



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 5, 2026 – 12:06 PM UTC

PDB ID : 1VQQ / pdb_00001vqq
Title : Structure of Penicillin binding protein 2a from methicillin resistant Staphylococcus aureus strain 27r at 1.80 Å resolution.
Authors : Lim, D.; Strynadka, N.C.J.
Deposited on : 2004-12-17
Resolution : 1.80 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

<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
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

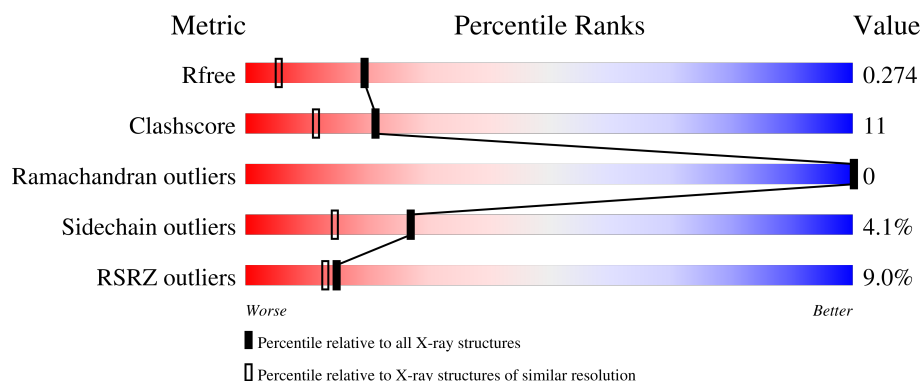
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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}	180053	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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 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	646	9% 73% 23% ..
1	B	646	9% 77% 18% ..

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10644 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 Beta-lactam-inducible penicillin-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	635	Total	C	N	O	S	0	0	0
			5102	3218	860	1009	15			
1	B	625	Total	C	N	O	S	0	0	0
			5023	3171	843	994	15			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	23	MET	TYR	engineered mutation	UNP O54286
B	23	MET	TYR	engineered mutation	UNP O54286

- Molecule 2 is CADMIUM ION (CCD ID: CD) (formula: Cd).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Cd	0	0
			2	2		
2	B	5	Total	Cd	0	0
			5	5		

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Cl	0	0
			1	1		
3	B	3	Total	Cl	0	0
			3	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	246	Total	O	0	0
			246	246		

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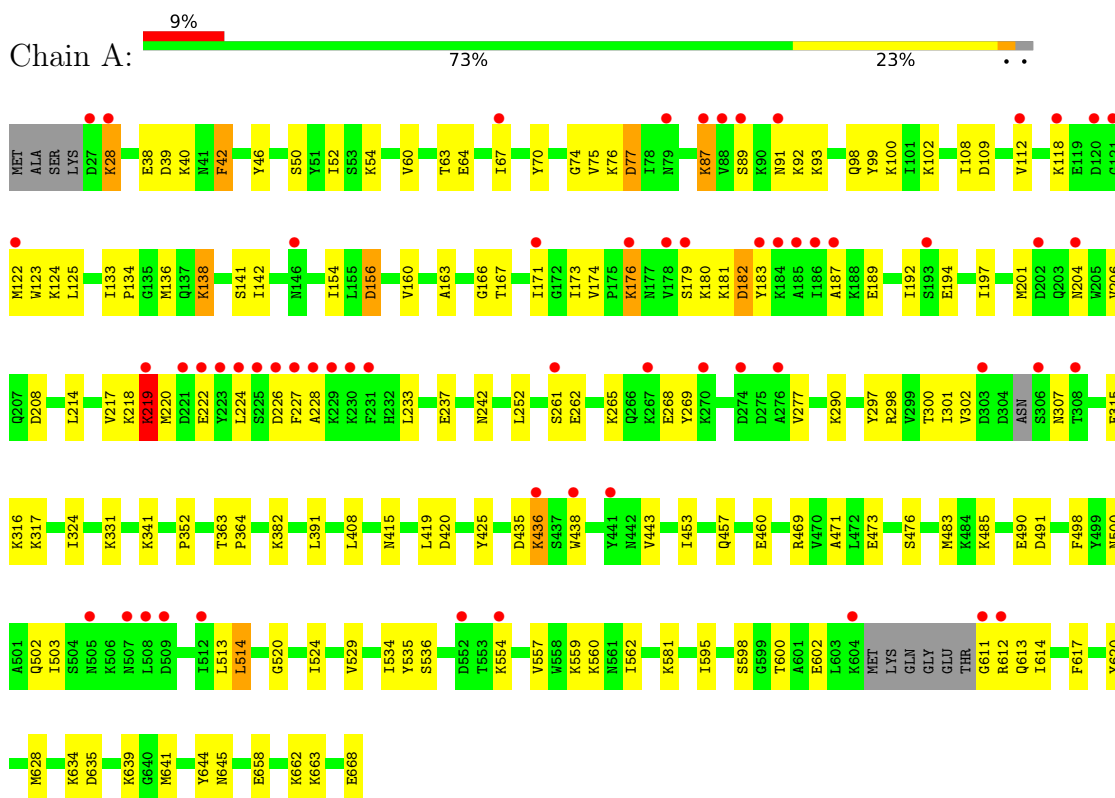
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	262	Total 262	O 262	0	0

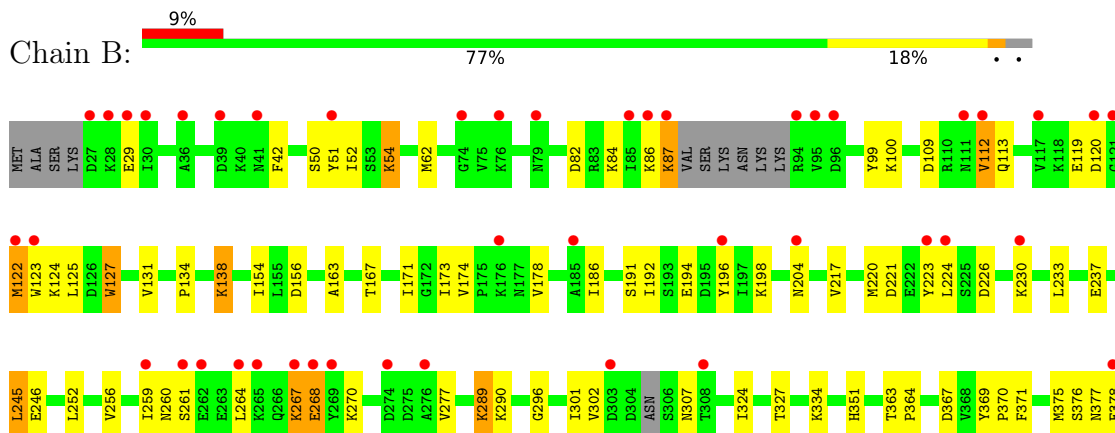
3 Residue-property plots

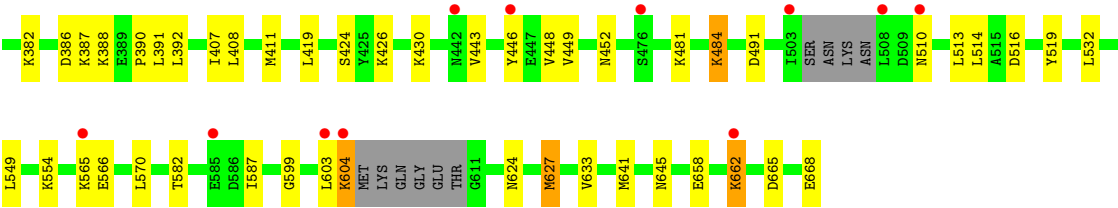
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Beta-lactam-inducible penicillin-binding protein



- Molecule 1: Beta-lactam-inducible penicillin-binding protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	80.87Å 100.61Å 186.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.92 – 1.80 24.92 – 1.80	Depositor EDS
% Data completeness (in resolution range)	97.3 (24.92-1.80) 97.4 (24.92-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.44 (at 1.80Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.237 , 0.274 0.237 , 0.274	Depositor DCC
R_{free} test set	6974 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	23.9	Xtriage
Anisotropy	0.804	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 40.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10644	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.47% 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: CL, CD

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	0.56	0/5186	0.98	14/6971 (0.2%)
1	B	0.54	0/5105	0.99	21/6862 (0.3%)
All	All	0.55	0/10291	0.99	35/13833 (0.3%)

There are no bond length outliers.

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	174	VAL	N-CA-C	-10.51	96.35	107.60
1	B	599	GLY	N-CA-C	9.69	123.15	112.29
1	B	174	VAL	N-CA-C	-9.58	97.35	107.60
1	A	341	LYS	N-CA-C	7.57	121.44	111.75
1	B	351	HIS	CA-C-N	7.20	127.53	119.32
1	B	351	HIS	C-N-CA	7.20	127.53	119.32
1	A	219	LYS	N-CA-C	6.87	120.72	109.72
1	B	443	VAL	N-CA-C	-6.60	100.17	108.89
1	B	246	GLU	CB-CA-C	-6.40	109.21	116.63
1	A	218	LYS	N-CA-C	6.26	117.77	111.07
1	B	173	ILE	N-CA-C	5.98	117.10	108.48
1	B	178	VAL	N-CA-C	5.94	116.43	108.11
1	B	296	GLY	N-CA-C	-5.79	104.82	112.81
1	B	42	PHE	N-CA-C	5.73	117.52	111.28
1	A	301	ILE	N-CA-C	-5.69	99.54	107.80
1	B	204	ASN	N-CA-C	5.69	118.25	111.71
1	A	52	ILE	N-CA-C	5.64	116.42	110.72
1	B	131	VAL	N-CA-C	-5.61	104.70	112.50
1	A	206	VAL	N-CA-C	5.52	116.13	108.84
1	A	194	GLU	N-CA-C	-5.49	105.38	111.36
1	B	624	ASN	CA-C-N	5.41	126.60	119.84
1	B	624	ASN	C-N-CA	5.41	126.60	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	443	VAL	N-CA-C	-5.35	101.83	108.89
1	B	491	ASP	N-CA-C	-5.34	102.37	110.28
1	A	520	GLY	N-CA-C	5.31	123.18	115.32
1	B	641	MET	CB-CA-C	-5.27	110.52	116.63
1	A	160	VAL	N-CA-C	-5.27	102.72	109.30
1	B	301	ILE	N-CA-C	-5.19	100.28	107.80
1	B	246	GLU	N-CA-C	5.18	116.84	108.08
1	B	127	TRP	N-CA-C	5.17	117.73	110.23
1	A	476	SER	N-CA-C	5.16	117.30	111.11
1	B	52	ILE	N-CA-C	5.08	115.85	110.72
1	A	42	PHE	N-CA-C	5.03	116.76	111.28
1	A	438	TRP	N-CA-C	-5.02	106.94	113.16
1	B	582	THR	N-CA-C	5.02	117.13	111.11

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5102	0	5099	125	0
1	B	5023	0	5008	96	0
2	A	2	0	0	0	0
2	B	5	0	0	0	0
3	A	1	0	0	0	0
3	B	3	0	0	0	0
4	A	246	0	0	5	0
4	B	262	0	0	6	0
All	All	10644	0	10107	220	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 11.

All (220) 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:87:LYS:HE3	1:A:87:LYS:H	1.14	1.08
1:B:267:LYS:H	1:B:267:LYS:HD3	1.21	1.06
1:A:611:GLY:HA3	1:A:635:ASP:OD1	1.56	1.04
1:A:436:LYS:H	1:A:436:LYS:HD3	1.23	1.03
1:A:136:MET:HE2	1:A:142:ILE:HD11	1.42	0.99
1:A:133:ILE:HB	1:A:136:MET:HE3	1.49	0.94
1:B:267:LYS:HD3	1:B:267:LYS:N	1.89	0.86
1:A:87:LYS:H	1:A:87:LYS:CE	1.92	0.82
1:B:87:LYS:HD3	1:B:87:LYS:N	1.93	0.82
1:A:138:LYS:H	1:A:138:LYS:HD2	1.43	0.81
1:A:122:MET:HG3	1:A:124:LYS:NZ	1.95	0.81
1:B:226:ASP:O	1:B:230:LYS:HG2	1.82	0.79
1:A:28:LYS:HD2	1:A:28:LYS:N	1.97	0.79
1:A:87:LYS:HE3	1:A:87:LYS:N	1.96	0.78
1:B:510:ASN:HB3	1:B:513:LEU:HD12	1.65	0.77
1:A:436:LYS:H	1:A:436:LYS:CD	1.96	0.77
1:A:70:TYR:HB3	1:A:75:VAL:CG2	2.15	0.76
1:A:138:LYS:H	1:A:138:LYS:CD	1.99	0.74
1:B:387:LYS:HA	1:B:387:LYS:HE2	1.69	0.73
1:A:70:TYR:HB3	1:A:75:VAL:HG21	1.69	0.73
1:A:557:VAL:HG11	1:A:560:LYS:HG2	1.71	0.73
1:A:112:VAL:CG1	1:A:134:PRO:HB3	2.19	0.72
1:A:485:LYS:HE3	4:A:1170:HOH:O	1.90	0.72
1:A:613:GLN:HG2	1:A:641:MET:SD	2.30	0.71
1:B:256:VAL:HG22	4:B:1059:HOH:O	1.90	0.71
1:B:426:LYS:HG2	1:B:452:ASN:OD1	1.91	0.70
1:B:267:LYS:HA	1:B:270:LYS:HE3	1.73	0.70
1:B:122:MET:HG3	1:B:124:LYS:HE2	1.73	0.70
1:A:77:ASP:HB2	1:A:102:LYS:HD2	1.74	0.68
1:B:138:LYS:HE2	1:B:138:LYS:H	1.58	0.68
1:B:378:GLU:HG3	1:B:382:LYS:NZ	2.08	0.68
1:B:112:VAL:CG1	1:B:134:PRO:HB3	2.25	0.67
1:B:603:LEU:HD23	1:B:604:LYS:N	2.11	0.66
1:A:268:GLU:HG3	1:A:269:TYR:CD1	2.31	0.65
1:A:176:LYS:HG3	1:A:208:ASP:O	1.98	0.64
1:B:167:THR:HG23	1:B:237:GLU:HG3	1.80	0.64
1:A:171:ILE:HD11	1:A:217:VAL:HG13	1.79	0.64
1:B:112:VAL:HG13	1:B:134:PRO:HB3	1.80	0.64
1:A:436:LYS:HD3	1:A:436:LYS:N	2.06	0.63
1:A:614:ILE:HD11	1:A:634:LYS:HG3	1.79	0.63
1:A:171:ILE:HD11	1:A:217:VAL:CG1	2.29	0.63
1:A:133:ILE:HB	1:A:136:MET:CE	2.27	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:176:LYS:HD2	1:B:113:GLN:NE2	2.15	0.62
1:B:627:MET:O	1:B:627:MET:HG3	1.98	0.62
1:A:420:ASP:OD1	4:A:1011:HOH:O	2.16	0.62
1:B:100:LYS:HG2	1:B:109:ASP:OD2	2.00	0.62
1:A:138:LYS:HD2	1:A:138:LYS:N	2.12	0.61
1:A:598:SER:HB3	1:A:617:PHE:CD1	2.36	0.61
1:B:220:MET:HE2	1:B:224:LEU:HD23	1.82	0.61
1:A:602:GLU:OE1	1:A:613:GLN:NE2	2.29	0.61
1:B:29:GLU:OE1	1:B:123:TRP:HD1	1.84	0.61
1:A:408:LEU:HD22	1:A:534:ILE:HG21	1.83	0.61
1:B:99:TYR:HD1	1:B:112:VAL:HG11	1.66	0.60
1:A:46:TYR:O	1:A:54:LYS:HD3	2.01	0.60
1:B:186:ILE:HD13	1:B:233:LEU:HD21	1.84	0.60
1:B:226:ASP:HB3	1:B:230:LYS:HE3	1.83	0.60
1:B:220:MET:CE	1:B:224:LEU:HD23	2.32	0.60
1:A:261:SER:O	1:A:265:LYS:HE3	2.02	0.59
1:A:658:GLU:HB2	1:A:662:LYS:HG2	1.85	0.59
1:B:50:SER:O	1:B:54:LYS:HD2	2.02	0.59
1:B:261:SER:HA	1:B:264:LEU:HD12	1.84	0.59
1:A:503:ILE:O	1:A:524:ILE:HG12	2.04	0.58
1:B:277:VAL:HG23	4:B:1057:HOH:O	2.04	0.58
1:B:370:PRO:HB2	1:B:375:MET:HE2	1.84	0.58
1:A:290:LYS:HG3	4:A:1144:HOH:O	2.03	0.58
1:B:378:GLU:HG3	1:B:382:LYS:HZ3	1.69	0.57
1:B:122:MET:HE2	1:B:124:LYS:NZ	2.18	0.57
1:B:260:ASN:O	1:B:264:LEU:HG	2.05	0.57
1:B:371:PHE:CE1	1:B:375:MET:HE1	2.40	0.57
1:B:449:VAL:HG23	1:B:449:VAL:O	2.04	0.56
1:A:89:SER:OG	1:A:92:LYS:HB3	2.05	0.56
1:A:612:ARG:HG3	1:A:612:ARG:HH11	1.69	0.56
1:A:70:TYR:HB3	1:A:75:VAL:HG22	1.87	0.56
1:A:425:TYR:OH	1:A:473:GLU:HG3	2.06	0.56
1:B:86:LYS:HG2	1:B:87:LYS:N	2.20	0.56
1:B:125:LEU:HD13	1:B:127:TRP:N	2.21	0.56
1:A:28:LYS:N	1:A:28:LYS:CD	2.69	0.56
1:A:112:VAL:HG11	1:A:134:PRO:HB3	1.86	0.55
1:B:554:LYS:HG2	4:B:1196:HOH:O	2.06	0.55
1:A:219:LYS:H	1:A:219:LYS:HD2	1.71	0.55
1:A:277:VAL:HG23	4:A:1055:HOH:O	2.06	0.55
1:A:262:GLU:H	1:A:262:GLU:CD	2.13	0.54
1:A:290:LYS:HB2	1:A:324:ILE:HD11	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:460:GLU:OE1	1:A:581:LYS:HE3	2.07	0.54
1:B:82:ASP:OD1	1:B:82:ASP:O	2.26	0.54
1:B:220:MET:HE2	1:B:220:MET:HA	1.90	0.54
1:B:217:VAL:HG21	1:B:220:MET:HE3	1.90	0.53
1:B:221:ASP:OD2	1:B:224:LEU:HB2	2.08	0.53
1:A:658:GLU:HG3	1:A:662:LYS:HD3	1.90	0.53
1:A:302:VAL:CG1	1:A:307:ASN:HA	2.39	0.53
1:A:122:MET:HG3	1:A:124:LYS:HZ1	1.71	0.53
1:B:484:LYS:NZ	1:B:484:LYS:HB2	2.22	0.53
1:A:614:ILE:CD1	1:A:634:LYS:HA	2.39	0.52
1:A:639:LYS:HD3	1:A:644:TYR:CE1	2.44	0.52
1:B:82:ASP:O	1:B:84:LYS:HG3	2.09	0.52
1:A:298:ARG:HD3	1:A:315:GLU:OE1	2.10	0.52
1:A:189:GLU:HG2	1:A:227:PHE:CE1	2.45	0.52
1:A:614:ILE:CD1	1:A:634:LYS:HG3	2.39	0.52
1:A:217:VAL:HG21	1:A:224:LEU:HD13	1.92	0.51
1:B:87:LYS:HD3	1:B:87:LYS:H	1.72	0.51
1:B:86:LYS:CG	1:B:87:LYS:N	2.73	0.51
1:B:424:SER:OG	1:B:452:ASN:HB3	2.10	0.51
1:A:64:GLU:O	1:A:67:ILE:HB	2.10	0.51
1:B:119:GLU:O	1:B:120:ASP:C	2.53	0.51
1:A:536:SER:HA	1:A:628:MET:CE	2.41	0.50
1:B:99:TYR:HB2	1:B:112:VAL:CG2	2.41	0.50
1:B:386:ASP:HB3	1:B:390:PRO:HD3	1.92	0.50
1:B:516:ASP:HA	1:B:519:TYR:CE2	2.46	0.50
1:A:70:TYR:CB	1:A:75:VAL:HG21	2.40	0.50
1:A:180:LYS:HA	1:A:183:TYR:CE1	2.47	0.50
1:A:268:GLU:HG3	1:A:269:TYR:CE1	2.47	0.50
1:A:598:SER:HB3	1:A:617:PHE:CE1	2.47	0.50
1:B:565:LYS:HE2	4:B:1262:HOH:O	2.11	0.50
1:B:125:LEU:CD1	1:B:127:TRP:HA	2.42	0.49
1:B:658:GLU:HG3	1:B:662:LYS:HE2	1.94	0.49
1:A:60:VAL:HG13	1:A:64:GLU:OE1	2.12	0.49
1:B:363:THR:HA	1:B:364:PRO:C	2.37	0.49
1:B:587:ILE:HG12	1:B:587:ILE:O	2.12	0.49
1:B:390:PRO:HD2	4:B:1085:HOH:O	2.10	0.49
1:A:189:GLU:HG2	1:A:227:PHE:CZ	2.48	0.49
1:A:220:MET:SD	1:A:224:LEU:HD23	2.52	0.49
1:B:192:ILE:HD12	1:B:196:TYR:HD2	1.78	0.49
1:A:98:GLN:NE2	1:A:109:ASP:OD2	2.42	0.49
1:A:180:LYS:HA	1:A:183:TYR:CD1	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:658:GLU:CD	1:B:662:LYS:HG2	2.38	0.48
1:B:86:LYS:CG	1:B:87:LYS:H	2.27	0.48
1:A:483:MET:HE2	1:A:524:ILE:HD13	1.95	0.48
1:B:289:LYS:HB2	1:B:289:LYS:NZ	2.28	0.48
1:B:194:GLU:O	1:B:198:LYS:HG3	2.14	0.48
1:A:261:SER:HB2	1:A:262:GLU:OE1	2.14	0.48
1:A:277:VAL:HG23	1:A:277:VAL:O	2.14	0.48
1:A:352:PRO:HA	1:A:536:SER:HB2	1.95	0.48
1:A:363:THR:HA	1:A:364:PRO:C	2.38	0.48
1:B:369:TYR:HB2	1:B:370:PRO:HD3	1.96	0.48
1:B:191:SER:HB3	1:B:376:SER:HB3	1.96	0.47
1:B:378:GLU:HG3	1:B:382:LYS:HZ1	1.80	0.47
1:B:604:LYS:HA	1:B:604:LYS:HE3	1.96	0.47
1:B:665:ASP:HB3	1:B:668:GLU:HB2	1.97	0.47
1:B:668:GLU:OXT	1:B:668:GLU:HG2	2.15	0.47
1:A:222:GLU:O	1:A:226:ASP:HB2	2.14	0.47
1:B:364:PRO:HG2	1:B:388:LYS:HD3	1.96	0.47
1:A:166:GLY:HA3	1:A:242:ASN:HB2	1.97	0.46
1:A:471:ALA:HB1	1:A:514:LEU:HD22	1.97	0.46
1:B:138:LYS:H	1:B:138:LYS:CE	2.26	0.46
1:A:182:ASP:OD1	1:A:182:ASP:N	2.49	0.46
1:A:133:ILE:CB	1:A:136:MET:HE3	2.32	0.46
1:A:173:ILE:CD1	1:A:214:LEU:HD11	2.45	0.46
1:B:290:LYS:HB2	1:B:324:ILE:HD11	1.98	0.46
1:A:187:ALA:HB1	1:A:192:ILE:O	2.16	0.46
1:A:197:ILE:O	1:A:201:MET:HG2	2.16	0.46
1:A:93:LYS:HE3	1:A:123:TRP:CZ2	2.51	0.45
1:B:387:LYS:HA	1:B:387:LYS:CE	2.41	0.45
1:A:167:THR:CG2	1:A:237:GLU:HG3	2.47	0.45
1:B:62:MET:HE3	1:B:127:TRP:CG	2.51	0.45
1:B:510:ASN:CB	1:B:513:LEU:HD12	2.40	0.45
1:B:407:ILE:O	1:B:411:MET:HG3	2.15	0.45
1:B:302:VAL:CG1	1:B:307:ASN:HA	2.47	0.45
1:B:603:LEU:HD23	1:B:603:LEU:C	2.41	0.45
1:A:262:GLU:OE1	1:A:262:GLU:N	2.40	0.45
1:A:614:ILE:HD12	1:A:634:LYS:HA	1.98	0.45
1:B:566:GLU:HG2	4:B:1188:HOH:O	2.16	0.45
1:A:297:TYR:CE1	1:A:316:LYS:HD3	2.51	0.45
1:A:100:LYS:HA	1:A:108:ILE:O	2.17	0.45
1:A:595:ILE:HD11	1:A:620:TYR:CZ	2.52	0.44
1:B:54:LYS:HB3	1:B:54:LYS:HE2	1.73	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:39:ASP:O	1:A:40:LYS:HB2	2.17	0.44
1:A:179:SER:C	1:A:181:LYS:H	2.26	0.44
1:A:382:LYS:HD3	1:A:382:LYS:C	2.42	0.44
1:B:245:LEU:HD13	1:B:334:LYS:HG3	1.99	0.44
1:A:557:VAL:HG11	1:A:560:LYS:CG	2.45	0.44
1:B:267:LYS:H	1:B:267:LYS:CD	2.08	0.44
1:B:125:LEU:HD11	1:B:127:TRP:HA	2.00	0.44
1:A:112:VAL:CG1	1:A:134:PRO:CB	2.94	0.43
1:A:490:GLU:OE1	1:A:559:LYS:NZ	2.39	0.43
1:B:267:LYS:HG2	1:B:268:GLU:HG2	1.99	0.43
1:A:138:LYS:H	1:A:138:LYS:CE	2.30	0.43
1:A:154:ILE:HB	1:A:163:ALA:HB3	2.00	0.43
1:A:559:LYS:HB3	1:A:562:ILE:HD11	2.00	0.43
1:B:217:VAL:CG2	1:B:220:MET:HE3	2.48	0.43
1:A:415:ASN:HD21	1:A:485:LYS:HE2	1.84	0.43
1:A:498:PHE:CE2	1:A:529:VAL:HG21	2.53	0.43
1:A:658:GLU:CG	1:A:662:LYS:HD3	2.49	0.43
1:A:99:TYR:HD1	1:A:112:VAL:HG11	1.84	0.43
1:A:298:ARG:HG2	1:A:300:THR:HG23	2.01	0.43
1:B:171:ILE:HD11	1:B:220:MET:HE1	2.00	0.43
1:A:70:TYR:CG	1:A:75:VAL:HG21	2.54	0.43
1:B:221:ASP:CG	1:B:223:TYR:HD2	2.26	0.43
1:A:93:LYS:HG3	1:A:123:TRP:CH2	2.54	0.42
1:B:167:THR:CG2	1:B:237:GLU:HG3	2.47	0.42
1:A:112:VAL:HG12	1:A:134:PRO:CB	2.49	0.42
1:A:453:ILE:HG23	1:A:457:GLN:CG	2.49	0.42
1:A:645:ASN:H	1:A:645:ASN:HD22	1.67	0.42
1:B:51:TYR:O	1:B:54:LYS:HG2	2.20	0.42
1:B:430:LYS:HG3	1:B:448:VAL:CG2	2.49	0.42
1:A:74:GLY:O	1:A:76:LYS:HD3	2.19	0.42
1:A:535:TYR:O	1:A:628:MET:HE1	2.20	0.42
1:A:50:SER:O	1:A:54:LYS:HG3	2.20	0.42
1:A:156:ASP:OD2	1:A:156:ASP:C	2.63	0.42
1:B:82:ASP:OD1	1:B:84:LYS:HD2	2.20	0.42
1:A:415:ASN:ND2	1:A:485:LYS:HE2	2.35	0.41
1:B:154:ILE:HB	1:B:163:ALA:HB3	2.02	0.41
1:A:138:LYS:HB3	1:A:138:LYS:HE3	1.86	0.41
1:A:435:ASP:HB2	1:A:436:LYS:HD3	2.02	0.41
1:A:112:VAL:HG12	1:A:134:PRO:HB3	2.01	0.41
1:A:502:GLN:O	1:A:502:GLN:HG3	2.19	0.41
1:A:222:GLU:OE2	1:A:222:GLU:HA	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:265:LYS:HD3	1:A:265:LYS:N	2.36	0.41
1:B:259:ILE:HG12	1:B:260:ASN:N	2.35	0.41
1:B:633:VAL:CG2	1:B:645:ASN:HD21	2.34	0.41
1:A:331:LYS:NZ	1:A:668:GLU:OXT	2.47	0.41
1:B:367:ASP:OD1	1:B:369:TYR:CD1	2.74	0.41
1:A:42:PHE:HB3	1:A:63:THR:HA	2.03	0.41
1:A:228:ALA:HA	1:A:233:LEU:HB2	2.02	0.41
1:B:327:THR:OG1	1:B:549:LEU:HA	2.22	0.40
1:A:663:LYS:HE2	4:A:1216:HOH:O	2.20	0.40
1:A:600:THR:HG21	1:A:641:MET:HB3	2.03	0.40
1:B:186:ILE:CD1	1:B:233:LEU:HD21	2.50	0.40
1:A:91:ASN:ND2	1:A:118:LYS:HB3	2.36	0.40
1:A:491:ASP:OD1	1:A:500:ASN:ND2	2.54	0.40
1:B:122:MET:HE2	1:B:124:LYS:HZ3	1.84	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	629/646 (97%)	612 (97%)	17 (3%)	0	100	100
1	B	615/646 (95%)	595 (97%)	20 (3%)	0	100	100
All	All	1244/1292 (96%)	1207 (97%)	37 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	567/576 (98%)	546 (96%)	21 (4%)	30	18
1	B	557/576 (97%)	532 (96%)	25 (4%)	24	12
All	All	1124/1152 (98%)	1078 (96%)	46 (4%)	27	15

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

Mol	Chain	Res	Type
1	A	28	LYS
1	A	38	GLU
1	A	77	ASP
1	A	87	LYS
1	A	125	LEU
1	A	138	LYS
1	A	141	SER
1	A	156	ASP
1	A	176	LYS
1	A	182	ASP
1	A	204	ASN
1	A	219	LYS
1	A	252	LEU
1	A	317	LYS
1	A	391	LEU
1	A	419	LEU
1	A	436	LYS
1	A	469	ARG
1	A	513	LEU
1	A	514	LEU
1	A	554	LYS
1	B	54	LYS
1	B	87	LYS
1	B	112	VAL
1	B	122	MET
1	B	138	LYS
1	B	156	ASP
1	B	245	LEU
1	B	252	LEU
1	B	267	LYS
1	B	268	GLU
1	B	289	LYS
1	B	377	ASN

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Mol	Chain	Res	Type
1	B	391	LEU
1	B	392	LEU
1	B	408	LEU
1	B	419	LEU
1	B	446	TYR
1	B	481	LYS
1	B	484	LYS
1	B	514	LEU
1	B	532	LEU
1	B	570	LEU
1	B	604	LYS
1	B	627	MET
1	B	662	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	ASN
1	A	79	ASN
1	A	81	GLN
1	A	91	ASN
1	A	113	GLN
1	A	115	ASN
1	A	159	ASN
1	A	203	GLN
1	A	207	GLN
1	A	342	ASN
1	A	415	ASN
1	A	555	ASN
1	A	645	ASN
1	A	659	ASN
1	B	32	ASN
1	B	81	GLN
1	B	104	ASN
1	B	111	ASN
1	B	159	ASN
1	B	177	ASN
1	B	377	ASN
1	B	415	ASN
1	B	433	GLN
1	B	502	GLN
1	B	645	ASN

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Mol	Chain	Res	Type
1	B	659	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 11 ligands modelled in this entry, 11 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 [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

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	635/646 (98%)	0.59	59 (9%) 14 12	17, 31, 51, 62	0
1	B	625/646 (96%)	0.53	55 (8%) 15 13	19, 31, 53, 62	0
All	All	1260/1292 (97%)	0.56	114 (9%) 15 13	17, 31, 52, 62	0

All (114) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	223	TYR	7.8
1	A	611	GLY	6.3
1	B	223	TYR	6.3
1	A	120	ASP	5.4
1	A	228	ALA	5.1
1	A	226	ASP	3.8
1	B	264	LEU	3.8
1	A	227	PHE	3.8
1	A	185	ALA	3.7
1	B	230	LYS	3.7
1	B	123	TRP	3.7
1	B	261	SER	3.6
1	B	122	MET	3.5
1	B	120	ASP	3.5
1	A	183	TYR	3.4
1	B	262	GLU	3.4
1	B	603	LEU	3.3
1	A	121	GLY	3.3
1	B	27	ASP	3.2
1	A	267	LYS	3.2
1	B	87	LYS	3.1
1	B	121	GLY	3.1
1	A	178	VAL	3.1
1	A	204	ASN	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	222	GLU	3.1
1	A	184	LYS	3.0
1	A	229	LYS	3.0
1	B	503	ILE	2.9
1	A	27	ASP	2.9
1	B	446	TYR	2.9
1	B	604	LYS	2.8
1	A	179	SER	2.8
1	A	186	ILE	2.8
1	B	267	LYS	2.8
1	B	269	TYR	2.8
1	B	185	ALA	2.8
1	A	176	LYS	2.7
1	A	508	LEU	2.7
1	B	30	ILE	2.7
1	B	112	VAL	2.7
1	A	231	PHE	2.7
1	B	204	ASN	2.7
1	B	29	GLU	2.7
1	B	565	LYS	2.7
1	A	612	ARG	2.7
1	B	476	SER	2.7
1	B	95	VAL	2.7
1	B	274	ASP	2.6
1	B	268	GLU	2.6
1	A	261	SER	2.6
1	A	554	LYS	2.6
1	A	79	ASN	2.6
1	A	91	ASN	2.6
1	A	224	LEU	2.6
1	B	224	LEU	2.6
1	A	303	ASP	2.6
1	B	378	GLU	2.6
1	A	308	THR	2.6
1	A	274	ASP	2.5
1	A	118	LYS	2.5
1	A	438	TRP	2.5
1	B	96	ASP	2.5
1	A	436	LYS	2.4
1	A	112	VAL	2.4
1	A	306	SER	2.4
1	B	36	ALA	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	94	ARG	2.4
1	B	111	ASN	2.4
1	A	225	SER	2.4
1	A	171	ILE	2.4
1	B	117	VAL	2.4
1	A	87	LYS	2.4
1	B	303	ASP	2.3
1	A	187	ALA	2.3
1	A	276	ALA	2.3
1	A	219	LYS	2.3
1	B	51	TYR	2.3
1	A	604	LYS	2.3
1	B	28	LYS	2.3
1	B	259	ILE	2.3
1	A	88	VAL	2.3
1	A	146	ASN	2.3
1	B	442	ASN	2.3
1	A	28	LYS	2.3
1	A	221	ASP	2.2
1	B	196	TYR	2.2
1	A	512	ILE	2.2
1	A	507	ASN	2.2
1	B	508	LEU	2.2
1	A	202	ASP	2.2
1	A	509	ASP	2.2
1	B	85	ILE	2.2
1	B	308	THR	2.2
1	A	270	LYS	2.2
1	A	122	MET	2.1
1	B	41	ASN	2.1
1	A	441	TYR	2.1
1	A	67	ILE	2.1
1	B	39	ASP	2.1
1	B	176	LYS	2.1
1	B	265	LYS	2.1
1	B	662	LYS	2.1
1	B	74	GLY	2.1
1	B	76	LYS	2.1
1	B	276	ALA	2.1
1	A	552	ASP	2.1
1	B	510	ASN	2.1
1	A	193	SER	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	505	ASN	2.0
1	B	79	ASN	2.0
1	B	585	GLU	2.0
1	A	230	LYS	2.0
1	B	86	LYS	2.0
1	A	89	SER	2.0

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
2	CD	B	1006	1/1	0.92	0.15	57,57,57,57	0
3	CL	B	1011	1/1	0.95	0.06	27,27,27,27	0
2	CD	B	1007	1/1	0.98	0.05	31,31,31,31	0
3	CL	A	1009	1/1	0.98	0.07	27,27,27,27	0
3	CL	B	1008	1/1	0.98	0.04	27,27,27,27	0
3	CL	B	1010	1/1	0.98	0.05	29,29,29,29	0
2	CD	A	1004	1/1	0.98	0.11	50,50,50,50	0
2	CD	A	1005	1/1	0.99	0.02	27,27,27,27	0
2	CD	B	1003	1/1	0.99	0.03	27,27,27,27	0
2	CD	B	1002	1/1	1.00	0.02	24,24,24,24	0
2	CD	B	1001	1/1	1.00	0.02	25,25,25,25	0

6.5 Other polymers [i](#)

There are no such residues in this entry.