



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 01:56 pm BST

PDB ID : 4WGV
Title : Crystal structure of Staphylococcus capitis divalent metal ion transporter (DMT) in complex with nanobody
Authors : Ehrnstorfer, I.A.; Geertsma, E.R.; Pardon, E.; Steyaert, J.; Dutzler, R.
Deposited on : 2014-09-19
Resolution : 3.10 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

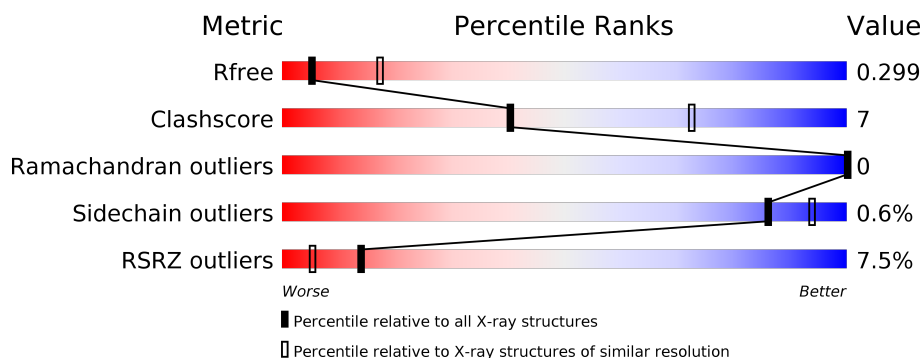
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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	1094 (3.10-3.10)
Clashscore	141614	1184 (3.10-3.10)
Ramachandran outliers	138981	1141 (3.10-3.10)
Sidechain outliers	138945	1141 (3.10-3.10)
RSRZ outliers	127900	1067 (3.10-3.10)

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	415	
1	C	415	
2	B	124	
2	D	124	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7946 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 Divalent metal cation transporter MntH.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	395	Total	C	N	O	S	0	0	0
			3050	2024	485	529	12			
1	C	395	Total	C	N	O	S	0	0	0
			3050	2024	485	529	12			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	41	MET	-	initiating methionine	UNP F9L8R0
A	42	SER	-	expression tag	UNP F9L8R0
A	449	ALA	-	expression tag	UNP F9L8R0
A	450	LEU	-	expression tag	UNP F9L8R0
A	451	GLU	-	expression tag	UNP F9L8R0
A	452	VAL	-	expression tag	UNP F9L8R0
A	453	LEU	-	expression tag	UNP F9L8R0
A	454	PHE	-	expression tag	UNP F9L8R0
A	455	GLN	-	expression tag	UNP F9L8R0
C	41	MET	-	initiating methionine	UNP F9L8R0
C	42	SER	-	expression tag	UNP F9L8R0
C	449	ALA	-	expression tag	UNP F9L8R0
C	450	LEU	-	expression tag	UNP F9L8R0
C	451	GLU	-	expression tag	UNP F9L8R0
C	452	VAL	-	expression tag	UNP F9L8R0
C	453	LEU	-	expression tag	UNP F9L8R0
C	454	PHE	-	expression tag	UNP F9L8R0
C	455	GLN	-	expression tag	UNP F9L8R0

- Molecule 2 is a protein called Camelid antibody fragment, nanobody.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	120	Total	C	N	O	S	0	0	0
			923	564	172	183	4			

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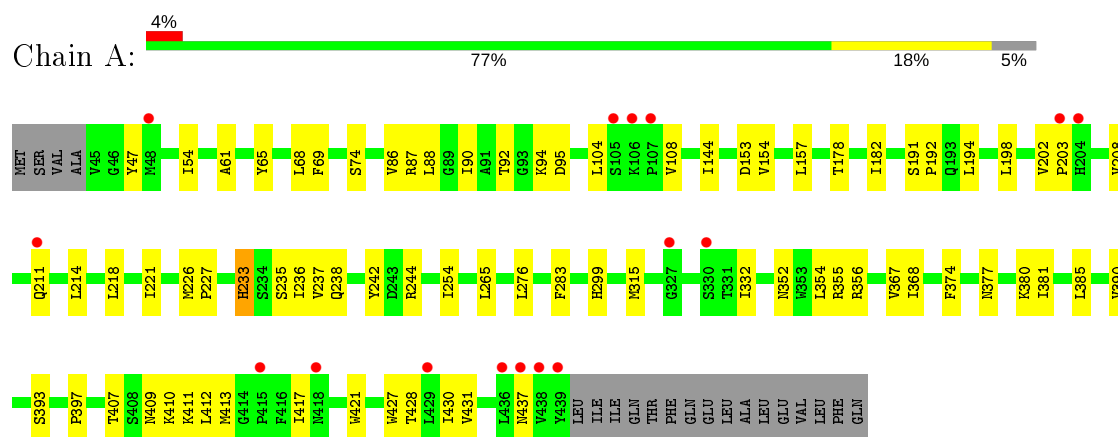
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	120	Total	C	N	O	S	0	0	0
			923	564	172	183	4			

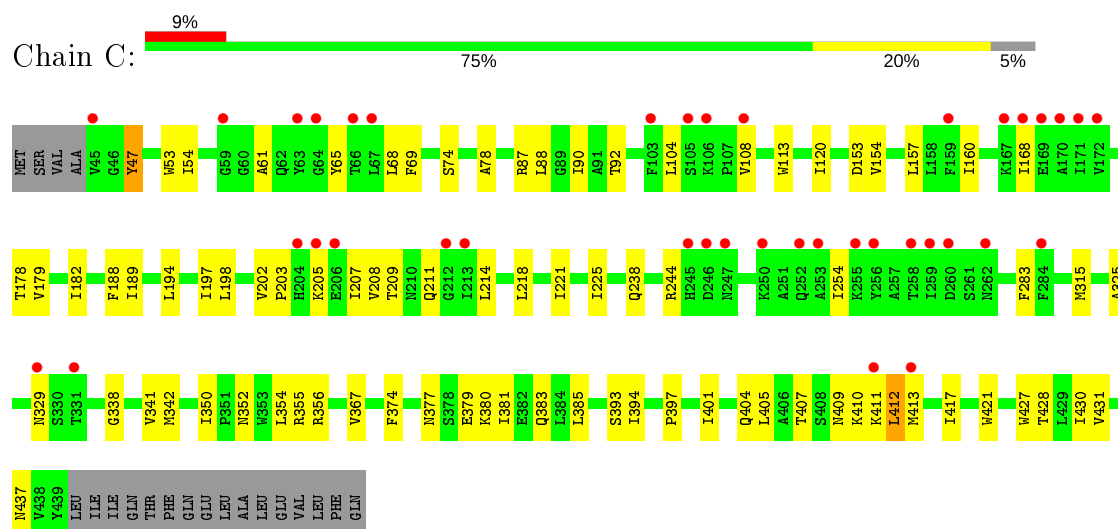
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

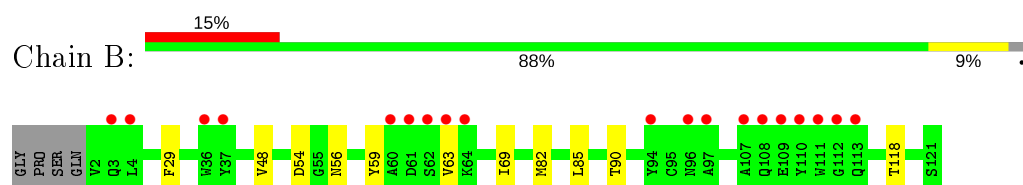
• Molecule 1: Divalent metal cation transporter MntH



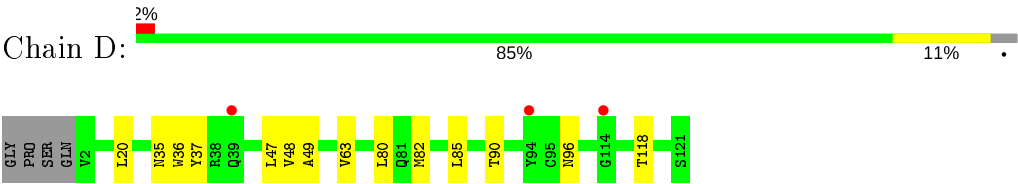
• Molecule 1: Divalent metal cation transporter MntH



• Molecule 2: Camelid antibody fragment, nanobody



● Molecule 2: Camelid antibody fragment, nanobody



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	114.35Å 114.35Å 257.87Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.93 – 3.10 46.22 – 3.10	Depositor EDS
% Data completeness (in resolution range)	99.4 (19.93-3.10) 99.4 (46.22-3.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.73 (at 3.12Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_1760)	Depositor
R, R_{free}	0.250 , 0.286 0.265 , 0.299	Depositor DCC
R_{free} test set	1784 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	118.3	Xtriage
Anisotropy	0.081	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 89.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.026 for -h,-k,l	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	7946	wwPDB-VP
Average B, all atoms (Å ²)	136.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.55% 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](#)

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	0.22	0/3112	0.41	0/4235
1	C	0.24	0/3112	0.44	0/4235
2	B	0.20	0/937	0.37	0/1268
2	D	0.20	0/937	0.37	0/1268
All	All	0.23	0/8098	0.42	0/11006

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 [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	3050	0	3222	49	0
1	C	3050	0	3222	56	0
2	B	923	0	897	6	0
2	D	923	0	897	8	0
All	All	7946	0	8238	116	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 7.

All (116) 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:94:LYS:HD3	1:A:238:GLN:HE22	1.49	0.78
1:C:238:GLN:O	1:C:244:ARG:NH2	2.18	0.77
1:C:417:ILE:HG22	1:C:421:TRP:CD1	2.24	0.72
1:C:205:LYS:O	1:C:209:THR:OG1	2.07	0.72
1:C:188:PHE:O	1:C:189:ILE:HG13	1.94	0.66
1:C:160:ILE:HD11	1:C:168:ILE:HG12	1.77	0.65
1:C:208:VAL:HA	1:C:214:LEU:HD13	1.79	0.65
1:C:104:LEU:HB3	1:C:108:VAL:HG13	1.78	0.65
1:A:354:LEU:HD11	1:C:354:LEU:HD11	1.79	0.65
2:B:48:VAL:HG13	2:B:63:VAL:HG21	1.78	0.64
2:D:48:VAL:HG13	2:D:63:VAL:HG21	1.80	0.64
1:A:238:GLN:O	1:A:244:ARG:NH2	2.31	0.62
1:A:54:ILE:HG21	1:A:385:LEU:HG	1.80	0.62
1:A:157:LEU:HD11	1:A:332:ILE:HD11	1.81	0.62
2:B:90:THR:HG23	2:B:118:THR:HA	1.82	0.62
1:A:211:GLN:NE2	1:A:437:ASN:OD1	2.34	0.60
1:C:211:GLN:NE2	1:C:437:ASN:OD1	2.34	0.60
1:C:74:SER:HB3	1:C:221:ILE:HG12	1.84	0.59
1:C:54:ILE:HG21	1:C:385:LEU:HG	1.84	0.59
2:D:20:LEU:HD12	2:D:80:LEU:HD23	1.85	0.58
2:D:90:THR:HG23	2:D:118:THR:HA	1.85	0.57
1:C:154:VAL:HG13	1:C:356:ARG:HE	1.69	0.57
1:C:417:ILE:HG22	1:C:421:TRP:HD1	1.69	0.57
1:A:74:SER:HB3	1:A:221:ILE:HG12	1.87	0.56
1:C:342:MET:HE1	1:C:355:ARG:HE	1.69	0.56
1:A:104:LEU:HB3	1:A:108:VAL:HG23	1.88	0.56
1:A:90:ILE:HD12	1:A:254:ILE:HG12	1.88	0.56
2:B:82:MET:HB3	2:B:85:LEU:HD21	1.88	0.56
1:C:411:LYS:O	1:C:413:MET:HG2	2.07	0.55
1:A:208:VAL:HA	1:A:214:LEU:HD13	1.87	0.55
1:A:409:ASN:O	1:A:410:LYS:NZ	2.39	0.55
1:A:407:THR:O	1:A:413:MET:HB3	2.06	0.55
1:C:407:THR:O	1:C:413:MET:HB3	2.06	0.54
1:C:409:ASN:O	1:C:410:LYS:NZ	2.39	0.54
1:A:411:LYS:O	1:A:413:MET:HG2	2.07	0.54
1:C:214:LEU:HD11	1:C:430:ILE:HG23	1.88	0.54
1:C:207:ILE:HG13	1:C:208:VAL:HG23	1.90	0.53
2:B:59:TYR:HE1	2:B:69:ILE:H	1.57	0.52
1:A:154:VAL:HG13	1:A:356:ARG:HE	1.72	0.52
1:A:88:LEU:O	1:A:92:THR:OG1	2.25	0.52
1:C:87:ARG:HG2	1:C:254:ILE:HD13	1.91	0.52
1:A:198:LEU:HD22	1:A:203:PRO:HG3	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:352:ASN:OD1	1:C:355:ARG:NH1	2.44	0.51
2:D:82:MET:HB3	2:D:85:LEU:HD21	1.93	0.51
1:A:218:LEU:O	1:A:393:SER:OG	2.29	0.51
1:C:88:LEU:O	1:C:92:THR:OG1	2.24	0.50
1:C:153:ASP:OD1	1:C:154:VAL:N	2.45	0.50
1:C:61:ALA:HB1	1:C:283:PHE:HB2	1.94	0.50
2:D:35:ASN:N	2:D:96:ASN:O	2.35	0.50
1:A:428:THR:HA	1:A:431:VAL:HG22	1.95	0.49
2:D:35:ASN:HB2	2:D:96:ASN:HB3	1.94	0.49
1:A:367:VAL:HG13	1:A:381:ILE:HD11	1.95	0.48
1:C:218:LEU:O	1:C:393:SER:OG	2.31	0.47
1:A:354:LEU:HD21	1:C:354:LEU:HD21	1.95	0.47
1:A:374:PHE:CG	1:A:380:LYS:HG2	2.51	0.46
1:A:182:ILE:HD13	1:A:315:MET:HB3	1.97	0.46
1:A:417:ILE:HG22	1:A:421:TRP:CD1	2.49	0.46
1:A:95:ASP:HB3	1:A:235:SER:HB2	1.97	0.46
1:A:191:SER:HA	1:A:192:PRO:HD3	1.82	0.46
1:A:191:SER:HB2	1:A:194:LEU:HB2	1.97	0.46
1:A:87:ARG:HH12	1:A:407:THR:HG23	1.80	0.46
1:A:61:ALA:HB1	1:A:283:PHE:HB2	1.96	0.46
1:C:350:ILE:O	1:C:355:ARG:NH2	2.48	0.46
1:A:377:ASN:HB3	1:A:380:LYS:HB2	1.97	0.46
1:A:214:LEU:HD11	1:A:430:ILE:HG23	1.96	0.46
1:A:153:ASP:OD1	1:A:154:VAL:N	2.48	0.46
2:B:54:ASP:OD1	2:B:56:ASN:ND2	2.25	0.46
1:C:404:GLN:NE2	1:C:417:ILE:HG23	2.31	0.45
2:D:36:TRP:HB2	2:D:49:ALA:HB3	1.97	0.45
1:A:68:LEU:HG	1:A:276:LEU:HD22	1.98	0.44
1:C:120:ILE:HD13	1:C:338:GLY:HA2	1.98	0.44
1:C:428:THR:HA	1:C:431:VAL:HG22	1.99	0.44
1:C:367:VAL:HG13	1:C:381:ILE:HD11	1.99	0.44
1:A:397:PRO:HG3	1:A:427:TRP:HD1	1.83	0.44
1:C:154:VAL:HG22	1:C:356:ARG:HE	1.82	0.44
1:C:78:ALA:HB2	1:C:225:ILE:HG12	2.00	0.43
1:A:299:HIS:HB3	2:B:29:PHE:HA	2.00	0.43
1:C:47:TYR:O	1:C:53:TRP:NE1	2.45	0.43
1:C:182:ILE:HD13	1:C:315:MET:HB3	1.99	0.43
1:C:397:PRO:HG3	1:C:427:TRP:HD1	1.83	0.43
1:C:401:ILE:O	1:C:405:LEU:HD13	2.19	0.43
1:C:90:ILE:HD12	1:C:254:ILE:HG12	2.01	0.43
1:A:242:TYR:O	1:A:244:ARG:NH1	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:69:PHE:CD1	1:A:202:VAL:HB	2.54	0.43
1:C:377:ASN:HB3	1:C:380:LYS:HB2	2.00	0.42
1:A:233:HIS:HA	1:A:236:ILE:HG22	2.00	0.42
1:C:178:THR:O	1:C:182:ILE:HG13	2.19	0.42
1:C:197:ILE:HG23	1:C:198:LEU:HG	2.00	0.42
1:C:69:PHE:CD1	1:C:202:VAL:HB	2.55	0.42
2:D:37:TYR:CZ	2:D:47:LEU:HD23	2.54	0.42
1:A:244:ARG:HA	1:A:244:ARG:HD3	1.75	0.42
1:A:352:ASN:HA	1:A:355:ARG:HG2	2.02	0.42
1:A:410:LYS:HA	1:A:410:LYS:NZ	2.35	0.42
1:A:86:VAL:HG13	1:A:237:VAL:HG21	2.02	0.42
1:C:157:LEU:O	1:C:160:ILE:HG22	2.18	0.42
1:C:203:PRO:O	1:C:205:LYS:N	2.53	0.42
1:C:394:ILE:HD13	1:C:431:VAL:HB	2.02	0.42
1:C:120:ILE:HD11	1:C:341:VAL:HG21	2.01	0.41
1:C:325:ALA:O	1:C:329:ASN:ND2	2.53	0.41
1:A:178:THR:O	1:A:182:ILE:HG13	2.21	0.41
1:C:188:PHE:C	1:C:189:ILE:HG13	2.40	0.41
1:C:113:TRP:CD1	1:C:341:VAL:HG12	2.56	0.41
1:C:374:PHE:CG	1:C:380:LYS:HG2	2.55	0.41
1:C:410:LYS:HA	1:C:410:LYS:NZ	2.35	0.41
1:A:393:SER:HA	1:A:427:TRP:HZ2	1.86	0.41
1:C:194:LEU:HA	1:C:194:LEU:HD23	1.93	0.41
1:A:144:ILE:HD11	1:A:368:ILE:HG23	2.02	0.41
1:C:179:VAL:HA	1:C:182:ILE:HD12	2.03	0.41
1:C:254:ILE:CD1	1:C:412:LEU:HD21	2.51	0.41
1:A:65:TYR:HB3	1:A:68:LEU:HD12	2.03	0.41
1:C:379:GLU:O	1:C:383:GLN:HG3	2.21	0.41
1:A:226:MET:HA	1:A:227:PRO:HD3	1.87	0.41
1:A:390:VAL:O	1:A:393:SER:HB2	2.21	0.41
1:C:65:TYR:HB3	1:C:68:LEU:HD12	2.03	0.41
1:A:94:LYS:HB3	1:A:95:ASP:H	1.77	0.41
1:A:265:LEU:HA	1:A:265:LEU:HD23	1.95	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	393/415 (95%)	372 (95%)	21 (5%)	0	100	100
1	C	393/415 (95%)	368 (94%)	25 (6%)	0	100	100
2	B	118/124 (95%)	118 (100%)	0	0	100	100
2	D	118/124 (95%)	118 (100%)	0	0	100	100
All	All	1022/1078 (95%)	976 (96%)	46 (4%)	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	333/351 (95%)	330 (99%)	3 (1%)	78	91
1	C	333/351 (95%)	331 (99%)	2 (1%)	86	94
2	B	98/101 (97%)	98 (100%)	0	100	100
2	D	98/101 (97%)	98 (100%)	0	100	100
All	All	862/904 (95%)	857 (99%)	5 (1%)	86	94

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	47	TYR
1	A	233	HIS

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Mol	Chain	Res	Type
1	A	412	LEU
1	C	47	TYR
1	C	412	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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](#)

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	395/415 (95%)	-0.01	16 (4%) 37 18	77, 116, 176, 270	0
1	C	395/415 (95%)	0.31	39 (9%) 7 2	83, 150, 231, 293	0
2	B	120/124 (96%)	0.57	19 (15%) 2 1	88, 115, 147, 178	0
2	D	120/124 (96%)	-0.00	3 (2%) 57 34	88, 118, 158, 191	0
All	All	1030/1078 (95%)	0.18	77 (7%) 14 5	77, 125, 206, 293	0

All (77) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	246	ASP	11.7
1	C	411	LYS	7.5
1	C	108	VAL	6.8
1	C	259	ILE	6.5
1	A	204	HIS	5.5
2	B	113	GLN	5.1
1	C	204	HIS	5.0
1	C	169	GLU	4.9
1	C	245	HIS	4.8
2	B	111	TRP	4.7
2	B	108	GLN	4.6
2	B	4	LEU	4.6
2	B	3	GLN	4.4
1	C	168	ILE	4.4
2	B	94	TYR	4.3
2	B	37	TYR	4.1
1	C	256	TYR	4.0
1	A	437	ASN	4.0
1	A	327	GLY	3.9
2	D	94	TYR	3.8
1	A	106	LYS	3.8

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Mol	Chain	Res	Type	RSRZ
1	C	159	PHE	3.8
1	A	211	GLN	3.7
1	C	253	ALA	3.7
1	A	415	PRO	3.6
2	B	112	GLY	3.6
1	C	63	TYR	3.6
2	B	110	TYR	3.5
2	B	62	SER	3.4
1	C	247	ASN	3.2
1	C	67	LEU	3.2
1	A	438	VAL	3.2
1	C	262	ASN	3.2
2	B	64	LYS	3.1
1	C	213	ILE	3.1
1	A	439	TYR	3.0
1	A	436	LEU	3.0
1	C	252	GLN	3.0
1	C	59	GLY	3.0
1	C	329	ASN	2.8
2	B	60	ALA	2.7
1	C	170	ALA	2.7
1	C	212	GLY	2.7
1	C	45	VAL	2.7
2	B	97	ALA	2.7
2	B	107	ALA	2.6
1	C	66	THR	2.6
1	C	103	PHE	2.6
1	C	64	GLY	2.6
2	B	109	GLU	2.6
1	C	260	ASP	2.6
2	B	61	ASP	2.5
1	A	418	ASN	2.5
1	C	205	LYS	2.5
1	C	331	THR	2.4
1	A	203	PRO	2.4
1	C	172	VAL	2.4
1	C	167	LYS	2.4
1	C	413	MET	2.4
1	A	429	LEU	2.3
2	D	114	GLY	2.3
1	C	206	GLU	2.3
2	B	96	ASN	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	105	SER	2.2
2	D	39	GLN	2.2
1	C	250	LYS	2.2
1	A	330	SER	2.2
1	A	48	MET	2.1
2	B	63	VAL	2.1
1	C	106	LYS	2.1
1	C	255	LYS	2.1
1	C	171	ILE	2.1
2	B	36	TRP	2.1
1	A	107	PRO	2.0
1	C	258	THR	2.0
1	C	284	PHE	2.0
1	A	105	SER	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 ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.