	2866	763	942	5			
1	C	604	Total	C	N	O	Se	0	0	0
			4576	2866	763	942	5			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	29	GLY	-	expression tag	UNP Q97SC2
B	29	GLY	-	expression tag	UNP Q97SC2
C	29	GLY	-	expression tag	UNP Q97SC2

- Molecule 2 is NICKEL (II) ION (CCD ID: NI) (formula: Ni).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	3	Total	Ni	0	0
			3	3		
2	B	3	Total	Ni	0	0
			3	3		
2	C	3	Total	Ni	0	0
			3	3		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	169	Total	O	0	0
			169	169		
3	B	141	Total	O	0	0
			141	141		

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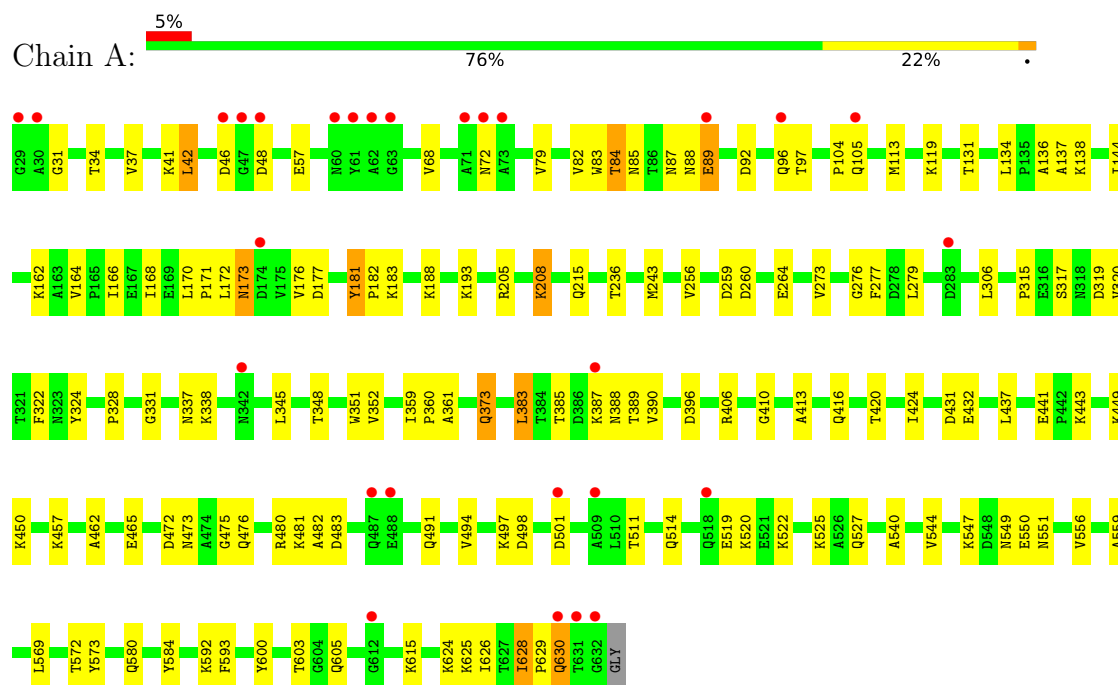
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	120	Total 120	O 120	0	0

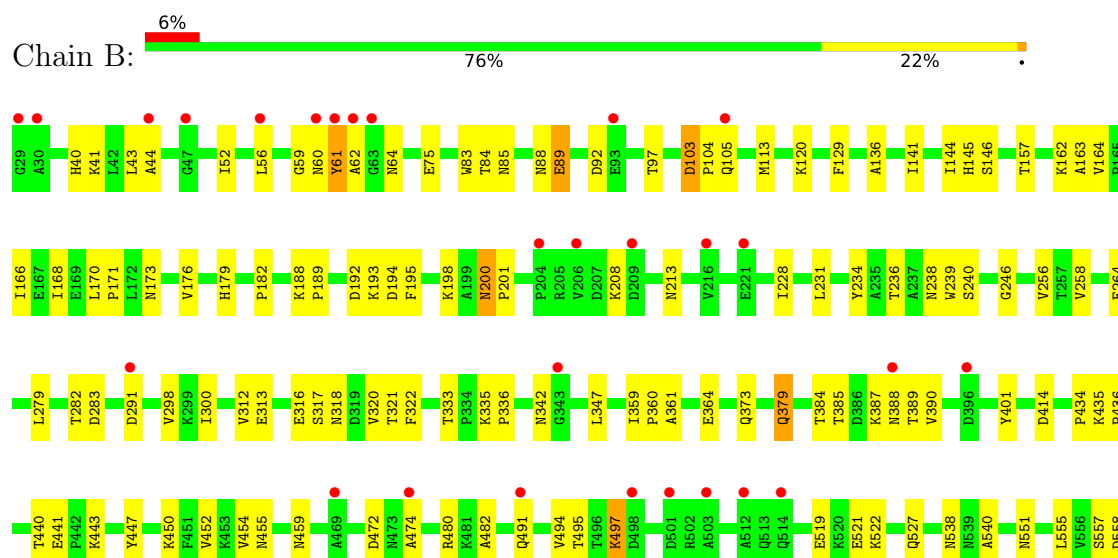
3 Residue-property plots

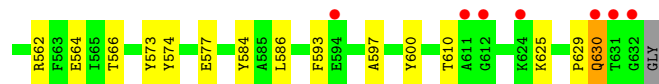
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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

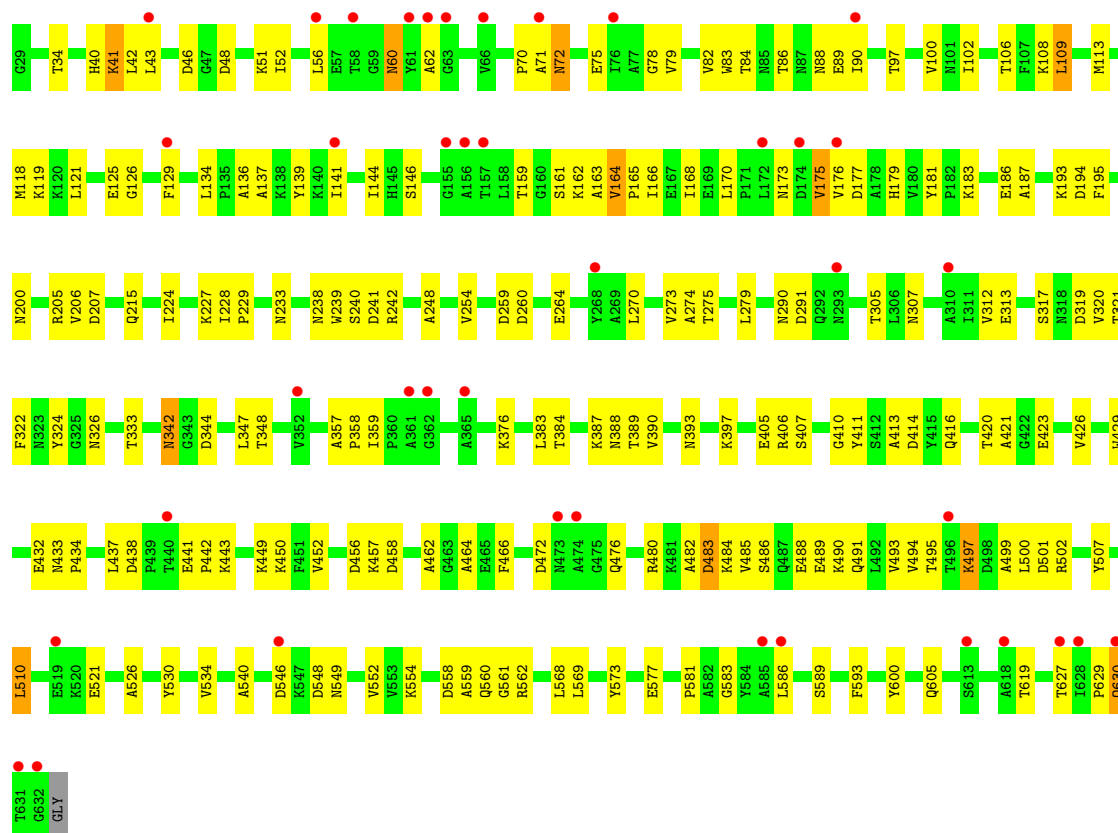


• Molecule 1: CELL WALL SURFACE ANCHOR FAMILY PROTEIN





● Molecule 1: CELL WALL SURFACE ANCHOR FAMILY PROTEIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	74.62Å 74.62Å 340.53Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	113.53 – 2.39 113.51 – 2.39	Depositor EDS
% Data completeness (in resolution range)	98.7 (113.53-2.39) 93.4 (113.51-2.39)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.25 (at 2.38Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.187 , 0.234 0.239 , 0.304	Depositor DCC
R_{free} test set	8066 reflections (10.26%)	wwPDB-VP
Wilson B-factor (Å ²)	41.6	Xtriage
Anisotropy	0.395	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 47.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.55$, $\langle L^2 \rangle = 0.39$	Xtriage
Estimated twinning fraction	0.000 for -h,-k,l 0.097 for h,-h-k,-l 0.002 for -k,-h,-l	Xtriage
Reported twinning fraction	0.592 for H, K, L 0.408 for K, H, -L	Depositor
Outliers	0 of 78179 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	14167	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.15% 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:
NI

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	0.59	0/4646	0.88	7/6315 (0.1%)
1	B	0.55	0/4646	0.84	4/6315 (0.1%)
1	C	0.52	0/4646	0.86	6/6315 (0.1%)
All	All	0.55	0/13938	0.86	17/18945 (0.1%)

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	175	VAL	N-CA-C	7.36	118.55	108.84
1	A	181	TYR	CA-C-N	6.25	126.22	119.78
1	A	181	TYR	C-N-CA	6.25	126.22	119.78
1	A	277	PHE	N-CA-C	6.05	118.16	109.14
1	C	200	ASN	CA-C-N	5.71	125.39	119.56
1	C	200	ASN	C-N-CA	5.71	125.39	119.56
1	A	188	LYS	CA-C-N	-5.62	114.79	120.85
1	A	188	LYS	C-N-CA	-5.62	114.79	120.85
1	A	628	ILE	CA-C-N	5.60	126.84	119.84
1	A	628	ILE	C-N-CA	5.60	126.84	119.84
1	B	200	ASN	CA-C-N	5.60	125.28	119.56
1	B	200	ASN	C-N-CA	5.60	125.28	119.56
1	C	173	ASN	CB-CA-C	-5.57	110.13	116.54
1	B	103	ASP	CA-C-N	5.56	125.23	119.56
1	B	103	ASP	C-N-CA	5.56	125.23	119.56
1	C	307	ASN	N-CA-C	5.32	117.61	109.95
1	C	568	LEU	N-CA-C	5.10	118.01	110.46

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	4576	0	4509	110	0
1	B	4576	0	4509	121	0
1	C	4576	0	4510	162	0
2	A	3	0	0	0	0
2	B	3	0	0	0	0
2	C	3	0	0	0	0
3	A	169	0	0	12	0
3	B	141	0	0	15	0
3	C	120	0	0	33	0
All	All	14167	0	13528	393	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 14.

All (393) 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)
3:A:2008:HOH:O	1:B:387:LYS:HB2	1.45	1.15
1:B:84:THR:HG21	1:B:88:ASN:HA	1.24	1.15
1:B:256:VAL:HG21	1:B:279:LEU:HD11	1.16	1.13
1:A:96:GLN:HB3	3:A:2012:HOH:O	1.48	1.12
1:C:168:ILE:HB	3:C:2026:HOH:O	1.47	1.11
1:C:248:ALA:HB3	3:C:2043:HOH:O	1.52	1.06
1:C:312:VAL:HG12	1:C:313:GLU:HG3	1.39	1.04
1:C:82:VAL:CG1	1:C:144:ILE:HD11	1.90	1.01
1:B:240:SER:HB2	3:B:2073:HOH:O	1.59	1.00
1:C:238:ASN:HB3	3:C:2039:HOH:O	1.63	0.98
1:B:84:THR:HG22	1:B:85:ASN:O	1.65	0.94
1:B:256:VAL:CG2	1:B:279:LEU:HD11	1.97	0.93
1:B:521:GLU:HG3	3:B:2117:HOH:O	1.68	0.92
1:A:113:MSE:HE1	1:A:144:ILE:CD1	1.99	0.92
1:B:84:THR:CG2	1:B:88:ASN:HA	1.98	0.92
1:B:256:VAL:HG21	1:B:279:LEU:CD1	2.00	0.92
1:C:491:GLN:O	1:C:494:VAL:HG12	1.70	0.90
1:C:348:THR:HG23	1:C:389:THR:HG22	1.53	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:82:VAL:HG11	1:C:144:ILE:HD11	1.53	0.89
1:C:70:PRO:HA	3:C:2002:HOH:O	1.73	0.88
1:C:464:ALA:HB2	1:C:581:PRO:HD3	1.53	0.87
1:B:629:PRO:C	1:B:630:GLN:HE21	1.83	0.85
1:C:416:GLN:HA	3:C:2083:HOH:O	1.76	0.85
1:A:84:THR:HG22	1:A:85:ASN:O	1.77	0.84
1:B:630:GLN:HE21	1:B:630:GLN:N	1.75	0.84
1:A:83:TRP:CH2	1:A:168:ILE:HD11	2.14	0.82
1:C:449:LYS:NZ	1:C:605:GLN:OE1	2.13	0.82
1:C:546:ASP:OD2	1:C:548:ASP:HB3	1.79	0.82
1:C:82:VAL:HG12	1:C:144:ILE:HD11	1.61	0.80
1:A:511:THR:HG22	1:A:514:GLN:HE21	1.47	0.80
1:C:102:ILE:HG12	1:C:109:LEU:HD12	1.62	0.80
1:C:462:ALA:HB2	1:C:559:ALA:HA	1.63	0.80
1:C:549:ASN:HB3	1:C:552:VAL:HG23	1.62	0.79
1:B:88:ASN:ND2	1:B:104:PRO:HG3	1.98	0.79
1:C:629:PRO:C	1:C:630:GLN:HE21	1.91	0.79
1:B:491:GLN:O	1:B:494:VAL:HG12	1.83	0.78
1:C:452:VAL:HA	1:C:561:GLY:O	1.82	0.78
1:A:113:MSE:HE1	1:A:144:ILE:HD13	1.68	0.75
1:B:387:LYS:O	1:B:387:LYS:HG3	1.88	0.74
1:B:40:HIS:NE2	1:B:179:HIS:HD2	1.85	0.74
1:C:583:GLY:HA3	3:C:2114:HOH:O	1.87	0.73
1:C:60:ASN:OD1	1:C:62:ALA:O	2.07	0.72
1:C:224:ILE:HD11	1:C:320:VAL:HG21	1.71	0.72
1:A:549:ASN:OD1	1:A:550:GLU:N	2.22	0.72
1:A:416:GLN:HB3	3:A:2102:HOH:O	1.90	0.71
1:B:246:GLY:HA3	1:B:316:GLU:CD	2.15	0.71
1:A:624:LYS:HG2	3:A:2168:HOH:O	1.92	0.70
1:C:376:LYS:HD2	3:C:2071:HOH:O	1.91	0.70
1:C:423:GLU:O	3:C:2084:HOH:O	2.08	0.70
1:A:82:VAL:HG11	1:A:144:ILE:HD11	1.72	0.70
1:A:84:THR:CG2	1:A:85:ASN:O	2.40	0.69
1:B:92:ASP:OD2	3:B:2013:HOH:O	2.11	0.69
1:A:84:THR:HG21	1:A:88:ASN:HA	1.73	0.69
1:C:489:GLU:OE1	3:C:2091:HOH:O	2.10	0.69
1:C:410:GLY:HA2	1:C:432:GLU:HA	1.76	0.68
1:B:192:ASP:OD1	1:B:208:LYS:NZ	2.24	0.68
1:B:61:TYR:OH	1:B:64:ASN:ND2	2.27	0.67
1:C:319:ASP:HA	1:C:441:GLU:CD	2.20	0.67
1:C:348:THR:HG23	1:C:389:THR:CG2	2.24	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:84:THR:CG2	1:B:88:ASN:CA	2.73	0.66
1:C:305:THR:HB	3:C:2043:HOH:O	1.94	0.66
1:B:573:TYR:HB2	1:B:593:PHE:CZ	2.31	0.66
1:C:554:LYS:NZ	3:C:2107:HOH:O	2.29	0.66
1:C:82:VAL:CG1	1:C:144:ILE:CD1	2.73	0.65
1:B:387:LYS:C	3:B:2084:HOH:O	2.39	0.65
1:A:603:THR:CG2	1:A:615:LYS:HG2	2.27	0.65
1:A:166:ILE:HG22	1:A:168:ILE:HG23	1.77	0.64
1:B:494:VAL:HG23	3:B:2114:HOH:O	1.96	0.64
1:C:72:ASN:HB3	3:C:2008:HOH:O	1.97	0.64
1:C:83:TRP:CH2	1:C:168:ILE:HD11	2.32	0.64
1:B:256:VAL:CG2	1:B:279:LEU:CD1	2.68	0.64
1:A:387:LYS:CG	1:A:387:LYS:O	2.47	0.63
1:C:238:ASN:CB	3:C:2039:HOH:O	2.30	0.63
1:A:113:MSE:CE	1:A:144:ILE:HD13	2.29	0.62
1:B:146:SER:HB2	3:B:2022:HOH:O	1.99	0.62
1:C:383:LEU:CD1	1:C:390:VAL:HB	2.30	0.62
1:B:480:ARG:HD3	1:B:538:ASN:OD1	2.00	0.62
1:A:361:ALA:HB3	1:A:385:THR:HG23	1.80	0.62
1:A:476:GLN:HB3	1:A:544:VAL:O	2.00	0.61
1:C:270:LEU:HD13	1:C:279:LEU:HD12	1.83	0.61
1:C:321:THR:HG21	3:C:2088:HOH:O	1.98	0.61
1:C:90:ILE:HD11	1:C:109:LEU:HD11	1.82	0.61
1:C:383:LEU:HD11	1:C:390:VAL:HB	1.82	0.61
1:C:433:ASN:HB3	1:C:434:PRO:CD	2.31	0.61
1:C:82:VAL:HG12	1:C:144:ILE:CD1	2.30	0.61
1:C:558:ASP:OD1	1:C:560:GLN:N	2.30	0.61
1:A:113:MSE:CE	1:A:144:ILE:CD1	2.76	0.61
1:C:41:LYS:HG3	1:C:79:VAL:HG21	1.82	0.61
1:B:519:GLU:HA	1:B:519:GLU:OE1	2.01	0.60
1:B:388:ASN:HA	3:B:2084:HOH:O	2.01	0.60
1:A:42:LEU:HD22	1:A:181:TYR:HB3	1.83	0.60
1:C:242:ARG:NH1	3:C:2041:HOH:O	2.32	0.60
1:A:215:GLN:HB3	1:A:569:LEU:HG	1.83	0.60
1:B:83:TRP:CH2	1:B:168:ILE:HD11	2.37	0.60
1:A:168:ILE:O	1:A:168:ILE:HG13	2.00	0.60
1:C:627:THR:O	1:C:629:PRO:HD3	2.02	0.60
1:B:258:VAL:HG22	1:B:298:VAL:HG22	1.84	0.60
1:C:240:SER:HB2	3:C:2060:HOH:O	2.02	0.60
1:C:241:ASP:HB2	1:C:320:VAL:HG22	1.83	0.59
1:A:629:PRO:C	1:A:630:GLN:HE21	2.09	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:603:THR:HG21	1:A:615:LYS:HG2	1.85	0.59
1:B:359:ILE:HB	1:B:360:PRO:HD2	1.84	0.59
1:B:452:VAL:HG12	1:B:454:VAL:HG13	1.84	0.59
1:B:40:HIS:NE2	1:B:179:HIS:CD2	2.69	0.59
1:A:37:VAL:HG23	1:A:131:THR:HG21	1.85	0.59
1:B:59:GLY:CA	3:B:2006:HOH:O	2.51	0.59
1:B:231:LEU:HA	1:B:291:ASP:HA	1.84	0.58
1:C:242:ARG:HD3	3:C:2041:HOH:O	2.02	0.58
1:C:450:LYS:HE3	1:C:600:TYR:CZ	2.38	0.58
1:C:546:ASP:OD2	1:C:548:ASP:CB	2.49	0.58
1:A:491:GLN:HA	1:A:494:VAL:HG12	1.83	0.58
1:C:166:ILE:HG12	3:C:2025:HOH:O	2.03	0.58
1:B:388:ASN:CA	3:B:2084:HOH:O	2.50	0.58
1:C:433:ASN:HB3	1:C:434:PRO:HD2	1.85	0.58
1:C:322:PHE:CE2	1:C:324:TYR:HB3	2.39	0.58
1:C:194:ASP:OD1	1:C:195:PHE:N	2.35	0.58
1:C:630:GLN:HE21	1:C:630:GLN:N	2.00	0.58
1:A:82:VAL:CG1	1:A:144:ILE:HD11	2.33	0.57
1:C:113:MSE:HG2	1:C:119:LYS:HE2	1.86	0.57
1:B:385:THR:HG22	1:B:385:THR:O	2.04	0.57
1:C:118:MSE:HE2	1:C:129:PHE:HB3	1.86	0.57
1:C:275:THR:HG23	3:C:2048:HOH:O	2.03	0.57
1:B:194:ASP:OD1	1:B:195:PHE:N	2.32	0.57
1:C:40:HIS:NE2	1:C:179:HIS:HD2	2.03	0.56
1:C:259:ASP:O	1:C:260:ASP:HB2	2.04	0.56
1:C:507:TYR:HA	1:C:510:LEU:HD12	1.87	0.56
1:B:497:LYS:HG2	3:B:2119:HOH:O	2.05	0.56
1:C:411:TYR:HA	1:C:429:TRP:O	2.04	0.56
1:A:256:VAL:HG21	1:A:279:LEU:HD11	1.86	0.56
1:C:141:ILE:HD11	1:C:168:ILE:HD11	1.87	0.56
1:C:507:TYR:HA	1:C:510:LEU:CD1	2.36	0.56
1:A:259:ASP:O	1:A:260:ASP:HB2	2.04	0.56
1:C:240:SER:O	1:C:320:VAL:HA	2.06	0.56
1:A:373:GLN:O	1:A:443:LYS:HD2	2.06	0.56
1:A:572:THR:HG21	1:A:592:LYS:HE3	1.87	0.55
1:C:90:ILE:HD12	1:C:100:VAL:HG12	1.87	0.55
1:C:347:LEU:HD11	1:C:426:VAL:HG23	1.88	0.55
1:B:317:SER:HB2	1:B:443:LYS:HD3	1.87	0.55
1:B:162:LYS:O	1:B:163:ALA:HB3	2.05	0.55
1:A:84:THR:CG2	1:A:88:ASN:HA	2.36	0.55
1:A:345:LEU:HD21	1:A:424:ILE:HD12	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:465:GLU:HG2	1:A:556:VAL:HG22	1.87	0.55
1:B:208:LYS:HD2	1:B:440:THR:OG1	2.07	0.55
1:C:146:SER:HA	1:C:326:ASN:O	2.06	0.55
1:C:573:TYR:HB2	1:C:593:PHE:CZ	2.42	0.55
1:A:162:LYS:NZ	3:A:2027:HOH:O	2.40	0.55
1:A:317:SER:HB2	1:A:443:LYS:HD3	1.88	0.55
1:C:89:GLU:HB2	3:C:2011:HOH:O	2.07	0.55
1:C:238:ASN:CA	3:C:2039:HOH:O	2.55	0.54
1:B:84:THR:CG2	1:B:85:ASN:O	2.49	0.54
1:B:472:ASP:HB3	1:B:574:TYR:CE2	2.42	0.54
1:A:164:VAL:HG12	3:A:2023:HOH:O	2.07	0.54
1:B:59:GLY:HA2	3:B:2006:HOH:O	2.07	0.54
1:B:312:VAL:HG12	1:B:313:GLU:HG3	1.88	0.54
1:C:548:ASP:HA	3:C:2104:HOH:O	2.07	0.54
1:A:193:LYS:HD3	1:A:441:GLU:HG3	1.90	0.53
1:A:193:LYS:HD2	1:A:320:VAL:CG2	2.38	0.53
1:C:141:ILE:HB	1:C:166:ILE:HD12	1.90	0.53
1:A:113:MSE:HE1	1:A:144:ILE:HD12	1.90	0.53
1:B:629:PRO:C	1:B:630:GLN:NE2	2.60	0.53
1:C:224:ILE:CD1	1:C:320:VAL:HG21	2.38	0.53
1:A:319:ASP:HA	1:A:441:GLU:CD	2.33	0.53
1:B:279:LEU:C	1:B:279:LEU:HD23	2.34	0.53
1:C:558:ASP:OD1	1:C:558:ASP:C	2.52	0.53
1:B:347:LEU:O	1:B:389:THR:HA	2.08	0.53
1:B:388:ASN:N	3:B:2084:HOH:O	2.42	0.53
1:C:484:LYS:O	1:C:484:LYS:HG3	2.08	0.53
1:A:164:VAL:CG1	3:A:2023:HOH:O	2.56	0.52
1:C:224:ILE:HD13	1:C:320:VAL:HG11	1.91	0.52
1:C:558:ASP:OD2	1:C:562:ARG:NH2	2.37	0.52
1:A:172:LEU:O	1:A:173:ASN:C	2.52	0.52
1:C:206:VAL:HG22	1:C:206:VAL:O	2.08	0.52
1:A:387:LYS:O	1:A:387:LYS:HG3	2.10	0.52
1:B:494:VAL:HG13	1:B:495:THR:N	2.24	0.52
1:A:462:ALA:HB2	1:A:559:ALA:HA	1.92	0.52
1:C:242:ARG:CZ	3:C:2041:HOH:O	2.58	0.52
1:A:481:LYS:HE3	3:A:2137:HOH:O	2.09	0.52
1:A:88:ASN:ND2	1:A:104:PRO:HG3	2.23	0.52
1:B:385:THR:O	1:B:385:THR:CG2	2.58	0.51
1:C:72:ASN:CG	1:C:72:ASN:O	2.53	0.51
1:A:83:TRP:CZ3	1:A:168:ILE:HD11	2.44	0.51
1:C:78:GLY:HA2	1:C:121:LEU:HD22	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:126:GLY:HA2	3:C:2017:HOH:O	2.10	0.51
1:B:84:THR:HG23	1:B:88:ASN:C	2.36	0.51
1:B:387:LYS:O	1:B:387:LYS:CG	2.57	0.51
1:A:472:ASP:OD1	1:A:472:ASP:C	2.54	0.51
1:C:546:ASP:OD1	1:C:548:ASP:HB2	2.11	0.51
1:B:113:MSE:HE1	1:B:144:ILE:HD12	1.92	0.51
1:B:282:THR:O	1:B:283:ASP:C	2.54	0.51
1:B:584:TYR:CD2	1:B:625:LYS:HB2	2.46	0.51
1:C:629:PRO:C	1:C:630:GLN:NE2	2.66	0.51
1:C:493:VAL:O	1:C:497:LYS:HB2	2.10	0.51
1:C:228:ILE:HG23	1:C:229:PRO:HD2	1.92	0.50
1:C:482:ALA:O	1:C:485:VAL:HG12	2.11	0.50
1:C:162:LYS:O	1:C:163:ALA:HB3	2.10	0.50
1:A:85:ASN:OD1	1:A:85:ASN:C	2.55	0.50
1:C:41:LYS:O	1:C:75:GLU:HA	2.12	0.50
1:C:136:ALA:O	1:C:137:ALA:HB2	2.11	0.50
1:C:164:VAL:HA	1:C:165:PRO:C	2.35	0.50
1:C:449:LYS:HZ1	1:C:619:THR:HG1	1.55	0.50
1:B:193:LYS:HD3	1:B:441:GLU:HG3	1.93	0.50
1:A:92:ASP:C	1:A:92:ASP:OD1	2.55	0.49
1:C:384:THR:H	1:C:387:LYS:HE3	1.77	0.49
1:B:494:VAL:CG1	1:B:495:THR:N	2.75	0.49
1:C:480:ARG:HA	1:C:540:ALA:O	2.11	0.49
1:C:420:THR:HG22	1:C:421:ALA:N	2.27	0.49
1:B:318:ASN:O	1:B:441:GLU:HG2	2.13	0.49
1:A:72:ASN:OD1	1:A:72:ASN:O	2.31	0.49
1:B:120:LYS:HB2	1:B:129:PHE:CZ	2.47	0.49
1:B:43:LEU:HD23	1:B:44:ALA:C	2.38	0.49
1:A:166:ILE:HD11	1:A:182:PRO:HD3	1.95	0.48
1:A:580:GLN:HG3	1:A:584:TYR:O	2.13	0.48
1:A:603:THR:HG22	1:A:615:LYS:HG2	1.95	0.48
1:C:42:LEU:HD22	1:C:181:TYR:HB3	1.95	0.48
1:C:500:LEU:HA	1:C:526:ALA:HB1	1.94	0.48
1:C:134:LEU:HB3	1:C:139:TYR:OH	2.14	0.48
1:A:348:THR:HG23	1:A:389:THR:HG22	1.94	0.48
1:B:188:LYS:HG2	1:B:189:PRO:O	2.14	0.48
1:B:335:LYS:HD2	1:B:335:LYS:N	2.28	0.48
1:A:482:ALA:HB2	1:A:551:ASN:HD21	1.79	0.48
1:A:628:ILE:HA	1:A:629:PRO:HD3	1.66	0.48
1:B:166:ILE:HD11	1:B:182:PRO:HG3	1.95	0.48
1:C:273:VAL:HG12	1:C:274:ALA:N	2.29	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:342:ASN:N	1:C:342:ASN:ND2	2.60	0.48
1:A:450:LYS:HE2	1:A:600:TYR:O	2.14	0.47
1:B:84:THR:HG23	1:B:89:GLU:N	2.28	0.47
1:B:312:VAL:CG1	1:B:313:GLU:HG3	2.43	0.47
1:C:499:ALA:HA	1:C:502:ARG:NH2	2.29	0.47
1:A:89:GLU:HG2	1:A:97:THR:CG2	2.44	0.47
1:C:46:ASP:OD1	1:C:72:ASN:OD1	2.32	0.47
1:C:414:ASP:OD1	1:C:414:ASP:C	2.55	0.47
1:A:359:ILE:HB	1:A:360:PRO:HD2	1.96	0.47
1:B:454:VAL:HA	1:B:459:ASN:O	2.14	0.47
1:C:450:LYS:HE3	1:C:600:TYR:CE1	2.49	0.47
1:B:141:ILE:HD11	1:B:168:ILE:HD11	1.97	0.47
1:C:161:SER:HB3	3:C:2024:HOH:O	2.14	0.47
1:A:41:LYS:HG3	1:A:79:VAL:HG11	1.96	0.47
1:A:136:ALA:O	1:A:137:ALA:HB2	2.14	0.47
1:B:44:ALA:HB2	1:B:52:ILE:CD1	2.45	0.47
1:B:145:HIS:HA	3:B:2024:HOH:O	2.14	0.47
1:C:240:SER:HB3	1:C:333:THR:HB	1.97	0.47
1:A:626:ILE:HA	3:A:2169:HOH:O	2.14	0.47
1:B:136:ALA:HA	1:B:170:LEU:O	2.14	0.47
1:B:494:VAL:CG2	3:B:2114:HOH:O	2.61	0.47
1:B:482:ALA:HB2	1:B:551:ASN:ND2	2.30	0.47
1:C:546:ASP:CG	1:C:548:ASP:CB	2.88	0.47
1:A:472:ASP:OD1	1:A:473:ASN:N	2.47	0.47
1:A:549:ASN:OD1	1:A:549:ASN:C	2.58	0.47
1:B:450:LYS:HE3	1:B:600:TYR:CE2	2.50	0.47
1:B:450:LYS:HE3	1:B:600:TYR:CZ	2.50	0.47
1:C:239:TRP:CZ2	1:C:322:PHE:CE1	3.03	0.47
1:C:472:ASP:OD1	1:C:472:ASP:C	2.58	0.47
1:A:162:LYS:HB3	1:A:183:LYS:HB2	1.96	0.46
1:A:322:PHE:HB3	1:A:437:LEU:HB2	1.97	0.46
1:A:472:ASP:OD1	1:A:473:ASN:HB2	2.15	0.46
1:B:347:LEU:HB3	1:B:390:VAL:HG12	1.97	0.46
1:C:52:ILE:HB	1:C:183:LYS:HD3	1.97	0.46
1:A:34:THR:HB	1:A:170:LEU:HD22	1.97	0.46
1:B:447:TYR:CE1	1:B:597:ALA:HA	2.50	0.46
1:C:420:THR:HG22	1:C:421:ALA:O	2.14	0.46
1:A:331:GLY:HA2	1:A:413:ALA:O	2.15	0.46
1:B:455:ASN:ND2	1:B:459:ASN:HB2	2.30	0.46
1:C:84:THR:CG2	1:C:102:ILE:HD12	2.46	0.46
1:C:205:ARG:HG2	1:C:207:ASP:OD1	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:527:GLN:OE1	1:A:527:GLN:C	2.59	0.46
1:C:321:THR:HG22	1:C:438:ASP:OD1	2.16	0.46
1:A:170:LEU:HA	1:A:171:PRO:C	2.40	0.46
1:A:396:ASP:OD1	1:A:396:ASP:C	2.59	0.46
1:B:61:TYR:CG	1:B:62:ALA:N	2.84	0.46
1:B:256:VAL:HG22	1:B:300:ILE:HG12	1.96	0.46
1:C:75:GLU:HG2	3:C:2017:HOH:O	2.14	0.46
1:A:46:ASP:CG	1:A:48:ASP:H	2.23	0.46
1:A:406:ARG:NH1	3:A:2099:HOH:O	2.38	0.46
1:B:482:ALA:HB2	1:B:551:ASN:HD21	1.80	0.46
1:C:464:ALA:CB	1:C:581:PRO:HD3	2.37	0.46
1:A:337:ASN:OD1	1:A:338:LYS:N	2.46	0.45
1:A:547:LYS:HG2	3:A:2139:HOH:O	2.15	0.45
1:B:44:ALA:HB2	1:B:52:ILE:HD11	1.97	0.45
1:C:193:LYS:HD3	1:C:441:GLU:HG2	1.98	0.45
1:B:562:ARG:NH1	1:B:610:THR:OG1	2.44	0.45
1:C:41:LYS:HD3	1:C:42:LEU:N	2.32	0.45
1:C:321:THR:CG2	3:C:2088:HOH:O	2.61	0.45
1:C:233:ASN:O	3:C:2037:HOH:O	2.21	0.45
1:A:383:LEU:HA	1:A:387:LYS:HE3	1.99	0.45
1:A:193:LYS:HD2	1:A:320:VAL:HG21	1.99	0.45
1:C:88:ASN:HB3	1:C:102:ILE:O	2.17	0.45
1:A:525:LYS:HB3	1:A:525:LYS:HE2	1.73	0.45
1:C:176:VAL:HG23	1:C:177:ASP:N	2.32	0.45
1:C:405:GLU:OE2	1:C:413:ALA:HB2	2.17	0.45
1:C:494:VAL:CG1	1:C:495:THR:N	2.80	0.45
1:A:256:VAL:CG2	1:A:279:LEU:HD11	2.47	0.44
1:A:236:THR:OG1	1:A:328:PRO:HA	2.17	0.44
1:A:388:ASN:OD1	1:A:389:THR:HG23	2.17	0.44
1:B:89:GLU:CG	1:B:97:THR:HB	2.46	0.44
1:B:103:ASP:HA	1:B:104:PRO:HD3	1.84	0.44
1:C:290:ASN:O	1:C:291:ASP:HB2	2.16	0.44
1:B:195:PHE:HB2	1:B:198:LYS:HD3	1.99	0.44
1:B:577:GLU:O	1:B:586:LEU:HD11	2.17	0.44
1:A:630:GLN:HE21	1:A:630:GLN:N	2.15	0.44
1:B:170:LEU:HA	1:B:171:PRO:C	2.43	0.44
1:B:361:ALA:HB3	1:B:385:THR:HG23	2.00	0.44
1:C:358:PRO:O	1:C:359:ILE:HG23	2.17	0.44
1:A:584:TYR:CD2	1:A:625:LYS:HB2	2.52	0.44
1:C:194:ASP:HA	1:C:442:PRO:HG3	1.99	0.44
1:B:239:TRP:CE2	1:B:322:PHE:CE1	3.06	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:97:THR:HG21	3:C:2013:HOH:O	2.18	0.44
1:A:457:LYS:HE3	1:A:457:LYS:HB2	1.68	0.44
1:B:435:LYS:HB2	1:B:436:PRO:HD2	2.00	0.44
1:C:577:GLU:HG2	1:C:586:LEU:HD12	2.00	0.44
1:C:344:ASP:OD1	1:C:393:ASN:OD1	2.36	0.43
1:C:494:VAL:HG13	1:C:495:THR:N	2.33	0.43
1:A:410:GLY:HA2	1:A:432:GLU:HA	2.01	0.43
1:B:447:TYR:HB3	3:B:2134:HOH:O	2.18	0.43
1:C:476:GLN:HE21	1:C:476:GLN:HB2	1.63	0.43
1:A:87:ASN:O	1:A:88:ASN:HB2	2.18	0.43
1:A:92:ASP:OD2	1:A:96:GLN:HB2	2.17	0.43
1:B:213:ASN:HB2	1:B:447:TYR:CZ	2.53	0.43
1:B:336:PRO:HA	1:B:401:TYR:O	2.18	0.43
1:B:480:ARG:HA	1:B:540:ALA:O	2.19	0.43
1:C:159:THR:HG21	1:C:187:ALA:HB3	2.00	0.43
1:B:157:THR:HG22	1:B:434:PRO:HG2	1.99	0.43
1:A:520:LYS:HG2	1:A:559:ALA:HB1	2.01	0.43
1:C:215:GLN:HB3	1:C:569:LEU:HG	2.01	0.43
1:C:317:SER:HB2	1:C:443:LYS:HD3	2.00	0.43
1:C:406:ARG:O	1:C:407:SER:C	2.62	0.43
1:A:82:VAL:HG23	1:A:119:LYS:HB3	2.00	0.43
1:A:138:LYS:HD2	3:A:2007:HOH:O	2.18	0.43
1:B:200:ASN:HA	1:B:201:PRO:HD3	1.89	0.43
1:B:364:GLU:HG2	1:B:384:THR:HA	2.01	0.43
1:C:48:ASP:CG	1:C:51:LYS:HB2	2.44	0.43
1:A:205:ARG:HB3	1:A:208:LYS:HG3	2.01	0.42
1:A:498:ASP:O	1:A:501:ASP:HB2	2.19	0.42
1:C:43:LEU:HD21	1:C:186:GLU:HG2	2.01	0.42
1:C:530:TYR:O	1:C:534:VAL:HG23	2.19	0.42
1:A:315:PRO:O	1:A:373:GLN:HB3	2.19	0.42
1:B:312:VAL:O	1:B:313:GLU:HB2	2.19	0.42
1:C:34:THR:HB	1:C:170:LEU:HD22	2.01	0.42
1:C:456:ASP:OD1	1:C:457:LYS:HG3	2.19	0.42
1:B:85:ASN:C	1:B:85:ASN:OD1	2.61	0.42
1:B:236:THR:CG2	1:B:238:ASN:ND2	2.82	0.42
1:C:40:HIS:NE2	1:C:179:HIS:CD2	2.84	0.42
1:A:215:GLN:O	1:A:306:LEU:HD23	2.19	0.42
1:B:519:GLU:OE1	1:B:522:LYS:HD3	2.20	0.42
1:C:357:ALA:HA	1:C:358:PRO:HD3	1.94	0.42
1:A:193:LYS:HD2	1:A:320:VAL:HG23	2.01	0.42
1:B:44:ALA:CB	1:B:52:ILE:HD11	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:320:VAL:HG12	1:B:321:THR:N	2.34	0.42
1:C:483:ASP:OD1	1:C:483:ASP:C	2.61	0.42
1:B:459:ASN:N	1:B:459:ASN:HD22	2.18	0.42
1:B:527:GLN:CG	1:B:558:ASP:HB3	2.50	0.42
1:A:573:TYR:HB2	1:A:593:PHE:CZ	2.54	0.42
1:B:557:SER:HA	1:B:562:ARG:O	2.19	0.42
1:C:500:LEU:O	1:C:501:ASP:C	2.62	0.42
1:C:227:LYS:HD2	3:C:2053:HOH:O	2.20	0.42
1:C:273:VAL:HG12	3:C:2048:HOH:O	2.19	0.41
1:B:89:GLU:HG2	1:B:97:THR:HB	2.03	0.41
1:C:388:ASN:O	1:C:389:THR:HG22	2.19	0.41
1:A:351:TRP:C	1:A:352:VAL:HG13	2.44	0.41
1:B:519:GLU:OE1	1:B:522:LYS:CD	2.68	0.41
1:B:555:LEU:HD22	1:B:564:GLU:O	2.21	0.41
1:B:239:TRP:CZ2	1:B:322:PHE:CE1	3.09	0.41
1:C:106:THR:C	1:C:108:LYS:H	2.29	0.41
1:A:383:LEU:HD11	1:A:390:VAL:HG12	2.02	0.41
1:A:449:LYS:NZ	1:A:605:GLN:OE1	2.38	0.41
1:C:70:PRO:C	1:C:72:ASN:H	2.29	0.41
1:C:206:VAL:O	1:C:206:VAL:CG2	2.68	0.41
1:C:486:SER:C	1:C:488:GLU:H	2.29	0.41
1:A:273:VAL:HG22	1:A:276:GLY:O	2.19	0.41
1:A:431:ASP:O	1:A:432:GLU:HB2	2.20	0.41
1:C:83:TRP:CZ2	1:C:141:ILE:HD11	2.56	0.41
1:A:629:PRO:C	1:A:630:GLN:NE2	2.77	0.41
1:B:312:VAL:HG23	1:B:566:THR:HB	2.02	0.41
1:A:31:GLY:HA2	1:A:134:LEU:O	2.20	0.41
1:A:480:ARG:HA	1:A:540:ALA:O	2.20	0.41
1:C:90:ILE:HD12	1:C:100:VAL:CG1	2.51	0.41
1:C:388:ASN:C	1:C:389:THR:CG2	2.93	0.41
1:C:466:PHE:CE2	1:C:577:GLU:HB2	2.56	0.41
1:A:322:PHE:CE2	1:A:324:TYR:HB3	2.56	0.41
1:B:41:LYS:O	1:B:75:GLU:HA	2.21	0.41
1:B:89:GLU:HG2	1:B:97:THR:CG2	2.51	0.41
1:B:320:VAL:CG1	1:B:321:THR:N	2.84	0.41
1:B:414:ASP:OD1	1:B:414:ASP:C	2.64	0.41
1:C:254:VAL:HB	3:C:2045:HOH:O	2.21	0.41
1:B:83:TRP:CH2	1:B:168:ILE:CD1	3.03	0.40
1:C:383:LEU:HD12	1:C:390:VAL:HB	2.04	0.40
1:B:92:ASP:OD1	1:B:92:ASP:C	2.64	0.40
1:B:228:ILE:HD12	1:B:234:TYR:HE2	1.86	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:240:SER:HB3	1:B:333:THR:HB	2.03	0.40
1:C:175:VAL:HA	3:C:2027:HOH:O	2.21	0.40
1:A:176:VAL:HG23	1:A:177:ASP:OD1	2.22	0.40
1:A:519:GLU:O	1:A:522:LYS:HB2	2.22	0.40
1:C:273:VAL:O	1:C:274:ALA:C	2.63	0.40
1:C:497:LYS:HG2	1:C:530:TYR:CE1	2.56	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	602/605 (100%)	566 (94%)	34 (6%)	2 (0%)	36	50
1	B	602/605 (100%)	573 (95%)	25 (4%)	4 (1%)	18	28
1	C	602/605 (100%)	569 (94%)	31 (5%)	2 (0%)	36	50
All	All	1806/1815 (100%)	1708 (95%)	90 (5%)	8 (0%)	30	43

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	475	GLY
1	B	61	TYR
1	A	173	ASN
1	B	173	ASN
1	B	379	GLN
1	C	458	ASP
1	B	474	ALA
1	C	71	ALA

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	489/484 (101%)	474 (97%)	15 (3%)	35	57
1	B	489/484 (101%)	477 (98%)	12 (2%)	42	64
1	C	489/484 (101%)	470 (96%)	19 (4%)	28	48
All	All	1467/1452 (101%)	1421 (97%)	46 (3%)	35	57

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	42	LEU
1	A	57	GLU
1	A	68	VAL
1	A	84	THR
1	A	89	GLU
1	A	105	GLN
1	A	208	LYS
1	A	243	MSE
1	A	264	GLU
1	A	373	GLN
1	A	383	LEU
1	A	420	THR
1	A	483	ASP
1	A	497	LYS
1	A	630	GLN
1	B	56	LEU
1	B	60	ASN
1	B	89	GLU
1	B	105	GLN
1	B	164	VAL
1	B	176	VAL
1	B	264	GLU
1	B	342	ASN
1	B	373	GLN
1	B	379	GLN
1	B	497	LYS

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Mol	Chain	Res	Type
1	B	630	GLN
1	C	41	LYS
1	C	56	LEU
1	C	60	ASN
1	C	72	ASN
1	C	86	THR
1	C	109	LEU
1	C	125	GLU
1	C	164	VAL
1	C	264	GLU
1	C	342	ASN
1	C	397	LYS
1	C	437	LEU
1	C	483	ASP
1	C	490	LYS
1	C	497	LYS
1	C	510	LEU
1	C	521	GLU
1	C	589	SER
1	C	630	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (39) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	ASN
1	A	87	ASN
1	A	94	ASN
1	A	101	ASN
1	A	105	GLN
1	A	179	HIS
1	A	293	ASN
1	A	327	ASN
1	A	393	ASN
1	A	416	GLN
1	A	514	GLN
1	A	539	ASN
1	A	551	ASN
1	A	630	GLN
1	B	64	ASN
1	B	101	ASN
1	B	105	GLN
1	B	179	HIS

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Mol	Chain	Res	Type
1	B	214	HIS
1	B	293	ASN
1	B	327	ASN
1	B	393	ASN
1	B	459	ASN
1	B	491	GLN
1	B	539	ASN
1	B	630	GLN
1	C	60	ASN
1	C	87	ASN
1	C	145	HIS
1	C	179	HIS
1	C	215	GLN
1	C	293	ASN
1	C	373	GLN
1	C	393	ASN
1	C	459	ASN
1	C	476	GLN
1	C	514	GLN
1	C	531	ASN
1	C	630	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 9 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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, 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	599/605 (99%)	0.60	28 (4%)	36	32	31, 47, 67, 91	0
1	B	599/605 (99%)	0.69	35 (5%)	29	25	35, 50, 73, 95	0
1	C	599/605 (99%)	0.97	40 (6%)	24	20	43, 59, 76, 101	0
All	All	1797/1815 (99%)	0.75	103 (5%)	29	25	31, 52, 74, 101	0

All (103) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	61	TYR	11.3
1	C	631	THR	6.6
1	B	632	GLY	6.5
1	A	62	ALA	6.4
1	A	632	GLY	6.2
1	C	61	TYR	6.0
1	A	631	THR	6.0
1	B	62	ALA	5.7
1	B	61	TYR	5.4
1	C	632	GLY	5.3
1	C	630	GLN	5.0
1	B	631	THR	4.6
1	C	62	ALA	4.6
1	A	29	GLY	3.7
1	B	630	GLN	3.7
1	C	546	ASP	3.6
1	A	63	GLY	3.6
1	C	613	SER	3.4
1	C	90	ILE	3.3
1	A	46	ASP	3.2
1	A	73	ALA	3.1
1	A	47	GLY	3.1
1	B	594	GLU	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	105	GLN	3.0
1	A	96	GLN	2.9
1	B	221	GLU	2.9
1	C	361	ALA	2.8
1	C	63	GLY	2.8
1	C	440	THR	2.8
1	A	487	GLN	2.8
1	A	71	ALA	2.7
1	B	209	ASP	2.7
1	B	474	ALA	2.7
1	B	396	ASP	2.7
1	A	342	ASN	2.7
1	C	174	ASP	2.6
1	C	76	ILE	2.6
1	C	43	LEU	2.6
1	C	474	ALA	2.6
1	A	488	GLU	2.6
1	C	156	ALA	2.6
1	A	72	ASN	2.6
1	C	473	ASN	2.6
1	A	630	GLN	2.6
1	B	63	GLY	2.5
1	C	56	LEU	2.5
1	B	512	ALA	2.5
1	C	352	VAL	2.5
1	B	611	ALA	2.5
1	B	388	ASN	2.4
1	A	509	ALA	2.4
1	B	44	ALA	2.4
1	C	66	VAL	2.4
1	C	268	TYR	2.4
1	C	58	THR	2.4
1	A	174	ASP	2.3
1	C	176	VAL	2.3
1	C	293	ASN	2.3
1	B	498	ASP	2.3
1	B	624	LYS	2.3
1	B	93	GLU	2.3
1	A	283	ASP	2.3
1	C	496	THR	2.3
1	C	627	THR	2.3
1	C	172	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	586	LEU	2.2
1	A	501	ASP	2.2
1	B	204	PRO	2.2
1	C	618	ALA	2.2
1	C	129	PHE	2.2
1	A	60	ASN	2.2
1	B	291	ASP	2.2
1	A	612	GLY	2.2
1	A	89	GLU	2.2
1	C	585	ALA	2.2
1	B	56	LEU	2.2
1	B	206	VAL	2.2
1	A	518	GLN	2.2
1	B	47	GLY	2.2
1	C	362	GLY	2.2
1	C	365	ALA	2.2
1	C	628	ILE	2.1
1	B	60	ASN	2.1
1	B	501	ASP	2.1
1	B	514	GLN	2.1
1	B	469	ALA	2.1
1	C	141	ILE	2.1
1	A	48	ASP	2.1
1	B	216	VAL	2.1
1	C	519	GLU	2.1
1	B	29	GLY	2.1
1	C	155	GLY	2.1
1	A	30	ALA	2.1
1	A	387	LYS	2.1
1	B	30	ALA	2.1
1	B	343	GLY	2.1
1	C	310	ALA	2.0
1	B	105	GLN	2.0
1	B	612	GLY	2.0
1	C	157	THR	2.0
1	B	503	ALA	2.0
1	C	71	ALA	2.0
1	B	491	GLN	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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($\text{\AA}^2$)	Q<0.9
2	NI	B	802	1/1	0.94	0.07	82,82,82,82	0
2	NI	C	802	1/1	0.95	0.07	78,78,78,78	0
2	NI	B	801	1/1	0.96	0.11	78,78,78,78	0
2	NI	A	801	1/1	0.96	0.08	78,78,78,78	0
2	NI	C	801	1/1	0.96	0.07	74,74,74,74	0
2	NI	A	802	1/1	0.96	0.06	77,77,77,77	0
2	NI	B	800	1/1	0.97	0.12	59,59,59,59	0
2	NI	A	800	1/1	0.98	0.04	62,62,62,62	0
2	NI	C	800	1/1	0.98	0.06	74,74,74,74	0

6.5 Other polymers [i](#)

There are no such residues in this entry.