



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:35 AM UTC

PDB ID : 5N4B / pdb_00005n4b
Title : Prolyl oligopeptidase B from *Galerina marginata* bound to 25mer macrocyclization substrate - S577A mutant
Authors : Czekster, C.M.; McMahon, S.A.; Ludewig, H.; Naismith, J.H.
Deposited on : 2017-02-10
Resolution : 1.44 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

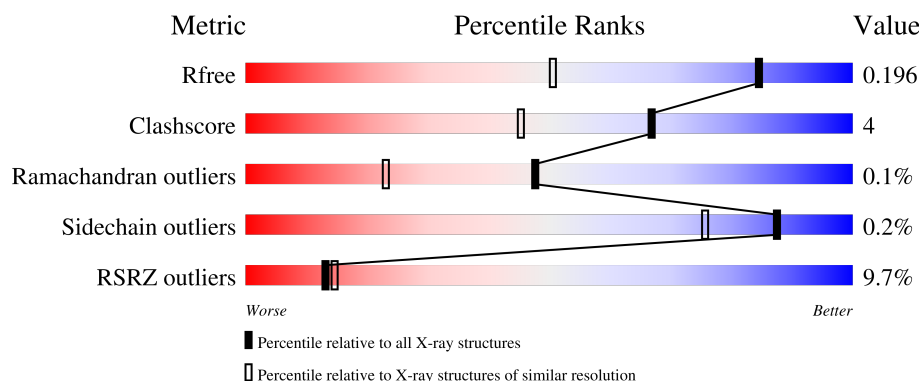
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.44 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	3234 (1.46-1.42)
Clashscore	190562	3289 (1.46-1.42)
Ramachandran outliers	187476	3248 (1.46-1.42)
Sidechain outliers	187428	3248 (1.46-1.42)
RSRZ outliers	180081	3234 (1.46-1.42)

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	731	8% 92% 6%
1	B	731	9% 92% 6%
2	C	25	28% 56% 12% 32%
2	D	25	24% 56% 12% 32%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13794 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 Prolyl oligopeptidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	722	Total	C	N	O	S	0	13	0
			5796	3717	971	1096	12			
1	B	722	Total	C	N	O	S	0	13	0
			5797	3719	971	1096	11			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP H2E7Q8
A	577	ALA	SER	engineered mutation	UNP H2E7Q8
B	0	GLY	-	expression tag	UNP H2E7Q8
B	577	ALA	SER	engineered mutation	UNP H2E7Q8

- Molecule 2 is a protein called Alpha-amanitin proprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	17	Total	C	N	O	S	0	0	0
			130	78	22	29	1			
2	C	17	Total	C	N	O	S	0	0	0
			130	78	22	29	1			

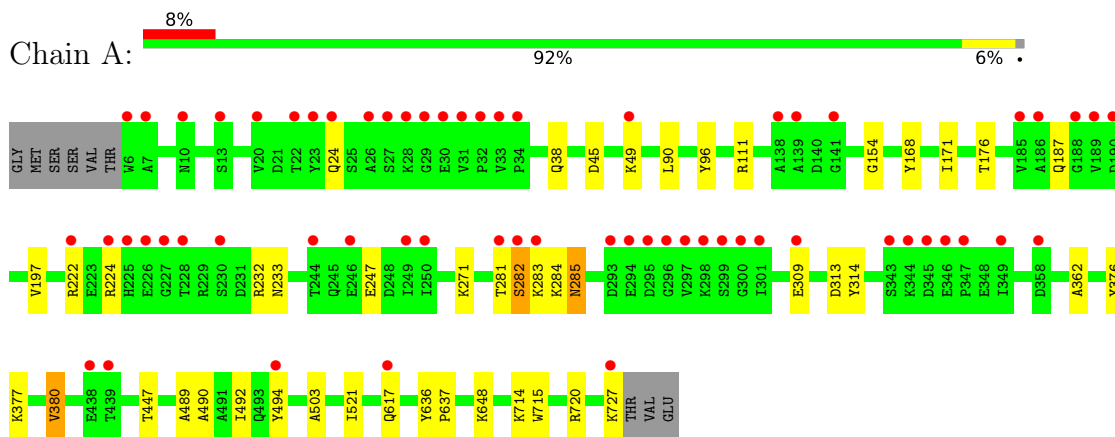
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	954	Total	O	0	0
			954	954		
3	B	939	Total	O	0	0
			939	939		
3	D	24	Total	O	0	0
			24	24		
3	C	24	Total	O	0	0
			24	24		

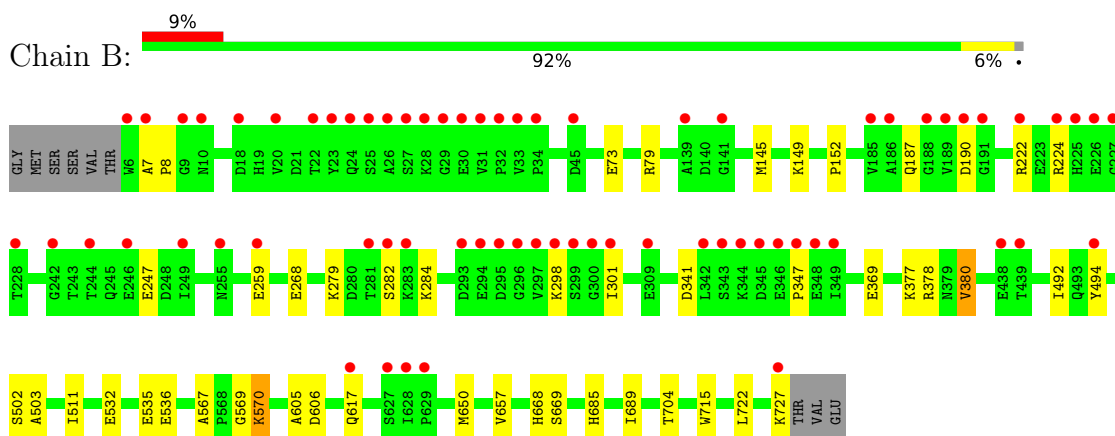
3 Residue-property plots [i](#)

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: Prolyl oligopeptidase



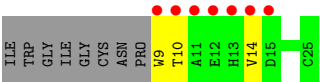
- Molecule 1: Prolyl oligopeptidase



- Molecule 2: Alpha-amanitin proprotein



- Molecule 2: Alpha-amanitin proprotein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	99.08Å 114.72Å 141.26Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.56 – 1.44 43.56 – 1.44	Depositor EDS
% Data completeness (in resolution range)	99.7 (43.56-1.44) 99.7 (43.56-1.44)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.20 (at 1.44Å)	Xtriage
Refinement program	PHENIX (1.11rc2_2522: ???)	Depositor
R, R_{free}	0.155 , 0.181 (Not available) , 0.196	Depositor DCC
R_{free} test set	14696 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	10.7	Xtriage
Anisotropy	0.125	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 36.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	13794	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 52.25 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.9927e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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 ⓘ

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	0.70	6/5993 (0.1%)	0.72	1/8139 (0.0%)
1	B	0.73	8/5994 (0.1%)	0.79	7/8141 (0.1%)
2	C	0.68	0/132	0.88	0/179
2	D	0.58	0/132	0.80	0/179
All	All	0.71	14/12251 (0.1%)	0.76	8/16638 (0.0%)

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	282	SER	N-CA	-12.01	1.30	1.45
1	A	111[A]	ARG	N-CA	9.40	1.57	1.46
1	A	111[B]	ARG	N-CA	9.40	1.57	1.46
1	B	377	LYS	C-N	-8.50	1.22	1.33
1	B	7	ALA	C-O	-8.21	1.15	1.24
1	A	282	SER	N-CA	-7.73	1.36	1.45
1	A	285	ASN	C-O	-6.64	1.15	1.23
1	A	284	LYS	C-O	-6.38	1.16	1.24
1	B	284	LYS	C-O	-6.31	1.16	1.23
1	B	282	SER	CA-C	6.21	1.61	1.53
1	B	570[A]	LYS	N-CA	6.21	1.53	1.46
1	B	570[B]	LYS	N-CA	6.21	1.53	1.46
1	B	8	PRO	C-O	-5.44	1.17	1.23
1	A	285	ASN	CG-ND2	-5.30	1.22	1.33

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	282	SER	N-CA-C	-13.64	92.14	110.55
1	B	377	LYS	CA-C-N	10.00	137.09	123.00
1	B	377	LYS	C-N-CA	10.00	137.09	123.00
1	B	378	ARG	O-C-N	8.52	133.16	123.27
1	A	282	SER	N-CA-C	-8.10	99.08	110.50
1	B	377	LYS	O-C-N	-7.25	114.00	123.13

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	569	GLY	CA-C-N	-5.45	113.90	122.79
1	B	569	GLY	C-N-CA	-5.45	113.90	122.79

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	5796	0	5618	42	0
1	B	5797	0	5621	41	0
2	C	130	0	112	3	0
2	D	130	0	112	4	0
3	A	954	0	0	27	5
3	B	939	0	0	29	5
3	C	24	0	0	1	0
3	D	24	0	0	1	0
All	All	13794	0	11463	87	5

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 (87) 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:490:ALA:O	3:A:801:HOH:O	1.65	1.14
1:B:145:MET:HG2	3:B:1552:HOH:O	1.64	0.96
1:A:24:GLN:NE2	3:A:802:HOH:O	1.96	0.95
1:B:145:MET:HE3	3:B:1552:HOH:O	1.66	0.95
2:C:9:TRP:O	3:C:101:HOH:O	1.91	0.89
1:B:145:MET:SD	3:B:1552:HOH:O	2.38	0.82
1:B:145:MET:CG	3:B:1552:HOH:O	2.23	0.81
1:A:521:ILE:N	3:A:801:HOH:O	2.13	0.81
1:B:145:MET:CE	3:B:1552:HOH:O	2.24	0.81
1:A:38:GLN:NE2	3:A:804:HOH:O	2.11	0.80

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:187:GLN:NE2	3:A:807:HOH:O	2.18	0.77
1:A:224:ARG:NH1	3:A:805:HOH:O	2.12	0.76
1:B:535[A]:GLU:HB2	3:B:1205:HOH:O	1.87	0.75
1:B:650:MET:HE1	3:B:1061:HOH:O	1.85	0.75
1:A:727:LYS:O	3:A:803:HOH:O	2.07	0.72
1:B:247:GLU:OE2	3:B:801:HOH:O	2.06	0.72
1:A:168:TYR:OH	1:A:222:ARG:NH1	2.24	0.70
1:A:377:LYS:N	3:A:809:HOH:O	2.24	0.70
1:A:285:ASN:HD21	1:A:314:TYR:H	1.38	0.69
1:A:720:ARG:HD3	3:A:1465:HOH:O	1.93	0.68
1:B:187[B]:GLN:OE1	3:B:802:HOH:O	2.12	0.68
1:B:259:GLU:OE1	1:B:617:GLN:NE2	2.27	0.67
1:B:268:GLU:OE1	3:B:803:HOH:O	2.12	0.67
1:B:492:ILE:HG13	1:B:715[B]:TRP:CZ3	2.30	0.66
1:B:190:ASP:HB3	3:B:858:HOH:O	1.98	0.64
1:A:281:THR:O	3:A:806:HOH:O	2.15	0.63
1:A:492:ILE:HG12	1:A:715[B]:TRP:CZ3	2.36	0.61
1:A:154:GLY:O	1:A:176[A]:THR:HG21	2.00	0.60
1:B:79[B]:ARG:NH2	1:B:704:THR:HG23	2.16	0.60
1:B:187[A]:GLN:NE2	3:B:808:HOH:O	2.34	0.59
1:B:606:ASP:N	3:B:810:HOH:O	2.35	0.59
1:B:279:LYS:HE2	3:B:1374:HOH:O	2.04	0.58
1:B:492:ILE:HG13	1:B:715[B]:TRP:HZ3	1.69	0.57
1:A:489:ALA:HB1	3:A:801:HOH:O	2.05	0.56
1:A:503:ALA:HB3	3:A:931:HOH:O	2.07	0.55
1:A:309:GLU:HG3	3:A:1451:HOH:O	2.06	0.55
1:B:689:ILE:HD11	3:B:1548:HOH:O	2.06	0.55
1:B:567:ALA:HB3	1:B:570[B]:LYS:HD2	1.89	0.54
1:A:492:ILE:HG12	1:A:715[B]:TRP:HZ3	1.72	0.54
1:B:152:PRO:HG3	3:B:1335:HOH:O	2.09	0.53
2:D:10:THR:HB	2:D:14:VAL:HG23	1.92	0.52
1:B:570[B]:LYS:NZ	1:B:722:LEU:O	2.42	0.51
1:A:282:SER:HB2	3:A:1234:HOH:O	2.10	0.51
1:A:283:LYS:O	1:A:313:ASP:HB2	2.11	0.51
1:B:369:GLU:HG3	3:B:1650:HOH:O	2.11	0.51
1:A:271:LYS:HD2	3:A:908:HOH:O	2.10	0.50
1:B:669:SER:HB2	3:B:1548:HOH:O	2.11	0.50
1:A:521:ILE:HG12	3:A:801:HOH:O	2.12	0.49
2:C:10:THR:HB	2:C:14:VAL:HG22	1.96	0.48
1:A:232:ARG:HH22	1:A:617:GLN:CD	2.21	0.48
1:B:605:ALA:HB3	3:B:810:HOH:O	2.14	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:521:ILE:HG23	3:A:801:HOH:O	2.13	0.47
1:A:362:ALA:HB2	3:A:809:HOH:O	2.14	0.47
1:B:149:LYS:CE	3:B:1340:HOH:O	2.62	0.47
1:B:502[B]:SER:OG	1:B:532:GLU:HG3	2.14	0.47
1:B:73:GLU:HG2	3:B:1281:HOH:O	2.16	0.46
1:B:503:ALA:HB3	3:B:885:HOH:O	2.14	0.45
1:A:362:ALA:CB	3:A:809:HOH:O	2.63	0.45
1:B:511:ILE:HG12	1:B:715[B]:TRP:CZ2	2.51	0.45
1:B:685:HIS:HE1	3:B:1589:HOH:O	1.98	0.45
1:B:657:VAL:HG11	3:B:1548:HOH:O	2.17	0.45
1:A:714:LYS:HE3	1:A:715[A]:TRP:CE2	2.52	0.45
1:B:668:HIS:HA	3:B:810:HOH:O	2.17	0.45
1:A:176[A]:THR:HG23	3:A:952:HOH:O	2.18	0.44
1:A:447[B]:THR:CG2	3:A:1533:HOH:O	2.65	0.44
1:A:447[B]:THR:HG21	3:A:1533:HOH:O	2.18	0.44
1:B:727:LYS:C	3:B:1418:HOH:O	2.61	0.44
2:D:10:THR:HA	3:D:112:HOH:O	2.17	0.44
1:A:376:TYR:C	3:A:809:HOH:O	2.58	0.43
1:A:224:ARG:HD2	1:B:224:ARG:HD2	2.00	0.43
1:A:247:GLU:OE2	3:A:808:HOH:O	2.21	0.43
1:B:535[B]:GLU:HB2	3:B:1205:HOH:O	2.18	0.43
1:A:90:LEU:HD12	1:A:96:TYR:CE1	2.53	0.43
1:B:298:LYS:O	1:B:301:ILE:HG13	2.19	0.42
1:A:492:ILE:HG21	1:A:715[B]:TRP:CH2	2.54	0.42
1:B:494:TYR:CE2	2:C:9:TRP:CZ3	3.08	0.42
1:A:233:ASN:ND2	3:A:842:HOH:O	2.53	0.42
1:A:648:LYS:HG2	3:A:817:HOH:O	2.19	0.42
1:B:222:ARG:NH1	3:B:814:HOH:O	2.38	0.42
1:B:341:ASP:O	1:B:347:PRO:HB3	2.20	0.41
1:A:45[B]:ASP:OD2	1:A:49:LYS:HE3	2.21	0.41
1:A:222:ARG:NE	3:A:812:HOH:O	2.53	0.41
1:A:494:TYR:CE2	2:D:9:TRP:CZ3	3.08	0.41
2:D:10:THR:HB	2:D:14:VAL:CG2	2.51	0.41
1:B:536:GLU:HG2	3:B:1523:HOH:O	2.22	0.40
1:A:636:TYR:HB3	1:A:637:PRO:HD3	2.03	0.40
1:A:171:ILE:HB	1:A:197:VAL:HB	2.03	0.40

All (5) 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)
3:A:840:HOH:O	3:B:1643:HOH:O[3_555]	1.88	0.32
3:A:1668:HOH:O	3:B:1132:HOH:O[4_455]	1.95	0.25
3:A:1523:HOH:O	3:B:1608:HOH:O[3_655]	2.00	0.20
3:A:1609:HOH:O	3:B:1433:HOH:O[3_555]	2.12	0.08
3:A:1622:HOH:O	3:B:1308:HOH:O[3_555]	2.12	0.08

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	733/731 (100%)	713 (97%)	19 (3%)	1 (0%)	48	23
1	B	733/731 (100%)	716 (98%)	16 (2%)	1 (0%)	48	23
2	C	15/25 (60%)	14 (93%)	1 (7%)	0	100	100
2	D	15/25 (60%)	14 (93%)	1 (7%)	0	100	100
All	All	1496/1512 (99%)	1457 (97%)	37 (2%)	2 (0%)	48	23

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	380	VAL
1	A	380	VAL

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	616/612 (101%)	615 (100%)	1 (0%)	87	76
1	B	616/612 (101%)	615 (100%)	1 (0%)	87	76
2	C	14/20 (70%)	14 (100%)	0	100	100
2	D	14/20 (70%)	14 (100%)	0	100	100
All	All	1260/1264 (100%)	1258 (100%)	2 (0%)	87	76

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	380	VAL
1	B	380	VAL

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

Mol	Chain	Res	Type
1	A	64	GLN
1	A	163	HIS
1	A	233	ASN
1	A	255	ASN
1	A	285	ASN
1	A	302	HIS
1	A	379	ASN
1	A	685	HIS
1	B	255	ASN
1	B	334	GLN
1	B	411	GLN
1	B	682	GLN
2	C	16	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	722/731 (98%)	0.38	62 (8%)	16 18	6, 16, 36, 64	13 (1%)
1	B	722/731 (98%)	0.47	69 (9%)	13 15	6, 17, 39, 78	13 (1%)
2	C	17/25 (68%)	3.04	7 (41%)	0 0	16, 24, 95, 112	0
2	D	17/25 (68%)	3.09	6 (35%)	1 1	14, 22, 119, 140	0
All	All	1478/1512 (97%)	0.48	144 (9%)	13 15	6, 17, 39, 140	26 (1%)

All (144) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	9	TRP	10.1
2	C	9	TRP	10.0
2	D	11	ALA	9.3
1	B	189	VAL	8.4
2	C	11	ALA	7.7
2	C	10	THR	7.6
2	C	14	VAL	6.7
2	D	14	VAL	6.5
2	D	13	HIS	6.4
2	D	10	THR	6.4
1	A	6	TRP	5.7
1	A	222	ARG	5.7
2	C	13	HIS	5.3
2	D	12	GLU	5.1
2	C	12	GLU	4.8
1	A	139	ALA	4.7
1	B	32	PRO	4.6
1	A	189	VAL	4.6
1	B	299	SER	4.5
1	B	190	ASP	4.5
1	B	31	VAL	4.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	6	TRP	4.4
1	B	727	LYS	4.4
1	B	295	ASP	4.3
1	B	139	ALA	4.2
1	B	347	PRO	4.2
1	B	228	THR	4.0
1	A	727	LYS	4.0
1	A	227	GLY	3.9
1	B	26	ALA	3.9
1	A	296	GLY	3.8
1	A	228	THR	3.7
1	B	7	ALA	3.7
1	B	349	ILE	3.7
1	B	281	THR	3.7
1	B	298	LYS	3.7
1	A	7	ALA	3.7
1	B	27	SER	3.5
1	B	188	GLY	3.5
1	A	26	ALA	3.4
1	B	348	GLU	3.4
1	A	345	ASP	3.4
1	B	222	ARG	3.4
1	B	294	GLU	3.4
1	A	439	THR	3.3
1	B	10	ASN	3.3
1	B	301	ILE	3.3
1	B	29	GLY	3.3
1	A	347	PRO	3.3
1	B	20	VAL	3.2
1	B	18	ASP	3.2
1	B	438	GLU	3.2
1	B	227	GLY	3.2
1	B	439	THR	3.2
1	B	282	SER	3.2
1	B	344	LYS	3.1
1	B	225	HIS	3.1
1	B	33	VAL	3.1
1	A	22	THR	3.1
1	B	22	THR	3.1
1	A	344	LYS	3.1
1	B	226	GLU	3.0
1	B	191	GLY	3.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	32	PRO	3.0
1	A	299	SER	3.0
1	B	28	LYS	3.0
1	B	224	ARG	2.9
1	A	281	THR	2.9
1	B	283	LYS	2.9
1	B	249	ILE	2.9
1	A	226	GLU	2.9
1	B	343	SER	2.9
1	B	345	ASP	2.8
1	B	346	GLU	2.8
1	A	298	LYS	2.8
1	B	300	GLY	2.8
1	A	20	VAL	2.8
1	A	301	ILE	2.8
1	B	259	GLU	2.8
1	A	29	GLY	2.7
1	A	494	TYR	2.7
1	A	438	GLU	2.7
1	B	23	TYR	2.7
1	A	224	ARG	2.7
1	B	25	SER	2.7
1	B	30	GLU	2.7
1	B	296	GLY	2.6
1	B	342	LEU	2.6
1	A	225	HIS	2.6
1	A	185	VAL	2.6
1	A	246	GLU	2.6
1	A	190	ASP	2.6
1	A	294	GLU	2.5
1	A	10	ASN	2.5
1	A	31	VAL	2.5
1	B	34	PRO	2.5
1	A	27	SER	2.5
1	A	617	GLN	2.5
1	B	297	VAL	2.5
1	B	617	GLN	2.5
1	B	628	ILE	2.5
1	A	300	GLY	2.5
1	B	246	GLU	2.5
1	A	24	GLN	2.5
1	B	185	VAL	2.5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	343	SER	2.4
1	A	23	TYR	2.4
1	A	295	ASP	2.4
1	A	28	LYS	2.4
1	A	244	THR	2.4
1	B	494	TYR	2.4
2	C	15	ASP	2.3
1	A	186	ALA	2.3
1	B	242	GLY	2.3
1	B	293	ASP	2.3
1	A	13	SER	2.3
1	A	33	VAL	2.3
1	B	629	PRO	2.2
1	A	358	ASP	2.2
1	A	141	GLY	2.2
1	A	346	GLU	2.2
1	A	138	ALA	2.2
1	A	283	LYS	2.2
1	A	230	SER	2.2
1	A	282	SER	2.2
1	A	293	ASP	2.2
1	B	186	ALA	2.2
1	A	34	PRO	2.1
1	B	9	GLY	2.1
1	A	30	GLU	2.1
1	B	244	THR	2.1
1	A	249	ILE	2.1
1	A	349	ILE	2.1
1	B	627	SER	2.1
1	A	188	GLY	2.1
1	A	297	VAL	2.1
1	B	24	GLN	2.1
1	A	309	GLU	2.1
1	A	49	LYS	2.0
1	A	250	ILE	2.0
1	B	141	GLY	2.0
1	B	309	GLU	2.0
1	B	45	ASP	2.0
1	B	255	ASN	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.