



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:15 PM GMT

PDB ID : 4E9V
Title : Multicopper Oxidase mgLAC (data1)
Authors : Komori, H.; Miyazaki, K.; Higuchi, Y.
Deposited on : 2012-03-21
Resolution : 1.80 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
<http://wwpdb.org/validation/2016/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
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

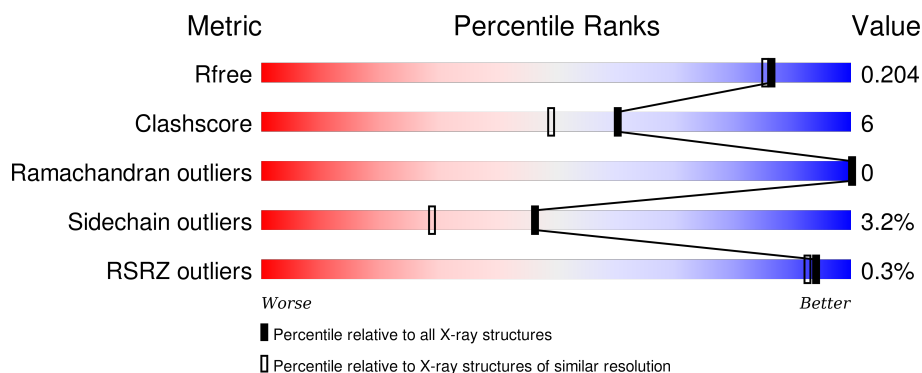
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



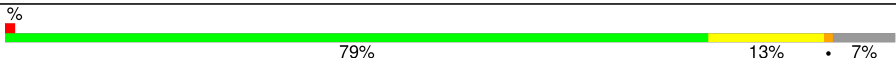
The reported resolution of this entry is 1.80 Å.

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}	91344	4533 (1.80-1.80)
Clashscore	102246	5383 (1.80-1.80)
Ramachandran outliers	100387	5320 (1.80-1.80)
Sidechain outliers	100360	5319 (1.80-1.80)
RSRZ outliers	91569	4547 (1.80-1.80)

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. 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	339	 81% 12% • 6%
1	B	339	 82% 11% 7%
1	C	339	 % 79% 13% • 7%

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
3	O	A	1403	-	-	-	X
3	O	B	2406	-	-	-	X
3	O	C	3406	-	-	-	X
4	OH	A	1406	-	-	X	X
4	OH	B	2403	-	-	-	X
4	OH	C	3403	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 8364 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 Multicopper oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	317	Total	C	N	O	S	0	7	0
			2539	1633	425	469	12			
1	B	316	Total	C	N	O	S	0	5	0
			2519	1617	423	468	11			
1	C	316	Total	C	N	O	S	0	7	0
			2533	1631	424	467	11			

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1001	MET	-	EXPRESSION TAG	UNP C0STU6
A	1327	LYS	-	EXPRESSION TAG	UNP C0STU6
A	1328	LEU	-	EXPRESSION TAG	UNP C0STU6
A	1329	ALA	-	EXPRESSION TAG	UNP C0STU6
A	1330	ALA	-	EXPRESSION TAG	UNP C0STU6
A	1331	ALA	-	EXPRESSION TAG	UNP C0STU6
A	1332	LEU	-	EXPRESSION TAG	UNP C0STU6
A	1333	GLU	-	EXPRESSION TAG	UNP C0STU6
A	1334	HIS	-	EXPRESSION TAG	UNP C0STU6
A	1335	HIS	-	EXPRESSION TAG	UNP C0STU6
A	1336	HIS	-	EXPRESSION TAG	UNP C0STU6
A	1337	HIS	-	EXPRESSION TAG	UNP C0STU6
A	1338	HIS	-	EXPRESSION TAG	UNP C0STU6
A	1339	HIS	-	EXPRESSION TAG	UNP C0STU6
B	2001	MET	-	EXPRESSION TAG	UNP C0STU6
B	2327	LYS	-	EXPRESSION TAG	UNP C0STU6
B	2328	LEU	-	EXPRESSION TAG	UNP C0STU6
B	2329	ALA	-	EXPRESSION TAG	UNP C0STU6
B	2330	ALA	-	EXPRESSION TAG	UNP C0STU6
B	2331	ALA	-	EXPRESSION TAG	UNP C0STU6
B	2332	LEU	-	EXPRESSION TAG	UNP C0STU6
B	2333	GLU	-	EXPRESSION TAG	UNP C0STU6
B	2334	HIS	-	EXPRESSION TAG	UNP C0STU6

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Chain	Residue	Modelled	Actual	Comment	Reference
B	2335	HIS	-	EXPRESSION TAG	UNP C0STU6
B	2336	HIS	-	EXPRESSION TAG	UNP C0STU6
B	2337	HIS	-	EXPRESSION TAG	UNP C0STU6
B	2338	HIS	-	EXPRESSION TAG	UNP C0STU6
B	2339	HIS	-	EXPRESSION TAG	UNP C0STU6
C	3001	MET	-	EXPRESSION TAG	UNP C0STU6
C	3327	LYS	-	EXPRESSION TAG	UNP C0STU6
C	3328	LEU	-	EXPRESSION TAG	UNP C0STU6
C	3329	ALA	-	EXPRESSION TAG	UNP C0STU6
C	3330	ALA	-	EXPRESSION TAG	UNP C0STU6
C	3331	ALA	-	EXPRESSION TAG	UNP C0STU6
C	3332	LEU	-	EXPRESSION TAG	UNP C0STU6
C	3333	GLU	-	EXPRESSION TAG	UNP C0STU6
C	3334	HIS	-	EXPRESSION TAG	UNP C0STU6
C	3335	HIS	-	EXPRESSION TAG	UNP C0STU6
C	3336	HIS	-	EXPRESSION TAG	UNP C0STU6
C	3337	HIS	-	EXPRESSION TAG	UNP C0STU6
C	3338	HIS	-	EXPRESSION TAG	UNP C0STU6
C	3339	HIS	-	EXPRESSION TAG	UNP C0STU6

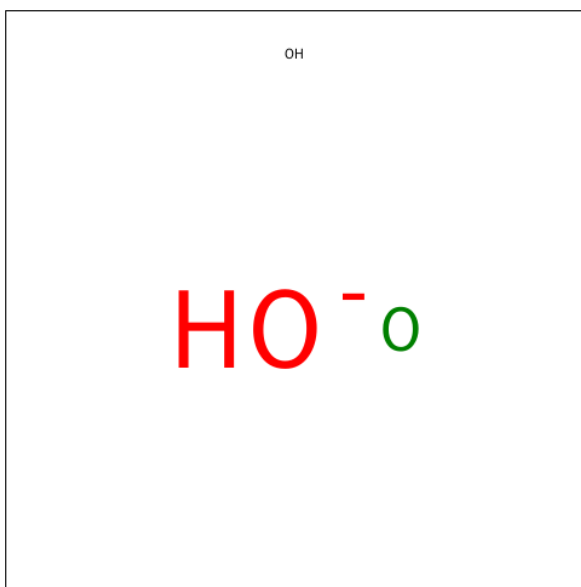
- Molecule 2 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	4	Total Cu 4 4	0	0
2	A	4	Total Cu 4 4	0	0
2	C	4	Total Cu 4 4	0	0

- Molecule 3 is OXYGEN ATOM (three-letter code: O) (formula: O).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total O 1 1	0	0
3	A	1	Total O 1 1	0	0
3	C	1	Total O 1 1	0	0

- Molecule 4 is HYDROXIDE ION (three-letter code: OH) (formula: HO).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total O 1 1	0	0
4	B	1	Total O 1 1	0	0
4	C	1	Total O 1 1	0	0

- Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total Cl 1 1	0	0

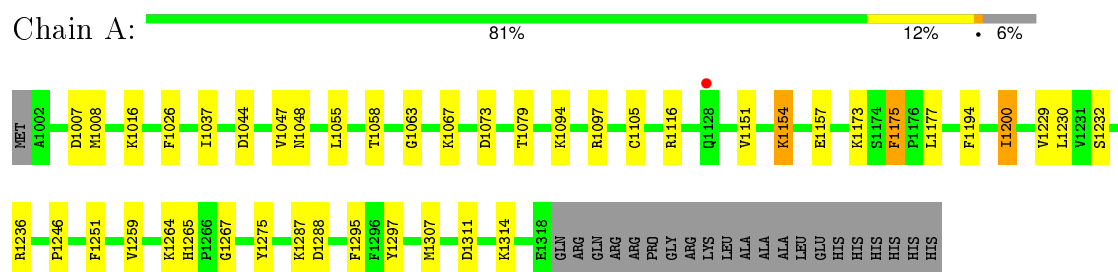
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	234	Total O 235 235	0	1
6	B	265	Total O 265 265	0	0
6	C	254	Total O 254 254	0	0

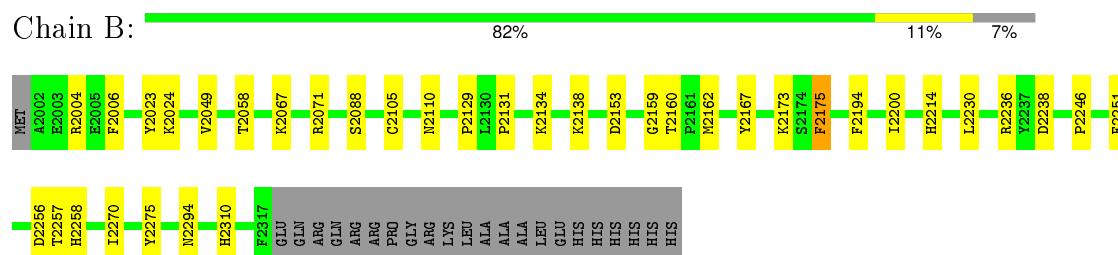
3 Residue-property plots [i](#)

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

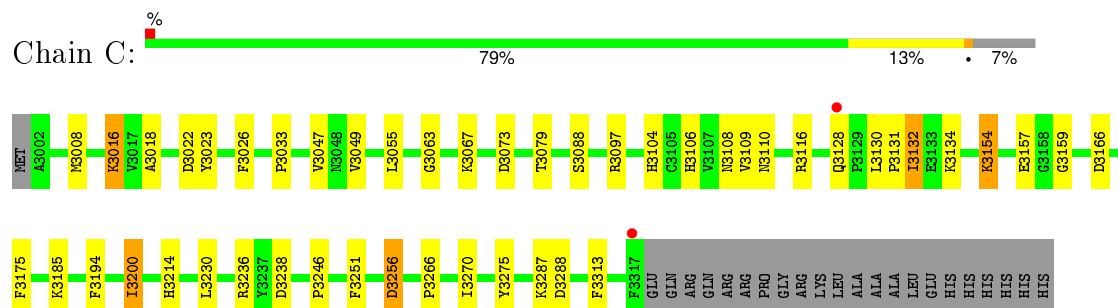
• Molecule 1: Multicopper oxidase



• Molecule 1: Multicopper oxidase



• Molecule 1: Multicopper oxidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	74.73Å 101.13Å 123.96Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 1.80 22.26 – 1.80	Depositor EDS
% Data completeness (in resolution range)	95.0 (10.00-1.80) 100.0 (22.26-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.17 (at 1.80Å)	Xtriage
Refinement program	SHELXL-97	Depositor
R, R_{free}	0.163 , (Not available) 0.155 , 0.204	Depositor DCC
R_{free} test set	4355 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	10.4	Xtriage
Anisotropy	0.415	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 67.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 87486 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8364	wwPDB-VP
Average B, all atoms (Å ²)	13.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.40% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: CL, O, CU, OH

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	0.37	0/2641	0.99	3/3594 (0.1%)
1	B	0.37	0/2615	0.99	4/3561 (0.1%)
1	C	0.37	0/2637	0.98	3/3590 (0.1%)
All	All	0.37	0/7893	0.99	10/10745 (0.1%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1116	ARG	NE-CZ-NH1	7.26	123.93	120.30
1	B	2004	ARG	NE-CZ-NH1	-6.45	117.07	120.30
1	C	3251	PHE	CB-CG-CD2	5.74	124.81	120.80
1	A	1116	ARG	NE-CZ-NH2	-5.65	117.48	120.30
1	B	2167	TYR	CB-CG-CD1	-5.58	117.65	121.00
1	A	1175	PHE	CB-CG-CD1	-5.52	116.93	120.80
1	C	3116	ARG	CD-NE-CZ	5.35	131.09	123.60
1	B	2071	ARG	NE-CZ-NH1	-5.30	117.65	120.30
1	B	2175	PHE	CB-CG-CD1	-5.22	117.14	120.80
1	C	3166	ASP	CB-CG-OD1	5.02	122.82	118.30

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	2539	0	2439	30	0
1	B	2519	0	2404	20	0
1	C	2533	0	2429	33	0
2	A	4	0	0	0	0
2	B	4	0	0	0	0
2	C	4	0	0	0	0
3	A	1	0	0	1	0
3	B	1	0	0	1	0
3	C	1	0	0	1	0
4	A	1	0	0	3	0
4	B	1	0	0	1	0
4	C	1	0	0	1	0
5	A	1	0	0	0	0
6	A	235	0	0	2	0
6	B	265	0	0	4	0
6	C	254	0	0	0	0
All	All	8364	0	7272	83	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 6.

All (83) 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:1311:ASP:HA	1:A:1314:LYS:HD2	1.66	0.78
1:C:3016:LYS:HE2	1:C:3018:ALA:O	1.84	0.77
1:B:2173:LYS:HE2	6:B:2654:HOH:O	1.85	0.75
1:A:1287:LYS:HE3	1:A:1288:ASP:OD2	1.87	0.74
1:A:1173[B]:LYS:HD3	1:A:1177:LEU:O	1.88	0.73
1:A:1151:VAL:HA	1:A:1154:LYS:HD3	1.69	0.73
1:C:3016:LYS:HE3	1:C:3022:ASP:OD1	1.89	0.71
1:C:3132:ILE:H	1:C:3132:ILE:HD13	1.57	0.69
1:A:1016:LYS:HE3	6:A:5165:HOH:O	1.93	0.69
1:C:3131:PRO:O	1:C:3134:LYS:HB2	1.94	0.68
1:A:1173[B]:LYS:HD2	1:A:1295:PHE:HE1	1.60	0.66
1:C:3200[A]:ILE:HG22	1:C:3230:LEU:HD11	1.78	0.66
3:B:2406:O:O	4:C:3403:OH:O	2.16	0.64
1:A:1154:LYS:HE2	1:A:1157:GLU:OE1	2.00	0.60
4:A:1406:OH:O	3:C:3406:O:O	2.18	0.60
1:B:2160:THR:HB	1:B:2162:MET:HE1	1.81	0.60
1:A:1008[A]:MET:HE2	1:A:1307:MET:HE1	1.84	0.60
3:A:1403:O:O	4:B:2403:OH:O	2.19	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2246:PRO:HA	1:B:2275:TYR:CG	2.40	0.57
1:A:1173[B]:LYS:HD2	1:A:1295:PHE:CE1	2.38	0.57
1:C:3130:LEU:HB3	1:C:3132:ILE:HG12	1.88	0.56
1:C:3026:PHE:HZ	1:C:3055:LEU:HD12	1.71	0.56
1:C:3185[B]:LYS:HD2	1:C:3246:PRO:CG	2.35	0.56
1:B:2129:PRO:HG2	1:B:2134:LYS:HE2	1.88	0.55
1:A:1063:GLY:O	1:A:1097:ARG:HG2	2.08	0.54
1:A:1246:PRO:HA	1:A:1275:TYR:CG	2.43	0.54
1:A:1044:ASP:OD1	1:A:1094:LYS:HE2	2.08	0.54
1:C:3026:PHE:CZ	1:C:3055:LEU:HD12	2.43	0.53
1:A:1232[A]:SER:OG	1:B:2230:LEU:HD13	2.08	0.53
1:C:3246:PRO:HA	1:C:3275:TYR:CG	2.45	0.52
1:B:2049:VAL:O	1:B:2088:SER:HA	2.10	0.52
1:B:2214:HIS:HB2	1:B:2238:ASP:HB2	1.92	0.51
1:C:3008[B]:MET:HG3	1:C:3047:VAL:HG13	1.91	0.51
1:B:2024:LYS:HE2	1:B:2153:ASP:OD1	2.11	0.51
1:A:1311:ASP:HA	1:A:1314:LYS:CD	2.40	0.50
1:C:3008[B]:MET:HG3	1:C:3047:VAL:CG1	2.42	0.50
1:C:3214:HIS:HB2	1:C:3238:ASP:HB2	1.92	0.50
1:C:3106:HIS:HA	1:C:3109:VAL:HG12	1.94	0.49
1:A:1007:ASP:OD1	1:A:1048:ASN:HB2	2.12	0.49
4:A:1406:OH:O	1:C:3104:HIS:CE1	2.66	0.49
1:B:2200[A]:ILE:HG22	1:B:2230:LEU:HD11	1.95	0.48
1:A:1265:HIS:CD2	1:C:3108:ASN:HA	2.48	0.48
1:A:1200[A]:ILE:HG22	1:A:1230:LEU:HD11	1.96	0.48
1:C:3049:VAL:O	1:C:3088:SER:HA	2.14	0.48
1:C:3130:LEU:CB	1:C:3132:ILE:HG12	2.44	0.47
1:B:2138:LYS:HB2	1:B:2138:LYS:HE2	1.57	0.47
1:B:2110:ASN:HB2	1:B:2159:GLY:O	2.16	0.46
1:A:1200[A]:ILE:CG2	1:A:1230:LEU:HD11	2.46	0.46
1:C:3154:LYS:HE2	1:C:3157:GLU:OE1	2.16	0.46
1:C:3063:GLY:O	1:C:3097:ARG:HG2	2.16	0.45
1:B:2162:MET:HB2	1:B:2162:MET:HE2	1.59	0.45
1:C:3131:PRO:HA	1:C:3134:LYS:HD2	1.98	0.45
1:B:2294:ASN:HB3	1:B:2310:HIS:CE1	2.52	0.44
1:B:2246:PRO:HA	1:B:2275:TYR:CD2	2.52	0.44
1:C:3185[B]:LYS:HD2	1:C:3246:PRO:HG3	1.99	0.44
1:A:1259:VAL:HG21	1:C:3109:VAL:CG2	2.47	0.44
4:A:1406:OH:O	1:C:3104:HIS:HE1	2.01	0.43
1:A:1175:PHE:O	1:A:1267:GLY:HA2	2.18	0.43
1:A:1008[A]:MET:HE2	1:A:1307:MET:CE	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1037:ILE:HG12	1:A:1307:MET:HE3	2.01	0.43
1:A:1200[A]:ILE:HD13	1:A:1232[A]:SER:HB2	2.00	0.42
1:A:1026:PHE:CZ	1:A:1055:LEU:HD12	2.54	0.42
1:A:1287:LYS:O	1:A:1288:ASP:HB2	2.20	0.42
1:A:1200[A]:ILE:HD13	1:A:1232[A]:SER:CB	2.50	0.42
1:A:1058:THR:O	1:A:1105:CYS:HA	2.20	0.42
6:B:2726:HOH:O	1:C:3256:ASP:HB3	2.20	0.41
1:C:3110:ASN:HB2	1:C:3159:GLY:O	2.20	0.41
1:B:2257:THR:HG23	6:B:2585:HOH:O	2.20	0.41
1:C:3073:ASP:HB3	1:C:3079:THR:HG21	2.03	0.41
1:A:1073:ASP:HB3	1:A:1079:THR:HG21	2.03	0.41
1:A:1229:VAL:HG13	6:A:5221:HOH:O	2.20	0.41
1:C:3033:PRO:HD2	1:C:3313:PHE:CG	2.56	0.41
1:C:3175:PHE:HB2	1:C:3270:ILE:HB	2.03	0.41
1:C:3200[A]:ILE:CG2	1:C:3230:LEU:HD11	2.50	0.40
1:C:3008[B]:MET:HE3	1:C:3008[B]:MET:HB3	1.92	0.40
1:A:1008[A]:MET:HE3	1:A:1047:VAL:HG11	2.03	0.40
1:C:3154:LYS:HE2	1:C:3157:GLU:CD	2.40	0.40
1:B:2131:PRO:O	1:B:2134:LYS:HB2	2.21	0.40
1:C:3287:LYS:O	1:C:3288:ASP:HB2	2.22	0.40
1:B:2175:PHE:HB2	1:B:2270:ILE:HB	2.04	0.40
1:B:2058:THR:O	1:B:2105:CYS:HA	2.21	0.40
1:B:2138:LYS:HD3	6:B:2730:HOH:O	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	322/339 (95%)	318 (99%)	4 (1%)	0	100	100
1	B	319/339 (94%)	312 (98%)	7 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	321/339 (95%)	313 (98%)	8 (2%)	0	100	100
All	All	962/1017 (95%)	943 (98%)	19 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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	276/287 (96%)	267 (97%)	9 (3%)	45	27
1	B	273/287 (95%)	266 (97%)	7 (3%)	54	37
1	C	275/287 (96%)	263 (96%)	12 (4%)	35	17
All	All	824/861 (96%)	796 (97%)	28 (3%)	46	26

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1067	LYS
1	A	1154	LYS
1	A	1194	PHE
1	A	1200[A]	ILE
1	A	1200[B]	ILE
1	A	1236	ARG
1	A	1251	PHE
1	A	1264	LYS
1	A	1297	TYR
1	B	2006	PHE
1	B	2023	TYR
1	B	2067	LYS
1	B	2194	PHE
1	B	2236	ARG
1	B	2251	PHE
1	B	2256	ASP
1	C	3016	LYS

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Mol	Chain	Res	Type
1	C	3023	TYR
1	C	3067	LYS
1	C	3128	GLN
1	C	3132	ILE
1	C	3154	LYS
1	C	3194	PHE
1	C	3200[A]	ILE
1	C	3200[B]	ILE
1	C	3236	ARG
1	C	3256	ASP
1	C	3266	PRO

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

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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 19 ligands modelled in this entry, 16 are monoatomic and 3 are modelled with single atom - 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	317/339 (93%)	-0.38	1 (0%) 94 92	5, 11, 26, 48	0
1	B	316/339 (93%)	-0.45	0 100 100	5, 11, 24, 45	0
1	C	316/339 (93%)	-0.39	2 (0%) 90 88	5, 11, 28, 69	0
All	All	949/1017 (93%)	-0.41	3 (0%) 94 92	5, 11, 26, 69	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1128	GLN	2.6
1	C	3317	PHE	2.3
1	C	3128	GLN	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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
4	OH	A	1406	1/1	0.86	0.27	28.50	19,19,19,19	1
4	OH	B	2403	1/1	0.84	0.27	20.12	18,18,18,18	1
4	OH	C	3403	1/1	0.79	0.23	10.00	14,14,14,14	1
3	O	A	1403	1/1	0.92	0.16	9.24	8,8,8,8	1
3	O	B	2406	1/1	0.85	0.19	7.39	6,6,6,6	1
3	O	C	3406	1/1	0.96	0.09	2.04	7,7,7,7	1
2	CU	C	3401	1/1	0.98	0.08	-0.31	16,16,16,16	1
2	CU	A	1402	1/1	0.99	0.07	-0.37	16,16,16,16	1
2	CU	A	1404	1/1	0.98	0.06	-2.21	18,18,18,18	1
5	CL	A	1407	1/1	1.00	0.04	-2.36	12,12,12,12	0
2	CU	B	2404	1/1	1.00	0.05	-2.58	9,9,9,9	0
2	CU	B	2401	1/1	0.99	0.05	-2.61	19,19,19,19	1
2	CU	A	1401	1/1	0.99	0.06	-2.96	11,11,11,11	0
2	CU	C	3402	1/1	0.99	0.04	-3.17	14,14,14,14	1
2	CU	B	2402	1/1	0.99	0.06	-3.47	18,18,18,18	1
2	CU	B	2405	1/1	0.99	0.04	-3.79	14,14,14,14	1
2	CU	C	3404	1/1	0.99	0.04	-5.12	10,10,10,10	0
2	CU	A	1405	1/1	1.00	0.03	-5.73	17,17,17,17	1
2	CU	C	3405	1/1	0.99	0.03	-6.03	14,14,14,14	1

6.5 Other polymers [i](#)

There are no such residues in this entry.