



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 29, 2020 – 06:27 PM BST

PDB ID : 6TFO
Title : Crystal structure of as isolated three-domain copper-containing nitrite reductase from *Hyphomicrobium denitrificans* strain 1NES1
Authors : Sasaki, D.; Watanabe, T.F.; Eady, R.R.; Garratt, R.C.; Antonyuk, S.V.; Hasnain, S.S.
Deposited on : 2019-11-14
Resolution : 2.05 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	:	2.13
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)	:	2.13

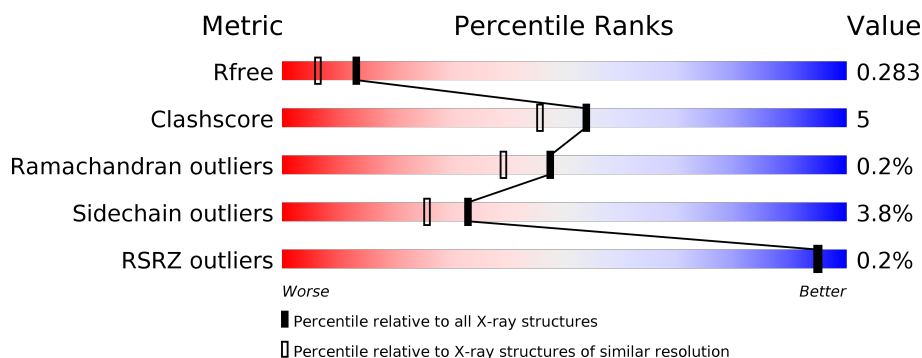
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION




The reported resolution of this entry is 2.05 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	1692 (2.04-2.04)
Clashscore	141614	1773 (2.04-2.04)
Ramachandran outliers	138981	1752 (2.04-2.04)
Sidechain outliers	138945	1752 (2.04-2.04)
RSRZ outliers	127900	1672 (2.04-2.04)

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	456	 81% 11% • 7%
1	B	456	 79% 12% • 7%
1	C	456	 82% 11% • 7%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10167 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 Copper-containing nitrite reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	423	Total	C	N	O	S	0	1	0
			3212	2042	550	610	10			
1	B	422	Total	C	N	O	S	0	1	0
			3206	2039	548	609	10			
1	C	424	Total	C	N	O	S	0	0	0
			3212	2042	550	610	10			

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP N0B9M5
A	451	GLU	-	expression tag	UNP N0B9M5
A	452	ASN	-	expression tag	UNP N0B9M5
A	453	LEU	-	expression tag	UNP N0B9M5
A	454	TYR	-	expression tag	UNP N0B9M5
A	455	PHE	-	expression tag	UNP N0B9M5
A	456	GLN	-	expression tag	UNP N0B9M5
B	1	MET	-	initiating methionine	UNP N0B9M5
B	451	GLU	-	expression tag	UNP N0B9M5
B	452	ASN	-	expression tag	UNP N0B9M5
B	453	LEU	-	expression tag	UNP N0B9M5
B	454	TYR	-	expression tag	UNP N0B9M5
B	455	PHE	-	expression tag	UNP N0B9M5
B	456	GLN	-	expression tag	UNP N0B9M5
C	1	MET	-	initiating methionine	UNP N0B9M5
C	451	GLU	-	expression tag	UNP N0B9M5
C	452	ASN	-	expression tag	UNP N0B9M5
C	453	LEU	-	expression tag	UNP N0B9M5
C	454	TYR	-	expression tag	UNP N0B9M5
C	455	PHE	-	expression tag	UNP N0B9M5
C	456	GLN	-	expression tag	UNP N0B9M5

- Molecule 2 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

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

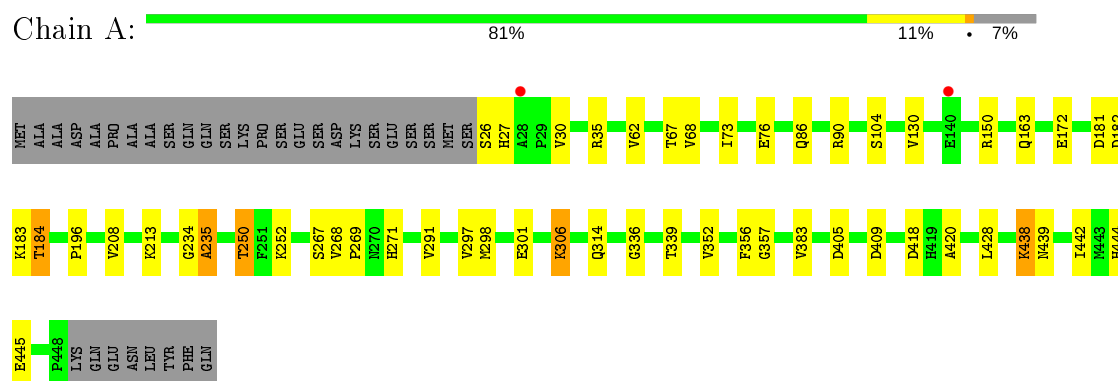
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	186	Total 186	O 186	0	0
3	B	173	Total 173	O 173	0	0
3	C	169	Total 169	O 169	0	0

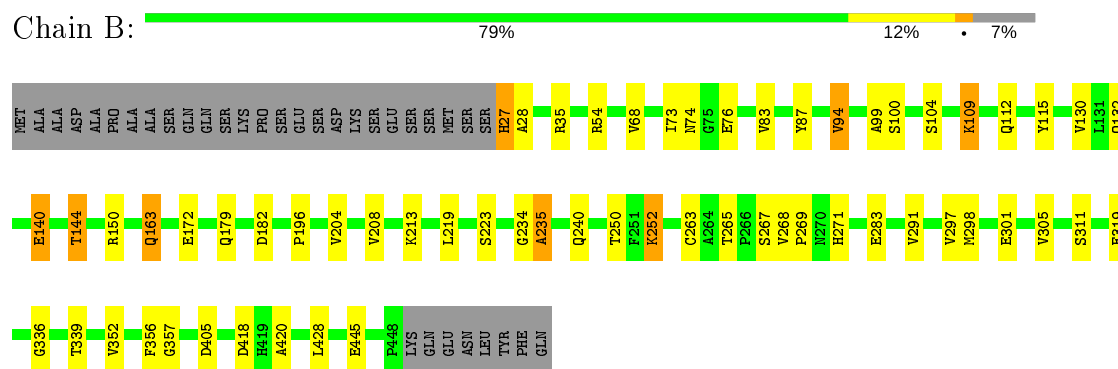
3 Residue-property plots [i](#)

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.

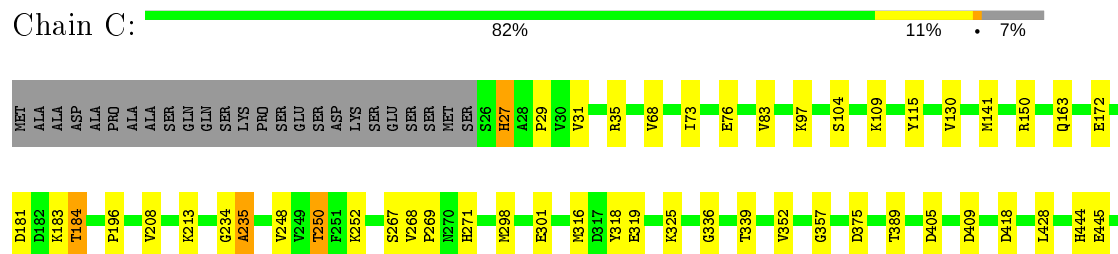
- Molecule 1: Copper-containing nitrite reductase



- Molecule 1: Copper-containing nitrite reductase



- Molecule 1: Copper-containing nitrite reductase



P448	GLN
K449	GLU
	ASN
	LEU
	TYR
	PHE
	GLN

4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	77.07Å 77.07Å 754.55Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	66.58 – 2.05 66.49 – 2.05	Depositor EDS
% Data completeness (in resolution range)	98.3 (66.58-2.05) 98.3 (66.49-2.05)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.39 (at 2.05Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.227 , 0.279 0.235 , 0.283	Depositor DCC
R_{free} test set	4195 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	38.4	Xtriage
Anisotropy	0.185	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 38.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10167	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.50% 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: CU

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	0/3292	0.84	0/4484
1	B	0.68	0/3286	0.84	1/4476 (0.0%)
1	C	0.71	0/3292	0.84	0/4484
All	All	0.70	0/9870	0.84	1/13444 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
All	All	0	2

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	182	ASP	CB-CA-C	5.11	120.63	110.40

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	420	ALA	Peptide
1	B	420	ALA	Peptide

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	3212	0	3138	29	0
1	B	3206	0	3131	44	0
1	C	3212	0	3139	33	0
2	A	3	0	0	0	0
2	B	3	0	0	0	0
2	C	3	0	0	0	0
3	A	186	0	0	4	0
3	B	173	0	0	4	0
3	C	169	0	0	4	0
All	All	10167	0	9408	101	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 5.

All (101) 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:445:GLU:O	1:C:248:VAL:O	1.87	0.92
1:C:181:ASP:HB3	1:C:184:THR:HG23	1.54	0.88
1:A:181:ASP:HB3	1:A:184:THR:HG23	1.53	0.88
1:B:240:GLN:OE1	1:B:265:THR:HG22	1.73	0.86
1:B:74:ASN:HB2	1:B:94:VAL:HG13	1.67	0.76
1:B:28:ALA:HA	3:B:726:HOH:O	1.89	0.72
1:B:204:VAL:HG23	1:B:283:GLU:O	1.95	0.67
1:B:163:GLN:NE2	1:B:163:GLN:H	1.93	0.65
1:B:27:HIS:O	1:B:27:HIS:ND1	2.30	0.64
1:C:27:HIS:ND1	1:C:27:HIS:O	2.31	0.63
1:A:442:ILE:HA	1:B:252:LYS:HG3	1.85	0.59
1:C:141:MET:HG3	1:C:183:LYS:HE2	1.84	0.59
1:B:172:GLU:HA	1:B:213:LYS:O	2.04	0.57
1:A:26:SER:HA	3:A:678:HOH:O	2.05	0.57
1:A:172:GLU:HA	1:A:213:LYS:O	2.05	0.57
1:B:319[A]:GLU:HG3	3:B:737:HOH:O	2.05	0.57
1:B:223:SER:OG	1:B:265:THR:HG23	2.05	0.56
1:C:172:GLU:HA	1:C:213:LYS:O	2.04	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:316:MET:CE	1:C:318:TYR:N	2.70	0.54
1:A:182:ASP:HA	3:A:601:HOH:O	2.06	0.53
1:A:439:ASN:HD21	1:A:445:GLU:HG2	1.74	0.53
1:B:94:VAL:HG22	1:B:99:ALA:HB1	1.91	0.53
1:B:94:VAL:HG22	1:B:99:ALA:CB	2.39	0.53
1:A:439:ASN:HB2	3:A:695:HOH:O	2.09	0.52
1:B:27:HIS:O	1:B:27:HIS:CG	2.62	0.51
1:B:298:MET:HA	1:B:357:GLY:O	2.10	0.51
1:A:30:VAL:HG11	1:A:62:VAL:HG22	1.93	0.50
1:C:29:PRO:HD2	3:C:725:HOH:O	2.11	0.50
1:A:298:MET:HA	1:A:357:GLY:O	2.11	0.49
1:C:298:MET:HA	1:C:357:GLY:O	2.10	0.49
1:A:35:ARG:HA	1:A:73:ILE:O	2.12	0.49
1:B:35:ARG:HA	1:B:73:ILE:O	2.12	0.49
1:C:27:HIS:O	1:C:27:HIS:CG	2.65	0.49
1:C:316:MET:CE	1:C:318:TYR:HA	2.42	0.49
1:B:74:ASN:HB2	1:B:94:VAL:CG1	2.41	0.48
1:C:27:HIS:HB3	3:C:714:HOH:O	2.14	0.48
1:C:35:ARG:HA	1:C:73:ILE:O	2.14	0.48
1:A:68:VAL:O	1:A:104:SER:HA	2.14	0.47
1:B:54:ARG:HD2	3:B:728:HOH:O	2.14	0.47
1:B:208:VAL:O	1:B:250:THR:HA	2.15	0.47
1:B:163:GLN:HE21	1:B:163:GLN:H	1.62	0.47
1:A:268:VAL:N	1:A:269:PRO:CD	2.78	0.47
1:C:316:MET:HE1	1:C:318:TYR:HA	1.97	0.47
1:A:208:VAL:O	1:A:250:THR:HA	2.15	0.47
1:A:267:SER:O	1:A:271:HIS:HD2	1.98	0.47
1:B:352:VAL:O	1:B:405:ASP:HA	2.14	0.47
1:C:208:VAL:O	1:C:250:THR:HA	2.15	0.47
1:A:183:LYS:N	3:A:601:HOH:O	2.20	0.46
1:B:68:VAL:O	1:B:104:SER:HA	2.14	0.46
1:C:68:VAL:O	1:C:104:SER:HA	2.15	0.46
1:C:268:VAL:N	1:C:269:PRO:CD	2.78	0.46
1:B:263:CYS:SG	1:B:265:THR:OG1	2.70	0.46
1:B:268:VAL:N	1:B:269:PRO:CD	2.79	0.46
1:B:267:SER:O	1:B:271:HIS:HD2	2.00	0.45
1:B:94:VAL:HG22	1:B:99:ALA:C	2.37	0.45
1:C:267:SER:O	1:C:271:HIS:HD2	1.99	0.45
1:C:352:VAL:O	1:C:405:ASP:HA	2.16	0.45
1:A:352:VAL:O	1:A:405:ASP:HA	2.17	0.44
1:A:444:HIS:HD2	1:A:445:GLU:O	2.00	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:87:TYR:CZ	1:B:109:LYS:HG3	2.51	0.44
1:C:150:ARG:HB3	1:C:196:PRO:HD2	2.00	0.44
1:B:418:ASP:HB3	1:B:428:LEU:HD23	1.99	0.44
1:B:150:ARG:HB3	1:B:196:PRO:HD2	2.00	0.44
1:B:94:VAL:CG2	1:B:99:ALA:C	2.85	0.44
1:C:316:MET:HE1	1:C:318:TYR:CA	2.48	0.44
1:C:444:HIS:HE1	3:C:707:HOH:O	2.00	0.44
1:B:112:GLN:HG2	1:B:132:GLN:HG3	1.99	0.44
1:B:305:VAL:HB	3:B:613:HOH:O	2.18	0.44
1:C:418:ASP:HB3	1:C:428:LEU:HD23	2.00	0.44
1:C:444:HIS:HD2	1:C:445:GLU:O	2.01	0.43
1:C:444:HIS:CE1	3:C:707:HOH:O	2.72	0.43
1:A:418:ASP:HB3	1:A:428:LEU:HD23	2.00	0.43
1:B:140:GLU:OE1	1:B:140:GLU:O	2.36	0.43
1:A:26:SER:N	1:A:67:THR:HG1	2.16	0.43
1:B:204:VAL:CG2	1:B:283:GLU:O	2.65	0.43
1:B:83:VAL:O	1:B:115:TYR:HA	2.19	0.43
1:C:375:ASP:O	1:C:389:THR:HA	2.19	0.42
1:B:144:THR:HG23	1:B:179:GLN:HE22	1.83	0.42
1:C:234:GLY:O	1:C:235:ALA:HB3	2.20	0.42
1:A:150:ARG:HB3	1:A:196:PRO:HD2	2.01	0.42
1:B:297:VAL:O	1:B:356:PHE:HA	2.20	0.42
1:A:27:HIS:N	1:A:27:HIS:CD2	2.87	0.42
1:C:316:MET:CE	1:C:318:TYR:CA	2.98	0.42
1:A:90:ARG:HD2	1:B:219:LEU:HD22	2.02	0.41
1:A:234:GLY:O	1:A:235:ALA:HB3	2.20	0.41
1:A:297:VAL:O	1:A:356:PHE:HA	2.21	0.41
1:B:100:SER:HB2	1:C:316:MET:HE2	2.03	0.41
1:B:336:GLY:HA2	1:B:339:THR:OG1	2.20	0.41
1:C:83:VAL:O	1:C:115:TYR:HA	2.21	0.41
1:B:112:GLN:HB3	1:B:112:GLN:HE21	1.67	0.41
1:C:336:GLY:HA2	1:C:339:THR:OG1	2.21	0.41
1:A:250:THR:HG21	1:C:448:PRO:CD	2.51	0.41
1:A:438:LYS:HG2	1:A:439:ASN:N	2.35	0.41
1:B:234:GLY:O	1:B:235:ALA:HB3	2.21	0.41
1:C:27:HIS:ND1	1:C:27:HIS:C	2.73	0.41
1:A:336:GLY:HA2	1:A:339:THR:OG1	2.21	0.41
1:B:223:SER:CB	1:B:265:THR:HG23	2.50	0.41
1:C:234:GLY:O	1:C:235:ALA:CB	2.69	0.41
1:A:86:GLN:H	1:A:86:GLN:CD	2.23	0.40
1:B:27:HIS:C	1:B:27:HIS:ND1	2.74	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:306:LYS:HB2	1:A:314:GLN:OE1	2.22	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	422/456 (92%)	408 (97%)	13 (3%)	1 (0%)	47	39
1	B	421/456 (92%)	407 (97%)	13 (3%)	1 (0%)	47	39
1	C	422/456 (92%)	408 (97%)	13 (3%)	1 (0%)	47	39
All	All	1265/1368 (92%)	1223 (97%)	39 (3%)	3 (0%)	47	39

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	235	ALA
1	B	235	ALA
1	C	235	ALA

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	343/371 (92%)	331 (96%)	12 (4%)	36	29

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	342/371 (92%)	330 (96%)	12 (4%)	36	29
1	C	343/371 (92%)	328 (96%)	15 (4%)	28	21
All	All	1028/1113 (92%)	989 (96%)	39 (4%)	33	26

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	76	GLU
1	A	130	VAL
1	A	163	GLN
1	A	184	THR
1	A	250	THR
1	A	252	LYS
1	A	291	VAL
1	A	301	GLU
1	A	306	LYS
1	A	383	VAL
1	A	409	ASP
1	A	438	LYS
1	B	27	HIS
1	B	76	GLU
1	B	94	VAL
1	B	109	LYS
1	B	130	VAL
1	B	140	GLU
1	B	144	THR
1	B	163	GLN
1	B	252	LYS
1	B	291	VAL
1	B	301	GLU
1	B	311	SER
1	C	27	HIS
1	C	31	VAL
1	C	76	GLU
1	C	97	LYS
1	C	109	LYS
1	C	130	VAL
1	C	163	GLN
1	C	184	THR
1	C	250	THR
1	C	252	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	301	GLU
1	C	319	GLU
1	C	325	LYS
1	C	409	ASP
1	C	449	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	27	HIS
1	A	112	GLN
1	A	132	GLN
1	A	274	ASN
1	A	290	GLN
1	A	444	HIS
1	B	112	GLN
1	B	163	GLN
1	B	274	ASN
1	C	112	GLN
1	C	132	GLN
1	C	274	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 9 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

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 [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	423/456 (92%)	-0.23	2 (0%) 91 92	35, 47, 68, 93	0
1	B	422/456 (92%)	-0.20	0 100 100	38, 50, 74, 105	0
1	C	424/456 (92%)	-0.20	0 100 100	36, 50, 73, 117	0
All	All	1269/1368 (92%)	-0.21	2 (0%) 95 95	35, 49, 72, 117	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	140	GLU	2.7
1	A	28	ALA	2.7

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 monosaccharides 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	CU	A	502	1/1	0.97	0.13	49,49,49,49	1

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	CU	B	503	1/1	0.99	0.10	46,46,46,46	0
2	CU	C	502	1/1	0.99	0.08	53,53,53,53	1
2	CU	C	501	1/1	0.99	0.08	45,45,45,45	0
2	CU	B	502	1/1	0.99	0.07	47,47,47,47	1
2	CU	C	503	1/1	0.99	0.11	44,44,44,44	0
2	CU	A	501	1/1	1.00	0.12	41,41,41,41	0
2	CU	A	503	1/1	1.00	0.12	48,48,48,48	0
2	CU	B	501	1/1	1.00	0.11	46,46,46,46	0

6.5 Other polymers [i](#)

There are no such residues in this entry.