



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:35 PM GMT

PDB ID : 4CPK
Title : Crystal structure of PBP2a double clinical mutant N146K-E150K from MRSA
Authors : Otero, L.H.; Rojas-Altuve, A.; Hermoso, J.A.
Deposited on : 2014-02-07
Resolution : 2.35 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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

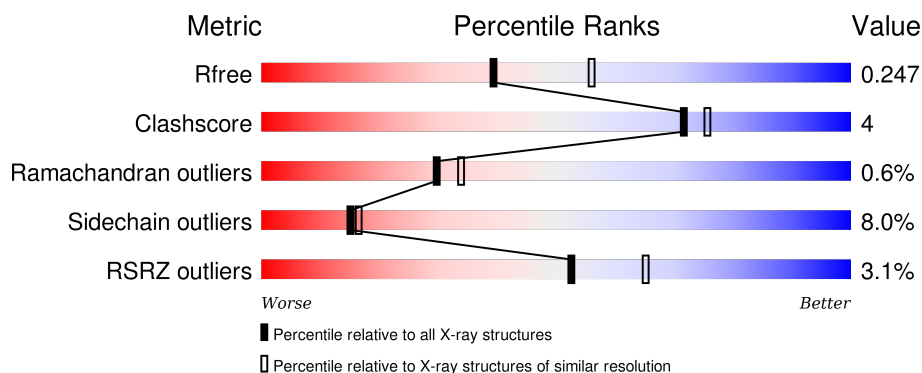
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

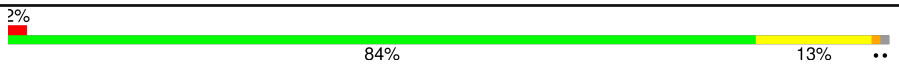

The reported resolution of this entry is 2.35 Å.

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}	91344	1352 (2.38-2.34)
Clashscore	102246	1456 (2.38-2.34)
Ramachandran outliers	100387	1435 (2.38-2.34)
Sidechain outliers	100360	1436 (2.38-2.34)
RSRZ outliers	91569	1358 (2.38-2.34)

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	643	
1	B	643	

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
4	MUR	A	1679	-	-	X	X
4	MUR	B	1677	-	-	X	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 11030 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 PENICILLIN BINDING PROTEIN 2 PRIME.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	637	Total	C	N	O	S	0	0	0
			5113	3226	864	1008	15			
1	B	636	Total	C	N	O	S	0	0	0
			5104	3220	862	1007	15			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	146	LYS	ASN	ENGINEERED MUTATION	UNP Q54113
A	150	LYS	GLU	ENGINEERED MUTATION	UNP Q54113
B	146	LYS	ASN	ENGINEERED MUTATION	UNP Q54113
B	150	LYS	GLU	ENGINEERED MUTATION	UNP Q54113

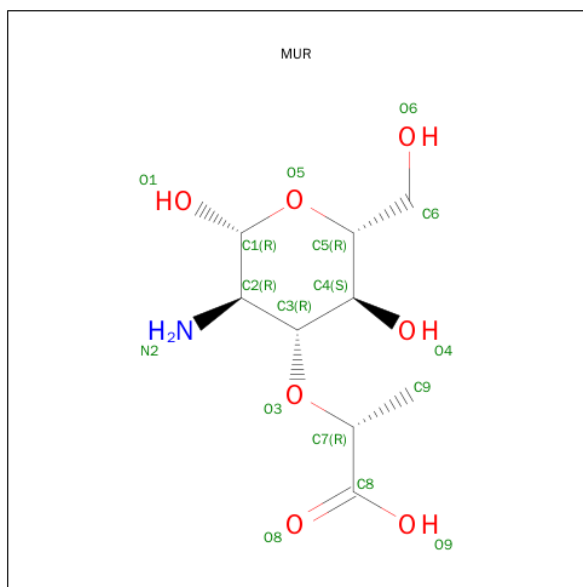
- Molecule 2 is CADMIUM ION (three-letter code: CD) (formula: Cd).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	4	Total	Cd	0	0
			4	4		
2	A	7	Total	Cd	0	0
			7	7		

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

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

- Molecule 4 is MURAMIC ACID (three-letter code: MUR) (formula: C₉H₁₇NO₇).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			17	9	1	7		
4	B	1	Total	C	N	O	0	0
			17	9	1	7		

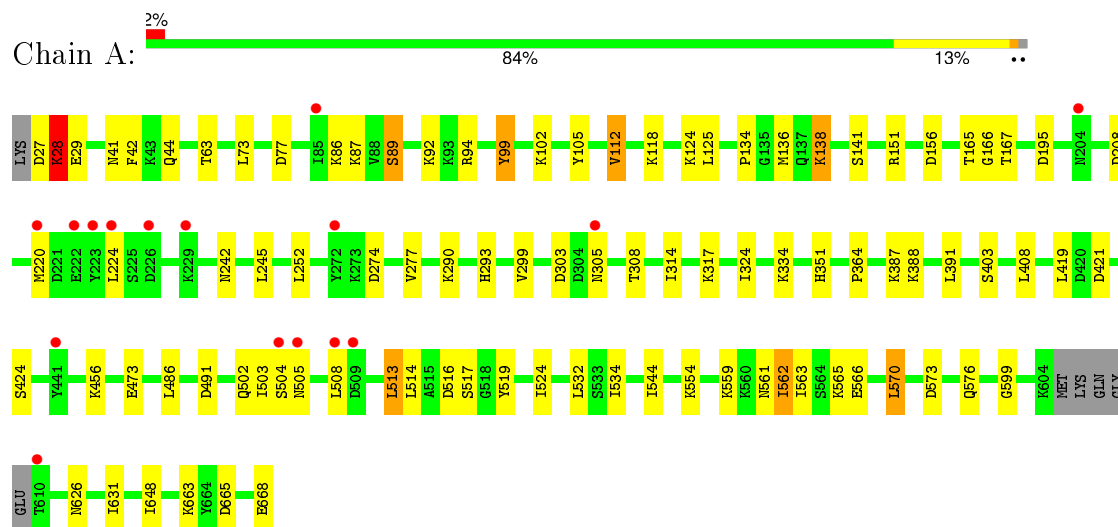
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	406	Total	O	0	0
			406	406		
5	B	358	Total	O	0	0
			358	358		

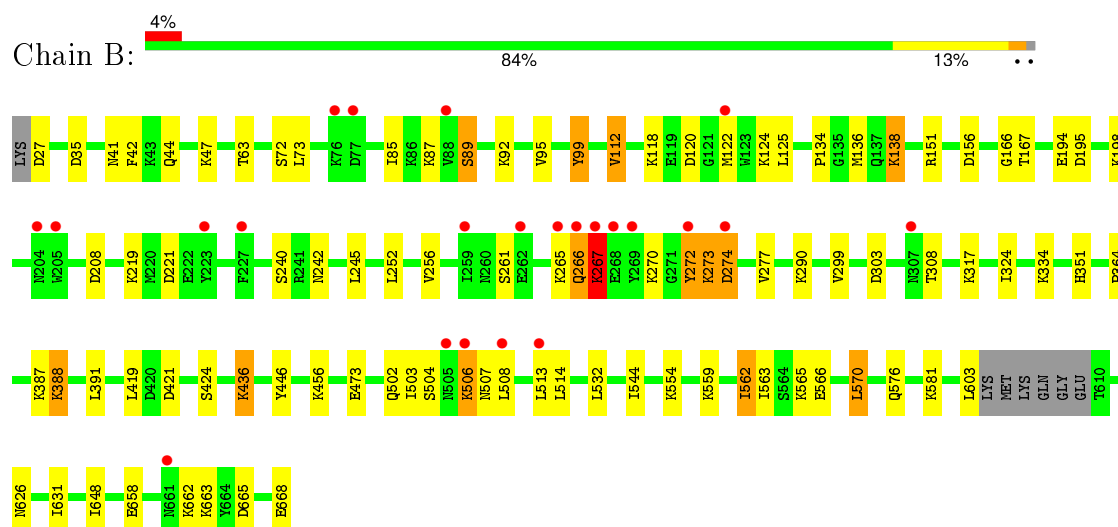
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: PENICILLIN BINDING PROTEIN 2 PRIME



• Molecule 1: PENICILLIN BINDING PROTEIN 2 PRIME



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	81.44Å 101.64Å 186.84Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.64 – 2.35 46.71 – 2.35	Depositor EDS
% Data completeness (in resolution range)	100.0 (46.64-2.35) 99.7 (46.71-2.35)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.15 (at 2.34Å)	Xtriage
Refinement program	BUSTER 2.10.0	Depositor
R, R_{free}	0.182 , 0.231 0.201 , 0.247	Depositor DCC
R_{free} test set	3319 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	32.7	Xtriage
Anisotropy	0.819	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 55.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 65497 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11030	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.70% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: MUR, CD, CL

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	0.49	0/5198	0.66	0/6987
1	B	0.49	0/5189	0.67	0/6976
All	All	0.49	0/10387	0.67	0/13963

There are no bond length outliers.

There are no bond angle outliers.

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	5113	0	5124	45	0
1	B	5104	0	5111	41	0
2	A	7	0	0	0	0
2	B	4	0	0	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	17	0	16	9	0
4	B	17	0	16	15	0
5	A	406	0	0	1	0
5	B	358	0	0	0	0
All	All	11030	0	10267	89	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 4.

All (89) 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:B:277:VAL:CG1	4:B:1677:MUR:H91	1.67	1.25
1:B:277:VAL:HG11	4:B:1677:MUR:H91	1.32	1.11
1:B:277:VAL:HG11	4:B:1677:MUR:C9	1.91	0.98
1:B:256:VAL:HG21	4:B:1677:MUR:H92	1.55	0.86
1:B:277:VAL:HG12	4:B:1677:MUR:H91	1.59	0.84
1:A:277:VAL:HB	4:A:1679:MUR:H91	1.65	0.78
1:A:220:MET:HA	1:A:224:LEU:HD23	1.67	0.77
1:B:277:VAL:CG1	4:B:1677:MUR:C9	2.54	0.75
1:B:138:LYS:H	1:B:138:LYS:HD2	1.51	0.74
1:A:277:VAL:HB	4:A:1679:MUR:C9	2.17	0.74
1:A:138:LYS:H	1:A:138:LYS:HD2	1.53	0.73
1:B:256:VAL:CG2	4:B:1677:MUR:H92	2.18	0.72
1:A:293:HIS:NE2	4:A:1679:MUR:O9	2.24	0.70
4:B:1677:MUR:H93	4:B:1677:MUR:N2	2.11	0.66
1:A:28:LYS:HG2	1:A:29:GLU:H	1.61	0.65
1:B:89:SER:HB3	1:B:92:LYS:HB2	1.80	0.64
1:A:559:LYS:HB3	1:A:562:ILE:HD13	1.80	0.63
4:B:1677:MUR:H3	4:B:1677:MUR:O8	1.98	0.63
1:B:559:LYS:HB3	1:B:562:ILE:HD13	1.81	0.62
1:A:305:ASN:HB3	1:B:72:SER:HB3	1.84	0.59
1:B:112:VAL:HG13	1:B:134:PRO:HB3	1.87	0.57
1:A:112:VAL:HG13	1:A:134:PRO:HB3	1.88	0.56
1:B:138:LYS:H	1:B:138:LYS:CD	2.19	0.56
1:A:165:THR:OG1	4:A:1679:MUR:O6	2.14	0.56
1:A:364:PRO:HG2	1:A:388:LYS:HB3	1.89	0.55
1:B:277:VAL:CB	4:B:1677:MUR:H91	2.35	0.54
1:A:73:LEU:HD21	1:A:299:VAL:HG11	1.90	0.54
1:A:138:LYS:H	1:A:138:LYS:CD	2.20	0.53
1:A:86:LYS:HE3	1:A:94:ARG:HD3	1.91	0.52
1:A:502:GLN:HE22	1:A:505:ASN:ND2	2.08	0.52
1:B:256:VAL:HG21	4:B:1677:MUR:C9	2.32	0.52
4:B:1677:MUR:HN22	4:B:1677:MUR:C9	2.24	0.51
4:B:1677:MUR:C9	4:B:1677:MUR:N2	2.72	0.51
1:B:151:ARG:HD2	4:B:1677:MUR:H7	1.92	0.51
1:A:151:ARG:HB2	4:A:1679:MUR:O9	2.11	0.51
1:A:166:GLY:HA3	1:A:242:ASN:HB2	1.94	0.50
1:B:166:GLY:HA3	1:B:242:ASN:HB2	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:504:SER:O	1:A:505:ASN:HB2	2.11	0.50
1:A:517:SER:HB3	1:A:524:ILE:HD11	1.93	0.50
1:A:508:LEU:HD23	1:A:513:LEU:HD23	1.94	0.49
1:B:73:LEU:HD21	1:B:299:VAL:HG11	1.94	0.49
1:A:277:VAL:HB	4:A:1679:MUR:C7	2.42	0.49
1:A:42:PHE:HB3	1:A:63:THR:HA	1.96	0.48
1:B:194:GLU:O	1:B:198:LYS:HD3	2.13	0.48
1:B:42:PHE:HB3	1:B:63:THR:HA	1.97	0.47
1:B:290:LYS:HB2	1:B:324:ILE:HD11	1.96	0.47
1:B:272:TYR:O	1:B:273:LYS:HB2	2.14	0.47
1:B:508:LEU:HD23	1:B:513:LEU:HD23	1.97	0.47
1:B:506:LYS:HD2	1:B:507:ASN:HB2	1.97	0.47
1:B:544:ILE:HB	1:B:559:LYS:HB2	1.97	0.46
1:B:240:SER:HA	4:B:1677:MUR:O1	2.15	0.46
1:B:436:LYS:HE3	1:B:436:LYS:H	1.81	0.46
1:A:99:TYR:HB2	1:A:112:VAL:HG22	1.98	0.45
1:B:41:ASN:ND2	1:B:44:GLN:HB2	2.32	0.45
1:B:99:TYR:HD2	1:B:112:VAL:HG11	1.82	0.45
1:A:277:VAL:CB	4:A:1679:MUR:C9	2.92	0.45
1:B:99:TYR:HB2	1:B:112:VAL:HG22	1.99	0.44
1:A:403:SER:CB	1:A:599:GLY:HA3	2.47	0.44
1:A:99:TYR:HD2	1:A:112:VAL:HG11	1.82	0.44
1:A:516:ASP:HA	1:A:519:TYR:CE1	2.53	0.44
1:A:665:ASP:HB3	1:A:668:GLU:HB2	2.00	0.44
1:A:277:VAL:O	4:A:1679:MUR:O8	2.36	0.43
1:A:503:ILE:O	1:A:524:ILE:HG12	2.17	0.43
1:B:658:GLU:HB2	1:B:662:LYS:HG2	2.00	0.43
1:A:486:LEU:HD23	1:A:562:ILE:HG13	2.00	0.43
1:B:665:ASP:HB3	1:B:668:GLU:HB2	1.99	0.43
1:A:290:LYS:HB2	1:A:324:ILE:HD11	2.00	0.43
1:A:561:ASN:ND2	5:A:2330:HOH:O	2.48	0.43
1:B:562:ILE:HG22	1:B:563:ILE:HG22	2.01	0.43
1:A:77:ASP:HB2	1:A:102:LYS:HD3	2.00	0.43
1:B:421:ASP:HB3	1:B:570:LEU:HD11	2.01	0.43
1:A:544:ILE:HB	1:A:559:LYS:HB2	2.00	0.43
1:B:85:ILE:HG12	1:B:95:VAL:HG22	2.00	0.43
1:A:351:HIS:HD2	1:A:626:ASN:O	2.02	0.43
1:A:41:ASN:ND2	1:A:44:GLN:HB2	2.33	0.42
1:A:562:ILE:HG22	1:A:563:ILE:HG22	2.02	0.42
1:A:421:ASP:HB3	1:A:570:LEU:HD11	2.00	0.42
1:A:245:LEU:HD13	1:A:334:LYS:HG3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:364:PRO:HG2	1:B:388:LYS:HB3	2.02	0.42
1:B:245:LEU:HD13	1:B:334:LYS:HG3	2.01	0.42
1:B:99:TYR:HB2	1:B:112:VAL:CG2	2.50	0.41
1:A:631:ILE:HD13	1:A:648:ILE:HG22	2.02	0.41
1:A:89:SER:HB3	1:A:92:LYS:HB2	2.02	0.41
1:B:631:ILE:HD13	1:B:648:ILE:HG22	2.03	0.41
1:A:277:VAL:CG1	4:A:1679:MUR:H7	2.51	0.41
1:B:266:GLN:HB3	1:B:267:LYS:H	1.65	0.41
1:A:105:TYR:O	1:A:314:ILE:HG21	2.22	0.40
1:A:408:LEU:HD22	1:A:534:ILE:HG21	2.02	0.40
1:B:351:HIS:HD2	1:B:626:ASN:O	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	633/643 (98%)	612 (97%)	20 (3%)	1 (0%)	52	63
1	B	632/643 (98%)	607 (96%)	19 (3%)	6 (1%)	21	22
All	All	1265/1286 (98%)	1219 (96%)	39 (3%)	7 (1%)	30	34

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	265	LYS
1	B	273	LYS
1	B	504	SER
1	B	267	LYS
1	B	506	LYS
1	A	28	LYS
1	B	274	ASP

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	568/573 (99%)	529 (93%)	39 (7%)	19	21
1	B	567/573 (99%)	515 (91%)	52 (9%)	11	11
All	All	1135/1146 (99%)	1044 (92%)	91 (8%)	15	16

All (91) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	ASP
1	A	28	LYS
1	A	87	LYS
1	A	89	SER
1	A	99	TYR
1	A	112	VAL
1	A	118	LYS
1	A	124	LYS
1	A	125	LEU
1	A	136	MET
1	A	138	LYS
1	A	141	SER
1	A	156	ASP
1	A	167	THR
1	A	195	ASP
1	A	208	ASP
1	A	252	LEU
1	A	274	ASP
1	A	303	ASP
1	A	308	THR
1	A	317	LYS
1	A	387	LYS
1	A	391	LEU
1	A	419	LEU
1	A	424	SER
1	A	456	LYS
1	A	473	GLU

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Mol	Chain	Res	Type
1	A	491	ASP
1	A	513	LEU
1	A	514	LEU
1	A	532	LEU
1	A	554	LYS
1	A	562	ILE
1	A	565	LYS
1	A	566	GLU
1	A	570	LEU
1	A	573	ASP
1	A	576	GLN
1	A	663	LYS
1	B	27	ASP
1	B	35	ASP
1	B	47	LYS
1	B	87	LYS
1	B	89	SER
1	B	99	TYR
1	B	112	VAL
1	B	118	LYS
1	B	120	ASP
1	B	122	MET
1	B	124	LYS
1	B	125	LEU
1	B	136	MET
1	B	138	LYS
1	B	156	ASP
1	B	167	THR
1	B	195	ASP
1	B	208	ASP
1	B	219	LYS
1	B	221	ASP
1	B	252	LEU
1	B	261	SER
1	B	266	GLN
1	B	267	LYS
1	B	270	LYS
1	B	272	TYR
1	B	274	ASP
1	B	303	ASP
1	B	308	THR
1	B	317	LYS

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Mol	Chain	Res	Type
1	B	387	LYS
1	B	388	LYS
1	B	391	LEU
1	B	419	LEU
1	B	424	SER
1	B	436	LYS
1	B	446	TYR
1	B	456	LYS
1	B	473	GLU
1	B	502	GLN
1	B	503	ILE
1	B	514	LEU
1	B	532	LEU
1	B	554	LYS
1	B	562	ILE
1	B	565	LYS
1	B	566	GLU
1	B	570	LEU
1	B	576	GLN
1	B	581	LYS
1	B	603	LEU
1	B	663	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	ASN
1	A	57	ASN
1	A	505	ASN
1	A	545	ASN
1	A	576	GLN
1	A	624	ASN
1	B	32	ASN
1	B	177	ASN
1	B	521	GLN
1	B	545	ASN
1	B	576	GLN
1	B	624	ASN

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 ⓘ

Of 17 ligands modelled in this entry, 15 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	MUR	A	1679	-	14,17,17	1.47	3 (21%)	17,24,24	3.95	11 (64%)
4	MUR	B	1677	-	14,17,17	1.08	2 (14%)	17,24,24	3.61	8 (47%)

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
4	MUR	A	1679	-	-	0/6/30/30	0/1/1/1
4	MUR	B	1677	-	-	0/6/30/30	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1679	MUR	O5-C1	-2.66	1.38	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1679	MUR	O5-C5	-2.39	1.38	1.44
4	A	1679	MUR	C4-C5	-2.32	1.48	1.53
4	B	1677	MUR	O3-C3	-2.17	1.38	1.43
4	B	1677	MUR	C4-C5	-2.16	1.48	1.53

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1679	MUR	C9-C7-C8	-10.24	100.07	113.18
4	B	1677	MUR	C9-C7-C8	-7.41	103.70	113.18
4	A	1679	MUR	O5-C1-C2	-6.93	101.47	109.62
4	B	1677	MUR	O5-C1-C2	-6.26	102.26	109.62
4	B	1677	MUR	C6-C5-C4	-5.50	99.46	113.02
4	A	1679	MUR	O4-C4-C3	-3.55	101.48	109.87
4	A	1679	MUR	O6-C6-C5	-3.21	100.73	111.33
4	A	1679	MUR	O4-C4-C5	-2.45	102.74	109.24
4	A	1679	MUR	C1-C2-N2	-2.07	107.04	111.10
4	A	1679	MUR	O1-C1-C2	2.76	115.33	109.02
4	A	1679	MUR	C1-O5-C5	2.84	118.73	113.47
4	B	1677	MUR	O5-C5-C4	3.56	116.36	109.68
4	B	1677	MUR	C1-C2-N2	3.64	118.25	111.10
4	A	1679	MUR	O5-C5-C4	3.68	116.60	109.68
4	B	1677	MUR	C3-C2-N2	4.42	118.84	110.73
4	B	1677	MUR	C1-O5-C5	4.56	121.90	113.47
4	A	1679	MUR	O3-C7-C9	4.66	118.33	107.17
4	B	1677	MUR	C4-C3-C2	4.74	118.94	111.26
4	A	1679	MUR	C4-C3-C2	4.88	119.16	111.26

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 24 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1679	MUR	9	0
4	B	1677	MUR	15	0

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	637/643 (99%)	-0.12	16 (2%) 61 73	19, 43, 77, 106	0
1	B	636/643 (98%)	0.01	23 (3%) 46 60	25, 43, 87, 128	0
All	All	1273/1286 (98%)	-0.05	39 (3%) 52 65	19, 43, 82, 128	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	266	GLN	5.9
1	B	505	ASN	5.7
1	B	506	LYS	5.3
1	B	267	LYS	5.2
1	A	223	TYR	3.8
1	A	508	LEU	3.5
1	B	122	MET	3.5
1	B	265	LYS	3.4
1	A	224	LEU	3.3
1	A	220	MET	2.9
1	A	509	ASP	2.9
1	A	226	ASP	2.7
1	B	269	TYR	2.7
1	B	508	LEU	2.6
1	A	229	LYS	2.6
1	A	504	SER	2.6
1	B	268	GLU	2.6
1	A	272	TYR	2.5
1	A	204	ASN	2.5
1	B	227	PHE	2.5
1	A	610	THR	2.5
1	A	222	GLU	2.4
1	B	204	ASN	2.4
1	B	88	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	305	ASN	2.3
1	B	205	TRP	2.3
1	B	77	ASP	2.2
1	B	259	ILE	2.2
1	B	262	GLU	2.2
1	A	505	ASN	2.2
1	B	223	TYR	2.2
1	B	272	TYR	2.1
1	A	85	ILE	2.1
1	B	661	ASN	2.1
1	B	274	ASP	2.1
1	B	307	ASN	2.0
1	A	441	TYR	2.0
1	B	76	LYS	2.0
1	B	513	LEU	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 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(Å ²)	Q<0.9
4	MUR	A	1679	17/17	0.80	0.28	11.20	20,33,61,63	0
4	MUR	B	1677	17/17	0.83	0.27	10.22	22,43,64,72	0
2	CD	B	1669	1/1	0.94	0.12	0.13	35,35,35,35	0
2	CD	B	1671	1/1	0.99	0.13	-0.42	33,33,33,33	0
3	CL	B	1675	1/1	0.96	0.11	-0.82	33,33,33,33	0
2	CD	A	1674	1/1	0.82	0.10	-0.89	160,160,160,160	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	CL	B	1674	1/1	0.96	0.12	-1.06	36,36,36,36	0
2	CD	A	1670	1/1	0.97	0.11	-1.20	34,34,34,34	0
2	CD	A	1672	1/1	0.93	0.09	-1.48	39,39,39,39	0
3	CL	A	1678	1/1	0.95	0.08	-1.67	35,35,35,35	0
3	CL	A	1677	1/1	0.98	0.09	-1.99	35,35,35,35	0
2	CD	A	1675	1/1	0.84	0.07	-	123,123,123,123	0
2	CD	A	1676	1/1	0.97	0.09	-	122,122,122,122	0
2	CD	B	1673	1/1	0.99	0.09	-	47,47,47,47	0
2	CD	A	1671	1/1	0.96	0.14	-	84,84,84,84	0
2	CD	B	1672	1/1	0.91	0.08	-	86,86,86,86	0
2	CD	A	1673	1/1	0.96	0.05	-	123,123,123,123	0

6.5 Other polymers [i](#)

There are no such residues in this entry.