Full wwPDB X-ray Structure Validation Report ⓘ

Oct 26, 2023 – 10:00 PM EDT

PDB ID	:	3NTM
Title	:	Crystal Structure of Tyrosinase from Bacillus megaterium crystallized in the absence of zinc, partial occupancy of CuB
Authors	:	Sendovski, M.; Kanteev, M.; Adir, N.; Fishman, A.
Deposited on	:	2010-07-05
Resolution	:	2.30 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:

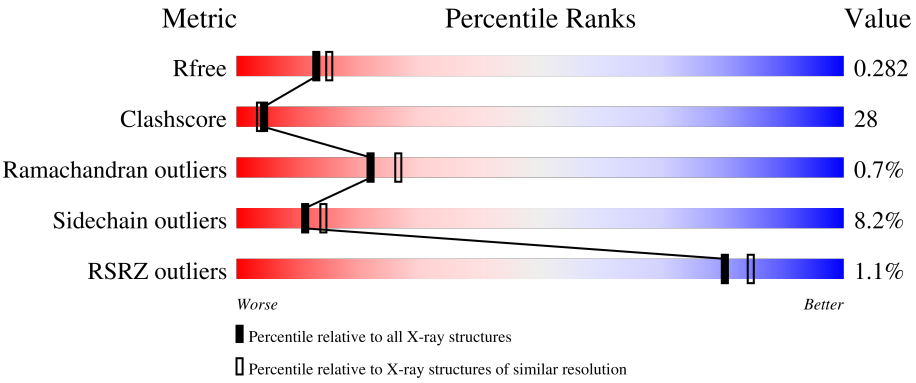
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 <=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	303	% 52% 35% 5% 8%
1	B	303	% 55% 33% 6% 6%

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
2	CU	A	502	-	-	-	X

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 4850 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 Tyrosinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	279	Total	C	N	O	S	0	0	0
			2282	1453	413	409	7			
1	B	284	Total	C	N	O	S	0	0	0
			2329	1483	419	419	8			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	298	HIS	-	expression tag	UNP B2ZB02
A	299	HIS	-	expression tag	UNP B2ZB02
A	300	HIS	-	expression tag	UNP B2ZB02
A	301	HIS	-	expression tag	UNP B2ZB02
A	302	HIS	-	expression tag	UNP B2ZB02
A	303	HIS	-	expression tag	UNP B2ZB02
B	298	HIS	-	expression tag	UNP B2ZB02
B	299	HIS	-	expression tag	UNP B2ZB02
B	300	HIS	-	expression tag	UNP B2ZB02
B	301	HIS	-	expression tag	UNP B2ZB02
B	302	HIS	-	expression tag	UNP B2ZB02
B	303	HIS	-	expression tag	UNP B2ZB02

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Cu	0	0
			2	2		
2	B	1	Total	Cu	0	0
			1	1		

- Molecule 3 is water.

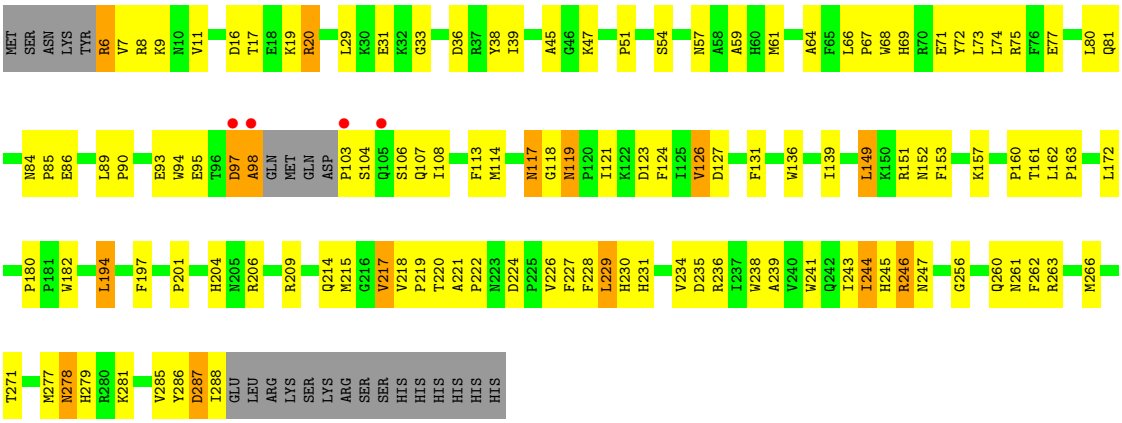
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	119	Total O 119 119	0	0
3	B	117	Total O 117 117	0	0

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.

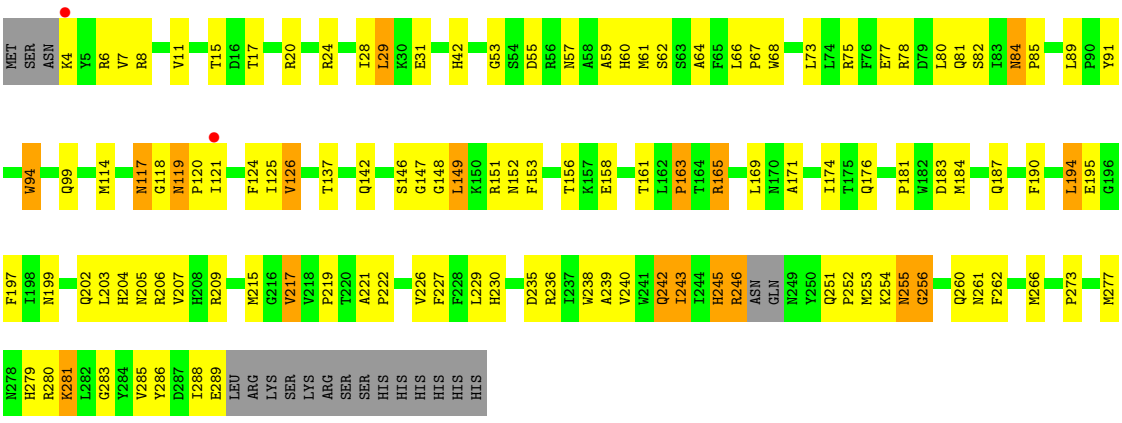
• Molecule 1: Tyrosinase

Chain A:



• Molecule 1: Tyrosinase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	46.01Å 78.48Å 74.55Å 90.00° 101.61° 90.00°	Depositor
Resolution (Å)	39.08 – 2.30 39.24 – 2.30	Depositor EDS
% Data completeness (in resolution range)	93.4 (39.08-2.30) 95.2 (39.24-2.30)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$< I/\sigma(I) >$ [1]	3.98 (at 2.29Å)	Xtriage
Refinement program	REFMAC, PHENIX 1.6.1_357	Depositor
R, R_{free}	0.193 , 0.289 0.197 , 0.282	Depositor DCC
R_{free} test set	1137 reflections (5.15%)	wwPDB-VP
Wilson B-factor (Å^2)	17.2	Xtriage
Anisotropy	0.507	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 54.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.43$, $< L^2 > = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.91	EDS
Total number of atoms	4850	wwPDB-VP
Average B, all atoms (Å^2)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.20% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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	0.57	2/2358 (0.1%)	0.66	1/3213 (0.0%)
1	B	0.53	2/2406 (0.1%)	0.65	2/3277 (0.1%)
All	All	0.55	4/4764 (0.1%)	0.65	3/6490 (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	B	0	2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	244	ILE	N-CA	5.88	1.58	1.46
1	B	245	HIS	C-N	-5.68	1.21	1.34
1	A	245	HIS	CA-CB	-5.64	1.41	1.53
1	B	242	GLN	C-N	-5.08	1.22	1.34

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	245	HIS	C-N-CA	6.54	138.06	121.70
1	B	245	HIS	O-C-N	-6.04	113.03	122.70
1	A	98	ALA	CB-CA-C	-5.53	101.81	110.10

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	245	HIS	Peptide
1	B	246	ARG	Sidechain

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	2282	0	2170	129	1
1	B	2329	0	2210	129	1
2	A	2	0	0	0	0
2	B	1	0	0	0	0
3	A	119	0	0	28	0
3	B	117	0	0	16	0
All	All	4850	0	4380	256	2

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 28.

All (256) 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:242:GLN:O	1:B:246:ARG:HA	1.34	1.20
1:A:6:ARG:N	1:A:7:VAL:HA	1.60	1.09
1:B:254:LYS:O	1:B:255:ASN:HB2	1.39	1.07
1:A:151:ARG:HD3	3:A:368:HOH:O	1.53	1.06
1:B:215:MET:HE2	1:B:221:ALA:HB1	1.32	1.04
1:B:215:MET:HE1	1:B:227:PHE:HD2	1.18	1.04
1:B:215:MET:CE	1:B:221:ALA:HB1	1.91	1.00
1:B:242:GLN:O	1:B:246:ARG:HG3	1.63	0.99
1:A:71:GLU:HA	3:A:362:HOH:O	1.69	0.92
1:A:215:MET:HE2	1:A:221:ALA:HB1	1.50	0.92
1:B:215:MET:HE1	1:B:227:PHE:CD2	2.04	0.92
1:A:72:TYR:HA	3:A:363:HOH:O	1.68	0.92
1:A:201:PRO:HG3	1:A:209:ARG:HD2	1.53	0.90
1:B:125:ILE:HD11	1:B:148:GLY:HA3	1.52	0.90
1:B:254:LYS:O	1:B:255:ASN:CB	2.18	0.89
1:B:68:TRP:N	1:B:266:MET:HE1	1.89	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:108:ILE:HD11	1:A:229:LEU:HD23	1.56	0.87
1:A:204:HIS:CD2	1:A:230:HIS:HE1	1.92	0.87
1:B:242:GLN:O	1:B:246:ARG:CA	2.22	0.87
1:B:227:PHE:CE2	3:B:360:HOH:O	2.29	0.84
1:B:255:ASN:HB2	3:B:337:HOH:O	1.77	0.84
1:B:24:ARG:HG2	1:B:24:ARG:HH11	1.42	0.84
1:A:75:ARG:HD3	3:A:363:HOH:O	1.77	0.83
1:B:8:ARG:HG2	1:B:89:LEU:O	1.80	0.82
1:A:215:MET:HE1	1:A:227:PHE:HD2	1.45	0.81
1:A:204:HIS:HD2	1:A:230:HIS:HE1	1.26	0.81
1:A:204:HIS:HD2	1:A:230:HIS:CE1	2.00	0.79
1:A:215:MET:CE	1:A:221:ALA:HB1	2.12	0.78
1:A:6:ARG:HG3	1:A:6:ARG:O	1.84	0.78
1:B:163:PRO:HD3	1:B:206:ARG:HG2	1.67	0.77
1:B:204:HIS:HD2	1:B:230:HIS:HE1	1.32	0.76
1:A:256:GLY:H	1:A:261:ASN:HD21	1.33	0.76
1:A:153:PHE:CZ	3:A:368:HOH:O	2.36	0.76
1:B:81:GLN:HE21	1:B:85:PRO:HA	1.49	0.75
1:A:214:GLN:HB2	3:A:372:HOH:O	1.86	0.75
1:B:117:ASN:HD21	1:B:153:PHE:H	1.35	0.74
1:A:215:MET:HE1	1:A:227:PHE:CD2	2.23	0.73
1:A:219:PRO:HB3	3:A:359:HOH:O	1.87	0.72
1:B:203:LEU:O	1:B:207:VAL:HG23	1.89	0.72
1:B:78:ARG:NH2	3:B:316:HOH:O	2.21	0.72
1:B:277:MET:HE1	3:B:306:HOH:O	1.90	0.72
1:A:103:PRO:HG3	3:A:376:HOH:O	1.89	0.71
1:B:6:ARG:HB2	1:B:283:GLY:O	1.90	0.71
1:B:256:GLY:N	1:B:261:ASN:HD21	1.89	0.71
1:B:222:PRO:HD3	3:B:360:HOH:O	1.90	0.71
1:A:278:ASN:HD22	1:A:281:LYS:H	1.39	0.70
1:A:114:MET:HG3	1:A:226:VAL:HG22	1.73	0.70
1:A:64:ALA:H	1:A:260:GLN:NE2	1.90	0.69
1:B:149:LEU:HG	3:B:382:HOH:O	1.92	0.69
1:B:204:HIS:HD2	1:B:230:HIS:CE1	2.11	0.69
1:B:62:SER:HB3	3:B:366:HOH:O	1.91	0.69
1:B:161:THR:HB	3:B:384:HOH:O	1.92	0.69
1:A:11:VAL:HG22	1:A:90:PRO:HB2	1.74	0.68
1:B:57:ASN:HD21	1:B:60:HIS:HD1	1.43	0.67
1:A:6:ARG:N	1:A:285:VAL:H	1.92	0.67
1:B:24:ARG:HG2	1:B:24:ARG:NH1	2.09	0.67
1:B:161:THR:HG23	1:B:206:ARG:NH1	2.10	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:242:GLN:O	1:B:246:ARG:CG	2.41	0.66
1:B:221:ALA:HB3	3:B:360:HOH:O	1.94	0.66
1:B:64:ALA:H	1:B:260:GLN:HE22	1.44	0.66
1:B:64:ALA:HB1	1:B:266:MET:CE	2.26	0.65
1:B:165:ARG:HG3	1:B:169:LEU:HD13	1.79	0.65
1:A:113:PHE:HD1	3:A:367:HOH:O	1.80	0.64
1:A:215:MET:CE	1:A:227:PHE:HD2	2.10	0.64
1:A:98:ALA:HB2	1:A:163:PRO:O	1.97	0.64
1:B:99:GLN:HA	3:B:376:HOH:O	1.98	0.64
1:A:126:VAL:HG13	1:A:149:LEU:HB3	1.80	0.63
1:A:8:ARG:CG	1:A:89:LEU:O	2.47	0.63
1:B:281:LYS:HG3	1:B:281:LYS:O	1.98	0.63
1:A:119:ASN:ND2	1:A:121:ILE:H	1.96	0.63
1:B:57:ASN:HD22	1:B:59:ALA:H	1.47	0.62
1:B:15:THR:OG1	1:B:17:THR:HG22	1.99	0.62
1:B:204:HIS:CD2	1:B:230:HIS:HE1	2.17	0.62
1:B:242:GLN:C	1:B:246:ARG:HA	2.14	0.62
1:A:19:LYS:HB3	3:A:367:HOH:O	2.00	0.62
1:A:8:ARG:HG3	1:A:89:LEU:O	2.00	0.61
1:A:217:VAL:HG23	1:A:220:THR:OG1	1.99	0.61
1:A:64:ALA:HB1	1:A:266:MET:CE	2.31	0.61
1:A:94:TRP:NE1	1:A:230:HIS:HD2	1.99	0.61
1:B:117:ASN:ND2	1:B:153:PHE:H	1.99	0.61
1:B:195:GLU:O	1:B:204:HIS:HB3	2.01	0.61
1:B:119:ASN:C	1:B:119:ASN:HD22	2.03	0.60
1:B:64:ALA:HB1	1:B:266:MET:HE3	1.83	0.60
1:A:119:ASN:HD22	1:A:121:ILE:H	1.48	0.60
1:B:68:TRP:H	1:B:266:MET:HE1	1.64	0.60
1:B:142:GLN:HG3	1:B:142:GLN:O	2.01	0.60
1:B:94:TRP:NE1	1:B:230:HIS:HD2	1.99	0.60
1:A:66:LEU:N	1:A:67:PRO:HD2	2.17	0.60
1:A:77:GLU:HA	1:A:80:LEU:HD12	1.83	0.60
1:B:84:ASN:C	1:B:84:ASN:HD22	2.06	0.59
1:B:199:ASN:HB2	1:B:202:GLN:NE2	2.17	0.59
1:B:176:GLN:NE2	1:B:251:GLN:HE22	2.01	0.59
1:A:64:ALA:H	1:A:260:GLN:HE22	1.49	0.59
1:B:215:MET:HE3	1:B:221:ALA:O	2.02	0.59
1:A:94:TRP:HE1	1:A:230:HIS:HD2	1.49	0.59
1:B:7:VAL:HG12	3:B:374:HOH:O	2.02	0.59
1:B:207:VAL:HG11	1:B:230:HIS:CD2	2.38	0.59
1:A:6:ARG:HE	1:A:6:ARG:HA	1.65	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:119:ASN:HD22	1:A:119:ASN:C	2.06	0.58
1:A:124:PHE:HD1	3:A:369:HOH:O	1.86	0.58
1:B:161:THR:O	1:B:206:ARG:HD3	2.03	0.58
1:A:68:TRP:N	1:A:266:MET:HE1	2.17	0.58
1:A:126:VAL:HG12	3:A:392:HOH:O	2.02	0.58
1:B:94:TRP:HE1	1:B:230:HIS:HD2	1.51	0.58
1:B:42:HIS:CE1	3:B:360:HOH:O	2.56	0.58
1:A:94:TRP:CE3	1:A:163:PRO:HG2	2.38	0.58
1:A:277:MET:HG2	3:A:382:HOH:O	2.04	0.58
1:A:94:TRP:HA	1:A:97:ASP:HB2	1.86	0.58
1:A:93:GLU:HG3	1:A:95:GLU:OE2	2.04	0.57
1:B:256:GLY:H	1:B:261:ASN:HD21	1.52	0.57
1:A:157:LYS:O	1:A:160:PRO:HD3	2.04	0.57
1:A:66:LEU:HB3	1:A:238:TRP:CE3	2.40	0.57
1:B:11:VAL:HG23	1:B:91:TYR:O	2.04	0.57
1:B:64:ALA:HA	1:B:67:PRO:HG2	1.86	0.57
1:A:218:VAL:HB	1:A:219:PRO:HD3	1.86	0.56
1:A:126:VAL:CG1	1:A:149:LEU:HB3	2.35	0.56
1:B:24:ARG:HH12	1:B:28:ILE:HD11	1.70	0.56
1:A:57:ASN:HD22	1:A:59:ALA:H	1.54	0.56
1:B:94:TRP:HE1	1:B:230:HIS:CD2	2.23	0.56
1:B:114:MET:SD	1:B:226:VAL:HG22	2.46	0.56
1:A:215:MET:HE2	1:A:221:ALA:CB	2.31	0.55
1:A:201:PRO:HG3	1:A:209:ARG:CD	2.31	0.55
1:A:104:SER:O	1:A:106:SER:N	2.38	0.55
1:A:119:ASN:HD21	1:A:121:ILE:HB	1.71	0.55
1:B:158:GLU:OE1	1:B:209:ARG:NH2	2.40	0.55
1:A:8:ARG:HD2	1:A:77:GLU:OE2	2.07	0.55
1:B:82:SER:HA	3:B:413:HOH:O	2.06	0.54
1:B:119:ASN:HD22	1:B:120:PRO:N	2.05	0.54
1:A:204:HIS:CD2	1:A:230:HIS:CE1	2.79	0.54
1:B:125:ILE:CD1	1:B:148:GLY:HA3	2.33	0.54
1:B:8:ARG:NH2	1:B:89:LEU:HD12	2.23	0.53
1:B:215:MET:CE	1:B:227:PHE:HD2	2.06	0.53
1:A:239:ALA:O	1:A:243:ILE:HG13	2.09	0.53
1:B:205:ASN:HB2	3:B:415:HOH:O	2.09	0.53
1:A:123:ASP:HA	3:A:345:HOH:O	2.09	0.53
3:A:385:HOH:O	1:B:53:GLY:HA2	2.09	0.53
1:A:61:MET:HB3	3:A:361:HOH:O	2.09	0.53
1:A:279:HIS:HD2	3:A:311:HOH:O	1.92	0.53
1:B:24:ARG:NH1	1:B:28:ILE:HD11	2.23	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:288:ILE:O	1:B:288:ILE:HG22	2.09	0.53
1:A:287:ASP:O	1:A:288:ILE:C	2.46	0.52
1:B:24:ARG:HH12	1:B:28:ILE:CD1	2.22	0.52
1:B:73:LEU:HD22	1:B:89:LEU:HD11	1.91	0.52
1:B:156:THR:HG21	1:B:209:ARG:HH12	1.73	0.52
1:A:39:ILE:HG21	1:A:139:ILE:HD11	1.90	0.52
1:A:51:PRO:HD2	1:A:54:SER:HB2	1.91	0.52
1:B:285:VAL:HB	1:B:289:GLU:HG2	1.92	0.52
1:B:156:THR:HG21	1:B:209:ARG:NH1	2.25	0.52
1:A:9:LYS:HA	1:A:287:ASP:OD2	2.09	0.51
1:A:61:MET:HG2	1:A:197:PHE:CZ	2.46	0.51
1:B:77:GLU:HA	1:B:80:LEU:HD12	1.93	0.51
1:B:255:ASN:CB	3:B:337:HOH:O	2.47	0.51
1:B:174:ILE:HG21	1:B:190:PHE:N	2.24	0.51
1:A:71:GLU:OE2	1:A:75:ARG:HD2	2.11	0.50
1:A:8:ARG:HG2	1:A:89:LEU:O	2.11	0.50
1:A:11:VAL:CG2	1:A:90:PRO:HB2	2.41	0.50
1:A:20:ARG:HD3	3:A:407:HOH:O	2.11	0.50
1:A:74:LEU:HB3	3:A:362:HOH:O	2.12	0.49
1:B:279:HIS:HE1	1:B:289:GLU:OE1	1.96	0.49
1:A:235:ASP:OD2	1:A:286:TYR:OH	2.27	0.49
1:A:68:TRP:H	1:A:266:MET:HE1	1.77	0.49
1:B:146:SER:OG	1:B:147:GLY:N	2.45	0.49
1:B:171:ALA:O	1:B:174:ILE:HG12	2.12	0.49
1:B:161:THR:CG2	1:B:206:ARG:NH1	2.75	0.49
1:A:6:ARG:N	1:A:7:VAL:CA	2.48	0.49
1:A:6:ARG:O	1:A:6:ARG:CG	2.59	0.49
1:A:89:LEU:HD11	1:A:228:PHE:HB3	1.95	0.48
1:B:217:VAL:HG23	1:B:219:PRO:HD2	1.96	0.48
1:B:240:VAL:O	1:B:243:ILE:HB	2.14	0.48
1:A:84:ASN:OD1	1:A:86:GLU:HG2	2.12	0.48
1:A:163:PRO:HD3	1:A:206:ARG:HD2	1.94	0.48
1:B:286:TYR:O	1:B:289:GLU:HB2	2.14	0.48
1:A:117:ASN:HD21	1:A:153:PHE:H	1.61	0.48
1:B:181:PRO:HG2	1:B:183:ASP:OD2	2.14	0.48
1:A:69:HIS:O	1:A:73:LEU:HG	2.14	0.48
1:B:165:ARG:HE	1:B:165:ARG:HB2	1.21	0.48
1:A:151:ARG:N	3:A:392:HOH:O	2.47	0.47
1:A:64:ALA:HB1	1:A:266:MET:HE2	1.96	0.47
1:B:253:MET:O	1:B:254:LYS:HG3	2.15	0.47
1:A:36:ASP:OD1	1:A:139:ILE:HD13	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:61:MET:O	1:B:62:SER:HB2	2.15	0.47
1:A:106:SER:OG	1:A:108:ILE:HG22	2.14	0.46
1:B:64:ALA:HB1	1:B:266:MET:HE2	1.98	0.46
1:B:57:ASN:ND2	1:B:60:HIS:HD1	2.11	0.46
1:A:113:PHE:CD1	3:A:367:HOH:O	2.56	0.45
1:B:57:ASN:ND2	1:B:60:HIS:N	2.64	0.45
1:A:117:ASN:ND2	1:A:118:GLY:H	2.13	0.45
1:A:47:LYS:HE3	1:B:142:GLN:C	2.37	0.45
1:B:119:ASN:ND2	1:B:121:ILE:H	2.14	0.45
1:B:6:ARG:HD3	1:B:81:GLN:OE1	2.16	0.45
1:B:66:LEU:HB3	1:B:238:TRP:CE3	2.52	0.45
1:A:71:GLU:OE1	1:A:271:THR:HG21	2.17	0.44
1:A:172:LEU:HD23	1:A:241:TRP:HB2	1.97	0.44
1:B:81:GLN:HE21	1:B:85:PRO:CA	2.22	0.44
1:A:94:TRP:HE1	1:A:230:HIS:CD2	2.32	0.44
1:B:114:MET:CG	1:B:226:VAL:HG22	2.47	0.44
1:B:280:ARG:HG3	1:B:285:VAL:HG12	1.99	0.44
1:A:194:LEU:HD23	1:A:194:LEU:HA	1.67	0.44
1:A:277:MET:HE1	3:A:350:HOH:O	2.17	0.44
1:B:161:THR:HG23	1:B:206:ARG:HH11	1.81	0.44
1:A:98:ALA:HB1	1:A:162:LEU:O	2.18	0.44
1:A:98:ALA:HA	1:A:162:LEU:HB2	1.99	0.44
1:B:119:ASN:HD22	1:B:120:PRO:CD	2.31	0.44
1:A:95:GLU:OE2	1:A:236:ARG:NH2	2.49	0.43
1:A:64:ALA:HB1	1:A:266:MET:HE3	1.99	0.43
1:B:253:MET:C	1:B:254:LYS:HG3	2.38	0.43
1:A:74:LEU:HD23	3:A:362:HOH:O	2.18	0.43
1:A:278:ASN:ND2	1:A:281:LYS:H	2.10	0.43
1:B:262:PHE:O	1:B:273:PRO:HD2	2.18	0.43
1:A:98:ALA:CB	1:A:163:PRO:O	2.66	0.43
1:B:66:LEU:HD21	1:B:194:LEU:HD13	1.99	0.43
1:B:235:ASP:HB3	1:B:286:TYR:OH	2.19	0.43
1:A:157:LYS:HG3	3:A:332:HOH:O	2.19	0.43
1:A:247:ASN:ND2	1:A:247:ASN:H	2.17	0.43
1:B:114:MET:HG3	1:B:226:VAL:HG22	2.01	0.43
1:B:119:ASN:C	1:B:119:ASN:ND2	2.72	0.43
1:A:47:LYS:HE3	1:B:142:GLN:O	2.18	0.42
1:A:119:ASN:O	1:A:124:PHE:N	2.38	0.42
1:A:215:MET:HE3	1:A:227:PHE:HB2	2.00	0.42
1:A:153:PHE:CE1	3:A:368:HOH:O	2.63	0.42
1:B:15:THR:HG1	1:B:17:THR:HG22	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:117:ASN:HD21	1:B:153:PHE:N	2.10	0.42
1:A:33:GLY:HA2	3:A:356:HOH:O	2.19	0.42
1:A:163:PRO:HD3	1:A:206:ARG:CD	2.50	0.42
1:A:180:PRO:HD3	1:A:182:TRP:CH2	2.54	0.42
1:A:84:ASN:HA	1:A:85:PRO:HD2	1.76	0.42
1:B:29:LEU:HD12	1:B:29:LEU:HA	1.84	0.42
1:A:139:ILE:HD13	1:A:139:ILE:HG21	1.66	0.42
1:A:151:ARG:HB2	3:A:392:HOH:O	2.20	0.42
1:A:29:LEU:HD12	1:A:29:LEU:HA	1.80	0.41
1:A:131:PHE:HD2	1:A:136:TRP:CE3	2.37	0.41
1:B:68:TRP:HB2	1:B:266:MET:HE2	2.01	0.41
1:A:224:ASP:OD1	1:A:226:VAL:HG23	2.21	0.41
1:B:42:HIS:HB3	3:B:322:HOH:O	2.20	0.41
1:B:236:ARG:O	1:B:239:ALA:HB3	2.20	0.41
1:A:161:THR:HB	3:A:377:HOH:O	2.20	0.41
1:A:38:TYR:HB3	1:A:72:TYR:OH	2.21	0.41
1:A:194:LEU:HD13	1:A:234:VAL:HG13	2.03	0.41
1:B:194:LEU:HA	1:B:203:LEU:HD13	2.02	0.41
1:B:252:PRO:HG2	1:B:261:ASN:HA	2.03	0.41
1:A:221:ALA:HB3	1:A:222:PRO:HD3	2.03	0.41
1:A:262:PHE:HD2	1:A:263:ARG:HE	1.68	0.41
1:B:4:LYS:HD2	1:B:4:LYS:N	2.36	0.41
1:B:126:VAL:HG13	1:B:149:LEU:HB3	2.03	0.41
1:B:126:VAL:HG11	1:B:149:LEU:HD13	2.03	0.41
1:A:64:ALA:O	1:A:266:MET:HE1	2.21	0.40
1:A:119:ASN:ND2	1:A:121:ILE:HB	2.34	0.40
1:B:57:ASN:ND2	1:B:60:HIS:H	2.19	0.40
1:B:194:LEU:HA	1:B:203:LEU:CD1	2.51	0.40
1:B:242:GLN:NE2	1:B:277:MET:HE3	2.36	0.40
1:A:127:ASP:OD1	1:A:127:ASP:N	2.51	0.40
1:A:218:VAL:HG12	1:A:219:PRO:N	2.36	0.40
1:A:45:ALA:O	1:A:57:ASN:HB2	2.21	0.40
1:A:69:HIS:CD2	1:A:231:HIS:CE1	3.10	0.40
1:B:117:ASN:ND2	1:B:118:GLY:H	2.19	0.40
1:B:124:PHE:HB3	1:B:151:ARG:O	2.21	0.40

All (2) 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)
1:B:137:THR:CG2	1:B:246:ARG:NH1[1_455]	2.05	0.15

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:104:SER:OG	1:A:246:ARG:CB[2_545]	2.18	0.02

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	275/303 (91%)	263 (96%)	10 (4%)	2 (1%)	22 26
1	B	280/303 (92%)	267 (95%)	11 (4%)	2 (1%)	22 26
All	All	555/606 (92%)	530 (96%)	21 (4%)	4 (1%)	22 26

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	255	ASN
1	B	256	GLY
1	A	287	ASP
1	A	244	ILE

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	242/266 (91%)	224 (93%)	18 (7%)	13 17
1	B	247/266 (93%)	225 (91%)	22 (9%)	9 11
All	All	489/532 (92%)	449 (92%)	40 (8%)	11 14

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	ARG
1	A	16	ASP
1	A	17	THR
1	A	20	ARG
1	A	31	GLU
1	A	81	GLN
1	A	97	ASP
1	A	107	GLN
1	A	117	ASN
1	A	119	ASN
1	A	126	VAL
1	A	149	LEU
1	A	152	ASN
1	A	194	LEU
1	A	217	VAL
1	A	229	LEU
1	A	246	ARG
1	A	278	ASN
1	B	20	ARG
1	B	29	LEU
1	B	31	GLU
1	B	55	ASP
1	B	75	ARG
1	B	84	ASN
1	B	94	TRP
1	B	117	ASN
1	B	119	ASN
1	B	126	VAL
1	B	149	LEU
1	B	152	ASN
1	B	163	PRO
1	B	165	ARG
1	B	184	MET
1	B	187	GLN
1	B	194	LEU
1	B	197	PHE
1	B	217	VAL
1	B	229	LEU
1	B	243	ILE
1	B	281	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	57	ASN
1	A	117	ASN
1	A	119	ASN
1	A	142	GLN
1	A	152	ASN
1	A	202	GLN
1	A	204	HIS
1	A	205	ASN
1	A	230	HIS
1	A	247	ASN
1	A	260	GLN
1	A	261	ASN
1	A	270	ASN
1	A	278	ASN
1	A	279	HIS
1	B	57	ASN
1	B	81	GLN
1	B	84	ASN
1	B	117	ASN
1	B	119	ASN
1	B	152	ASN
1	B	176	GLN
1	B	187	GLN
1	B	204	HIS
1	B	230	HIS
1	B	260	GLN
1	B	261	ASN
1	B	279	HIS

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 monosaccharides in this entry.

5.6 Ligand geometry

Of 3 ligands modelled in this entry, 3 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

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, 95^{th} 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	279/303 (92%)	-0.02	4 (1%)	75	80	10, 18, 39, 61	0
1	B	284/303 (93%)	-0.10	2 (0%)	87	91	10, 18, 32, 60	0
All	All	563/606 (92%)	-0.06	6 (1%)	80	85	10, 18, 35, 61	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	98	ALA	6.0
1	A	105	GLN	4.0
1	A	97	ASP	2.8
1	A	103	PRO	2.8
1	B	4	LYS	2.7
1	B	121	ILE	2.6

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 monosaccharides in this entry.

6.4 Ligands

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, 95^{th} 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	0.50	0.59	30,30,30,30	1
2	CU	B	501	1/1	0.84	0.14	20,20,20,20	0
2	CU	A	501	1/1	0.96	0.10	19,19,19,19	0

6.5 Other polymers (i)

There are no such residues in this entry.