



Full wwPDB X-ray Structure Validation Report ⓘ

May 28, 2020 – 09:18 pm BST

PDB ID : 2C6W
Title : PENICILLIN-BINDING PROTEIN 1A (PBP-1A) FROM STREPTOCOCCUS PNEUMONIAE
Authors : Contreras-Martel, C.; Job, V.; Di Guilmi, A.-M.; Vernet, T.; Dideberg, O.; Dessen, A.
Deposited on : 2005-11-14
Resolution : 2.61 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

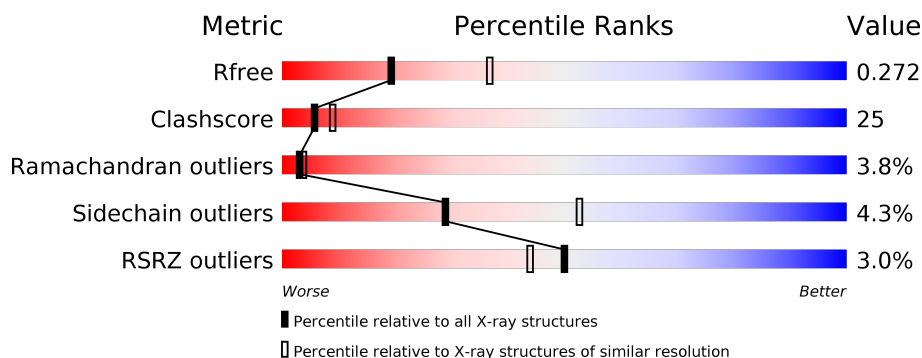
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.61 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}	130704	3797 (2.64-2.60)
Clashscore	141614	4168 (2.64-2.60)
Ramachandran outliers	138981	4093 (2.64-2.60)
Sidechain outliers	138945	4093 (2.64-2.60)
RSRZ outliers	127900	3731 (2.64-2.60)

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 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	16	<div> <div></div> <div>38%</div> <div>56%</div> <div>6%</div> </div>
2	B	384	<div> <div>3%</div> <div>59%</div> <div>37%</div> <div>.</div> </div>

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 3167 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 1A.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	16	Total	C	N	O	0	0	0
			119	73	19	27			

- Molecule 2 is a protein called PENICILLIN-BINDING PROTEIN 1A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	384	Total	C	N	O	S	0	0	0
			3006	1904	494	595	13			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	388	ASP	GLU	conflict	UNP Q8DR59
B	540	THR	SER	conflict	UNP Q8DR59

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	3	Total	Zn	0	0
			3	3		

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	4	Total	Cl	0	0
			4	4		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	35	Total	O	0	0
			35	35		

3 Residue-property plots [i](#)

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.

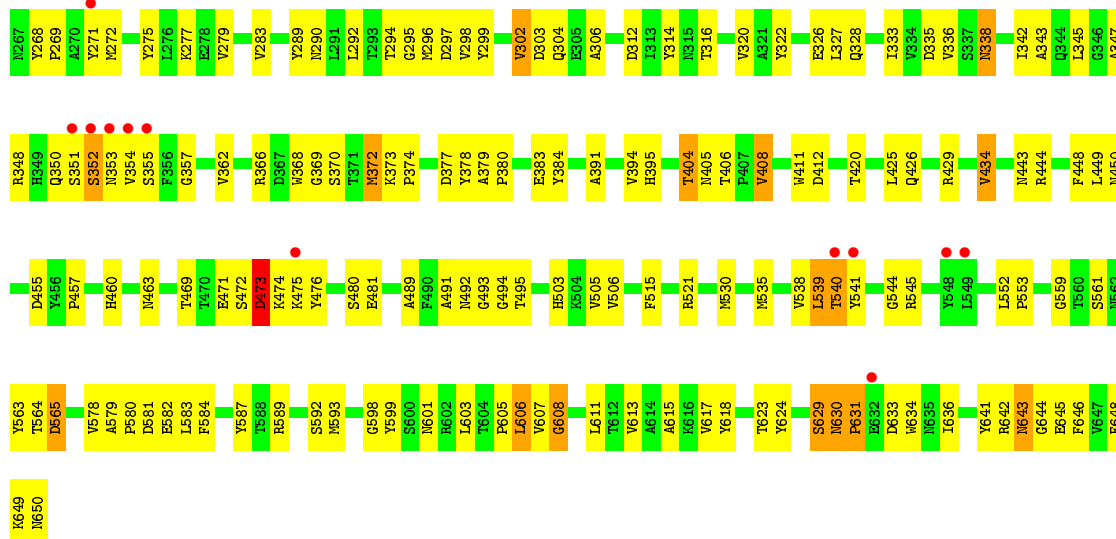
• Molecule 1: PENICILLIN-BINDING PROTEIN 1A

Chain A: 



• Molecule 2: PENICILLIN-BINDING PROTEIN 1A

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	109.78Å 187.26Å 51.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.99 – 2.61 44.99 – 2.61	Depositor EDS
% Data completeness (in resolution range)	96.0 (44.99-2.61) 96.1 (44.99-2.61)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.70 (at 2.61Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.240 , 0.276 0.234 , 0.272	Depositor DCC
R_{free} test set	779 reflections (4.89%)	wwPDB-VP
Wilson B-factor (Å ²)	62.5	Xtriage
Anisotropy	0.777	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 54.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.026 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.027 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3167	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.56% 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: ZN, 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	0.34	0/119	0.51	0/159
2	B	0.40	0/3080	0.68	1/4196 (0.0%)
All	All	0.39	0/3199	0.67	1/4355 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	408	VAL	N-CA-C	-5.02	97.44	111.00

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	119	0	100	14	0
2	B	3006	0	2833	146	1
3	B	3	0	0	0	0
4	B	4	0	0	1	0
5	B	35	0	0	1	0
All	All	3167	0	2933	154	1

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 25.

All (154) 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)
2:B:370:SER:HB2	2:B:559:GLY:HA2	1.47	0.96
2:B:474:LYS:CB	2:B:474:LYS:CD	2.51	0.87
2:B:275:TYR:CZ	2:B:302:VAL:HG22	2.15	0.82
2:B:404:THR:CG2	2:B:406:THR:H	1.95	0.80
2:B:297:ASP:HB2	2:B:506:VAL:HG22	1.63	0.80
2:B:290:ASN:HD21	2:B:292:LEU:HB2	1.46	0.80
2:B:404:THR:HG23	2:B:406:THR:H	1.53	0.74
2:B:384:TYR:CD1	2:B:444:ARG:HD2	2.26	0.70
2:B:443:ASN:HB2	4:B:705:CL:CL	2.29	0.70
2:B:563:TYR:HE1	2:B:578:VAL:HG12	1.56	0.70
2:B:335:ASP:HB3	2:B:338:ASN:HD21	1.57	0.69
2:B:290:ASN:ND2	2:B:292:LEU:HB2	2.06	0.69
2:B:336:VAL:HG12	2:B:492:ASN:HB3	1.78	0.66
2:B:578:VAL:HG13	2:B:605:PRO:CB	2.26	0.65
2:B:297:ASP:HB2	2:B:506:VAL:CG2	2.27	0.65
2:B:373:LYS:HB2	2:B:374:PRO:HD3	1.78	0.65
2:B:327:LEU:HD23	2:B:327:LEU:C	2.17	0.64
2:B:455:ASP:O	2:B:476:TYR:HA	1.98	0.64
2:B:578:VAL:HG13	2:B:605:PRO:HB2	1.80	0.64
2:B:642:ARG:HG3	2:B:643:ASN:N	2.12	0.64
2:B:379:ALA:HB3	2:B:380:PRO:HD3	1.78	0.63
2:B:457:PRO:HG3	2:B:471:GLU:HB2	1.81	0.63
2:B:271:TYR:CE1	2:B:272:MET:HG2	2.34	0.63
2:B:489:ALA:HA	2:B:492:ASN:HD21	1.62	0.62
2:B:354:VAL:O	2:B:354:VAL:HG22	1.99	0.62
2:B:489:ALA:HA	2:B:492:ASN:ND2	2.14	0.61
2:B:391:ALA:O	2:B:420:THR:HG22	2.00	0.61
2:B:368:TRP:CD2	2:B:583:LEU:HD22	2.36	0.61
2:B:366:ARG:HD2	2:B:581:ASP:OD1	2.00	0.61
2:B:601:ASN:HD21	2:B:603:LEU:HB2	1.66	0.60
2:B:298:VAL:HG13	2:B:505:VAL:HG12	1.83	0.60
2:B:642:ARG:HG3	2:B:643:ASN:H	1.66	0.60
1:A:52:SER:O	2:B:295:GLY:HA2	2.01	0.59
2:B:275:TYR:CE2	2:B:302:VAL:HG22	2.38	0.59
2:B:362:VAL:HG12	2:B:481:GLU:OE2	2.02	0.59
2:B:564:THR:O	2:B:565:ASP:HB2	2.00	0.59
2:B:493:GLY:O	2:B:521:ARG:NH2	2.35	0.59
2:B:333:ILE:HG13	2:B:593:MET:HG3	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:589:ARG:HH11	2:B:589:ARG:HG2	1.67	0.58
2:B:491:ALA:HB2	2:B:587:TYR:CD1	2.39	0.58
2:B:312:ASP:O	2:B:316:THR:HG23	2.04	0.57
2:B:335:ASP:HB3	2:B:338:ASN:ND2	2.19	0.57
2:B:553:PRO:HD2	2:B:631:PRO:HB3	1.86	0.57
2:B:378:TYR:CD1	2:B:425:LEU:HD13	2.39	0.56
2:B:275:TYR:CE1	2:B:302:VAL:HG22	2.40	0.56
2:B:384:TYR:HB3	2:B:444:ARG:NH1	2.20	0.56
2:B:578:VAL:CG1	2:B:605:PRO:HB3	2.36	0.55
2:B:539:LEU:HD11	2:B:584:PHE:HZ	1.72	0.55
2:B:539:LEU:HD12	2:B:544:GLY:O	2.07	0.55
2:B:404:THR:HG22	2:B:406:THR:H	1.72	0.55
2:B:335:ASP:CB	2:B:338:ASN:HD21	2.18	0.54
2:B:336:VAL:HG22	2:B:592:SER:HB2	1.90	0.54
2:B:314:TYR:HB3	2:B:348:ARG:HB2	1.89	0.54
2:B:408:VAL:HG11	2:B:434:VAL:HG11	1.90	0.54
2:B:368:TRP:CE3	2:B:583:LEU:HD22	2.44	0.53
2:B:613:VAL:O	2:B:617:VAL:HG23	2.08	0.53
2:B:314:TYR:HB3	2:B:348:ARG:CB	2.39	0.53
2:B:475:LYS:HE3	2:B:476:TYR:CZ	2.44	0.53
2:B:369:GLY:O	2:B:372:MET:HB2	2.10	0.52
2:B:563:TYR:HE1	2:B:578:VAL:CG1	2.20	0.52
2:B:598:GLY:HA2	2:B:606:LEU:HD21	1.91	0.52
2:B:299:TYR:O	2:B:503:HIS:HB3	2.09	0.52
2:B:268:TYR:CD1	2:B:268:TYR:C	2.83	0.52
2:B:271:TYR:CD1	2:B:272:MET:HG2	2.44	0.52
2:B:471:GLU:HB3	2:B:476:TYR:HD1	1.75	0.52
2:B:643:ASN:HB3	2:B:648:PHE:CE1	2.45	0.51
2:B:505:VAL:HG22	2:B:515:PHE:HE1	1.76	0.51
2:B:623:THR:HG22	2:B:624:TYR:N	2.26	0.51
2:B:563:TYR:CE2	2:B:580:PRO:HB3	2.46	0.51
2:B:336:VAL:HG11	2:B:491:ALA:HB3	1.93	0.50
2:B:404:THR:HG23	2:B:405:ASN:N	2.26	0.50
2:B:338:ASN:ND2	2:B:338:ASN:N	2.60	0.49
2:B:564:THR:OG1	2:B:565:ASP:N	2.44	0.49
2:B:450:ASN:N	2:B:450:ASN:HD22	2.10	0.49
2:B:589:ARG:HG2	2:B:589:ARG:NH1	2.27	0.49
2:B:615:ALA:O	2:B:618:TYR:HB3	2.13	0.48
2:B:335:ASP:CG	2:B:338:ASN:HD21	2.17	0.48
2:B:351:SER:O	2:B:352:SER:HB3	2.13	0.48
2:B:582:GLU:C	2:B:583:LEU:HG	2.34	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:54:LYS:C	1:A:55:ILE:HD12	2.34	0.48
2:B:336:VAL:HG11	2:B:491:ALA:CB	2.42	0.48
2:B:327:LEU:HD23	2:B:328:GLN:N	2.29	0.48
1:A:53:SER:HA	2:B:296:MET:O	2.14	0.47
2:B:641:TYR:CZ	2:B:648:PHE:HB2	2.49	0.47
2:B:289:TYR:N	2:B:289:TYR:CD1	2.82	0.47
2:B:298:VAL:HG13	2:B:505:VAL:CG1	2.44	0.47
2:B:629:SER:O	2:B:630:ASN:HB3	2.15	0.47
2:B:327:LEU:HD23	2:B:328:GLN:O	2.15	0.47
2:B:505:VAL:O	2:B:505:VAL:HG23	2.15	0.47
2:B:368:TRP:O	2:B:583:LEU:HD13	2.14	0.47
2:B:599:TYR:CE1	2:B:606:LEU:HD22	2.50	0.47
2:B:327:LEU:C	2:B:327:LEU:CD2	2.83	0.46
2:B:469:THR:HG23	2:B:471:GLU:O	2.15	0.46
2:B:578:VAL:HG13	2:B:605:PRO:HB3	1.94	0.46
2:B:299:TYR:HE1	2:B:506:VAL:HG13	1.81	0.46
2:B:535:MET:HA	2:B:538:VAL:HG23	1.97	0.46
2:B:636:ILE:O	2:B:636:ILE:HG23	2.15	0.46
2:B:473:ASP:HB3	2:B:474:LYS:H	1.44	0.46
2:B:275:TYR:O	2:B:279:VAL:HG23	2.16	0.45
2:B:411:TRP:CZ3	2:B:412:ASP:HB3	2.51	0.45
2:B:303:ASP:O	2:B:306:ALA:HB3	2.16	0.45
2:B:277:LYS:HE2	2:B:357:GLY:O	2.16	0.45
2:B:314:TYR:CD2	2:B:345:LEU:HD23	2.52	0.45
2:B:383:GLU:HG3	2:B:384:TYR:CD2	2.52	0.45
2:B:489:ALA:O	2:B:494:GLY:N	2.49	0.45
2:B:378:TYR:CE1	2:B:425:LEU:HD13	2.52	0.45
2:B:492:ASN:OD1	2:B:495:THR:HB	2.16	0.45
2:B:563:TYR:CE1	2:B:578:VAL:HG12	2.45	0.45
2:B:314:TYR:HA	2:B:322:TYR:OH	2.17	0.44
1:A:54:LYS:HB3	1:A:62:LEU:CD1	2.47	0.44
2:B:268:TYR:HD1	2:B:269:PRO:O	2.01	0.44
1:A:63:ILE:HD11	2:B:271:TYR:OH	2.18	0.44
2:B:368:TRP:CH2	2:B:480:SER:HA	2.53	0.44
2:B:342:ILE:O	2:B:343:ALA:HB2	2.18	0.44
2:B:530:MET:CE	2:B:649:LYS:HE2	2.48	0.44
1:A:59:LYS:O	1:A:61:GLN:N	2.51	0.43
1:A:57:ASP:OD2	2:B:304:GLN:NE2	2.51	0.43
2:B:404:THR:HG23	2:B:406:THR:N	2.25	0.43
2:B:644:GLY:O	2:B:646:PHE:N	2.51	0.43
2:B:383:GLU:HG3	2:B:384:TYR:CE2	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:457:PRO:HG3	2:B:471:GLU:CB	2.48	0.43
2:B:642:ARG:CG	2:B:643:ASN:N	2.82	0.43
2:B:394:VAL:HG22	2:B:395:HIS:H	1.84	0.43
2:B:338:ASN:N	2:B:338:ASN:HD22	2.16	0.43
2:B:641:TYR:HD2	2:B:650:ASN:ND2	2.17	0.43
2:B:391:ALA:HB2	2:B:646:PHE:CG	2.54	0.43
2:B:607:VAL:HG12	2:B:608:GLY:H	1.84	0.43
2:B:633:ASP:OD1	2:B:634:TRP:N	2.41	0.42
1:A:54:LYS:HB3	1:A:62:LEU:HD12	2.01	0.42
2:B:380:PRO:HG3	2:B:448:PHE:CD1	2.54	0.42
1:A:63:ILE:HD13	2:B:304:GLN:CD	2.40	0.42
2:B:540:THR:O	2:B:545:ARG:HG2	2.20	0.42
1:A:55:ILE:HD12	1:A:55:ILE:N	2.35	0.42
1:A:61:GLN:O	1:A:62:LEU:C	2.57	0.42
2:B:347:ALA:HB1	2:B:350:GLN:HE21	1.85	0.42
2:B:578:VAL:CG1	2:B:579:ALA:N	2.83	0.42
2:B:643:ASN:ND2	5:B:2034:HOH:O	2.53	0.42
1:A:57:ASP:OD2	1:A:61:GLN:HG3	2.18	0.42
2:B:404:THR:CG2	2:B:405:ASN:N	2.83	0.42
1:A:63:ILE:HD13	2:B:304:GLN:OE1	2.20	0.41
2:B:275:TYR:C	2:B:275:TYR:CD2	2.94	0.41
2:B:378:TYR:HD1	2:B:425:LEU:HD13	1.82	0.41
2:B:449:LEU:HD23	2:B:449:LEU:HA	1.88	0.41
2:B:530:MET:HE1	2:B:649:LYS:HE2	2.02	0.41
2:B:326:GLU:O	2:B:327:LEU:C	2.59	0.41
2:B:561:SER:OG	2:B:581:ASP:HB2	2.20	0.41
1:A:59:LYS:C	1:A:61:GLN:H	2.24	0.41
2:B:460:HIS:O	2:B:463:ASN:HB2	2.20	0.41
2:B:450:ASN:N	2:B:450:ASN:ND2	2.68	0.41
2:B:582:GLU:OE1	2:B:611:LEU:HD22	2.21	0.41
2:B:601:ASN:ND2	2:B:603:LEU:HB2	2.33	0.41
2:B:552:LEU:HA	2:B:553:PRO:HD2	1.99	0.40
2:B:606:LEU:HD23	2:B:606:LEU:N	2.36	0.40
2:B:279:VAL:O	2:B:283:VAL:HG23	2.21	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:472:SER:OG	2:B:472:SER:OG[3_756]	2.17	0.03

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	14/16 (88%)	10 (71%)	2 (14%)	2 (14%)	0	0
2	B	382/384 (100%)	325 (85%)	44 (12%)	13 (3%)	3	5
All	All	396/400 (99%)	335 (85%)	46 (12%)	15 (4%)	3	4

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	60	ASN
2	B	352	SER
2	B	355	SER
2	B	473	ASP
2	B	630	ASN
2	B	353	ASN
2	B	540	THR
2	B	565	ASP
2	B	631	PRO
2	B	643	ASN
2	B	645	GLU
2	B	294	THR
2	B	629	SER
1	A	62	LEU
2	B	608	GLY

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	11/15 (73%)	10 (91%)	1 (9%)	9	17
2	B	315/324 (97%)	302 (96%)	13 (4%)	30	55
All	All	326/339 (96%)	312 (96%)	14 (4%)	29	53

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	62	LEU
2	B	302	VAL
2	B	320	VAL
2	B	338	ASN
2	B	372	MET
2	B	377	ASP
2	B	404	THR
2	B	426	GLN
2	B	429	ARG
2	B	434	VAL
2	B	473	ASP
2	B	539	LEU
2	B	541	TYR
2	B	606	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	290	ASN
2	B	304	GLN
2	B	338	ASN
2	B	350	GLN
2	B	426	GLN
2	B	427	GLN
2	B	438	ASN
2	B	443	ASN
2	B	450	ASN
2	B	503	HIS
2	B	571	HIS
2	B	601	ASN
2	B	650	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 7 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 [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	16/16 (100%)	0.17	0	100	100	80, 92, 103, 109	0
2	B	384/384 (100%)	0.09	12 (3%)	49	42	40, 63, 97, 121	5 (1%)
All	All	400/400 (100%)	0.10	12 (3%)	50	44	40, 64, 98, 121	5 (1%)

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	541	TYR	3.8
2	B	353	ASN	2.9
2	B	549	LEU	2.9
2	B	352	SER	2.9
2	B	354	VAL	2.9
2	B	355	SER	2.7
2	B	540	THR	2.4
2	B	475	LYS	2.2
2	B	632	GLU	2.2
2	B	548	TYR	2.2
2	B	271	TYR	2.1
2	B	351	SER	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. 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
4	CL	B	706	1/1	0.76	0.31	105,105,105,105	0
4	CL	B	705	1/1	0.87	0.14	99,99,99,99	0
4	CL	B	707	1/1	0.89	0.46	100,100,100,100	0
3	ZN	B	701	1/1	0.92	0.19	100,100,100,100	0
4	CL	B	704	1/1	0.93	0.20	73,73,73,73	0
3	ZN	B	703	1/1	0.99	0.16	61,61,61,61	0
3	ZN	B	702	1/1	0.99	0.22	51,51,51,51	0

6.5 Other polymers [i](#)

There are no such residues in this entry.