



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:12 PM UTC

PDB ID : 5IKX / pdb_00005ikx
Title : Crystal structure of the alpha-esterase-7 carboxyl esterase (dimer), E3, from *Lucilia cuprina*
Authors : Jackson, C.; Fraser, N.
Deposited on : 2016-03-04
Resolution : 2.19 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

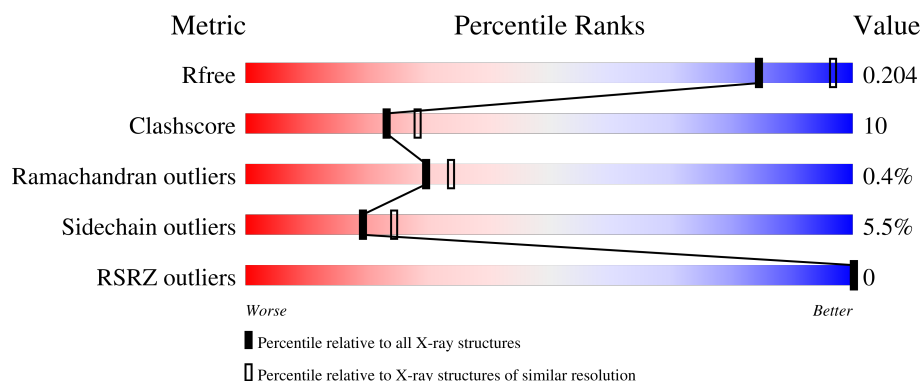
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.19 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	577	 79% 17% ..
1	B	577	 76% 21% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9516 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 Carboxylic ester hydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	569	Total	C	N	O	S	0	0	0
			4577	2923	770	850	34			
1	B	569	Total	C	N	O	S	0	0	0
			4563	2916	767	847	33			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	MET	-	initiating methionine	UNP Q25252
A	-5	HIS	-	expression tag	UNP Q25252
A	-4	HIS	-	expression tag	UNP Q25252
A	-3	HIS	-	expression tag	UNP Q25252
A	-2	HIS	-	expression tag	UNP Q25252
A	-1	HIS	-	expression tag	UNP Q25252
A	0	HIS	-	expression tag	UNP Q25252
A	83	ALA	ASP	conflict	UNP Q25252
A	364	LEU	MET	conflict	UNP Q25252
A	419	PHE	ILE	conflict	UNP Q25252
A	472	THR	ALA	conflict	UNP Q25252
A	505	THR	ILE	conflict	UNP Q25252
A	530	GLU	LYS	conflict	UNP Q25252
A	554	GLY	ASP	conflict	UNP Q25252
B	-6	MET	-	initiating methionine	UNP Q25252
B	-5	HIS	-	expression tag	UNP Q25252
B	-4	HIS	-	expression tag	UNP Q25252
B	-3	HIS	-	expression tag	UNP Q25252
B	-2	HIS	-	expression tag	UNP Q25252
B	-1	HIS	-	expression tag	UNP Q25252
B	0	HIS	-	expression tag	UNP Q25252
B	83	ALA	ASP	conflict	UNP Q25252
B	364	LEU	MET	conflict	UNP Q25252
B	419	PHE	ILE	conflict	UNP Q25252
B	472	THR	ALA	conflict	UNP Q25252

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Chain	Residue	Modelled	Actual	Comment	Reference
B	505	THR	ILE	conflict	UNP Q25252
B	530	GLU	LYS	conflict	UNP Q25252
B	554	GLY	ASP	conflict	UNP Q25252

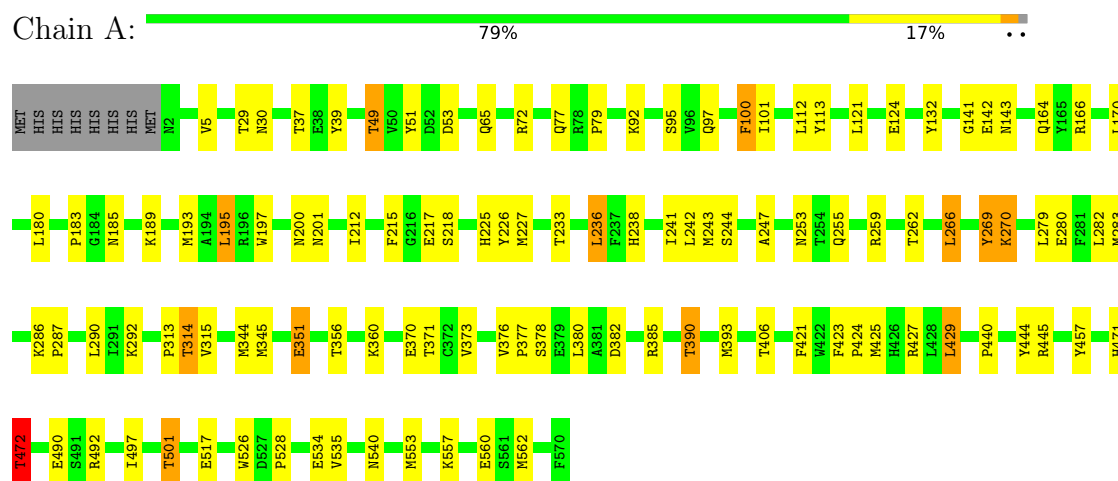
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	200	Total 200	O 200	0	0
2	B	176	Total 176	O 176	0	0

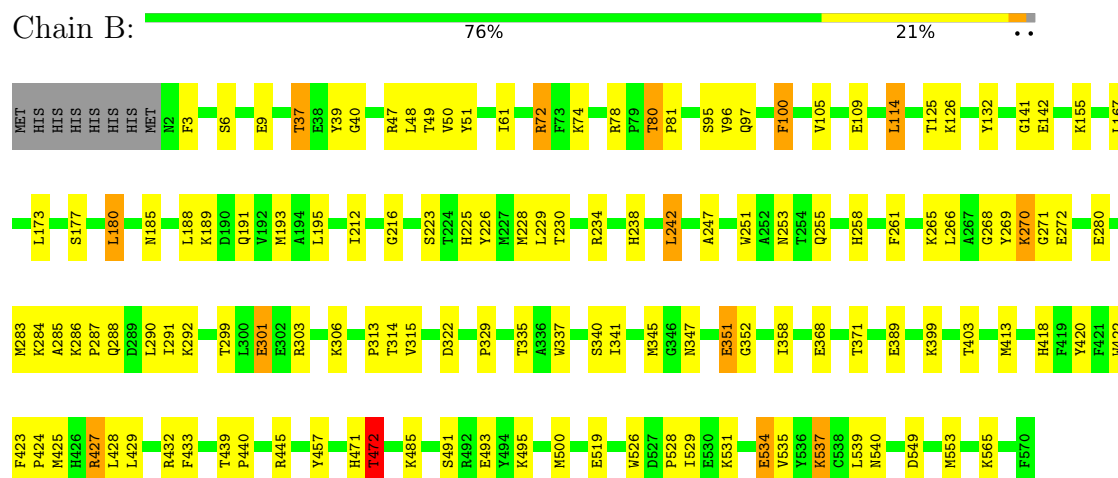
3 Residue-property plots

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: Carboxylic ester hydrolase



• Molecule 1: Carboxylic ester hydrolase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	61.77Å 108.96Å 92.17Å 90.00° 90.36° 90.00°	Depositor
Resolution (Å)	47.06 – 2.19 47.06 – 2.19	Depositor EDS
% Data completeness (in resolution range)	97.9 (47.06-2.19) 98.0 (47.06-2.19)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.05 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.171 , 0.209 0.170 , 0.204	Depositor DCC
R_{free} test set	3145 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	33.8	Xtriage
Anisotropy	0.438	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 17.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	0.420 for h,-k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	9516	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.99% 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

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	0.63	0/4693	0.85	3/6351 (0.0%)
1	B	0.64	0/4679	0.85	0/6334
All	All	0.64	0/9372	0.85	3/12685 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	269	TYR	CA-C-N	5.44	131.50	121.70
1	A	269	TYR	C-N-CA	5.44	131.50	121.70
1	A	236	LEU	N-CA-C	5.09	117.50	111.33

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	4577	0	4483	88	0
1	B	4563	0	4462	99	0
2	A	200	0	0	11	0
2	B	176	0	0	10	0
All	All	9516	0	8945	187	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 10.

All (187) 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:269:TYR:HA	1:A:270:LYS:HB2	1.16	1.09
1:B:427:ARG:HH11	1:B:427:ARG:HG2	1.19	1.03
1:A:351:GLU:HG3	1:A:471:HIS:HB2	1.42	0.99
1:B:286:LYS:HD2	1:B:287:PRO:HD2	1.46	0.95
1:B:78:ARG:HA	1:B:193:MET:HE1	1.48	0.94
1:B:471:HIS:O	1:B:472:THR:HB	1.66	0.93
1:A:471:HIS:O	1:A:472:THR:HB	1.67	0.92
1:B:72:ARG:HD2	1:B:109:GLU:OE2	1.70	0.91
1:A:49:THR:HG21	2:A:698:HOH:O	1.69	0.91
1:A:501:THR:HG21	2:A:633:HOH:O	1.73	0.88
1:A:269:TYR:HA	1:A:270:LYS:CB	2.03	0.88
1:A:185:ASN:HD21	1:A:314:THR:H	1.22	0.87
1:A:457:TYR:CB	1:A:472:THR:HG21	2.10	0.82
1:A:393:MET:HB3	1:A:562:MET:HE3	1.65	0.77
1:B:286:LYS:CD	1:B:287:PRO:HD2	2.15	0.77
1:B:72:ARG:NH2	1:B:285:ALA:O	2.20	0.75
1:B:37:THR:HG23	1:B:39:TYR:H	1.51	0.73
1:A:283:MET:HG3	2:A:632:HOH:O	1.88	0.73
1:B:427:ARG:HG2	1:B:427:ARG:NH1	1.98	0.73
1:B:225:HIS:HD2	1:B:247:ALA:H	1.36	0.73
1:A:344:MET:HE3	1:A:526:TRP:HH2	1.55	0.72
1:B:212:ILE:H	1:B:238:HIS:HD2	1.36	0.72
1:B:96:VAL:HG12	1:B:291:ILE:HG13	1.71	0.71
1:A:344:MET:HE3	1:A:526:TRP:CH2	2.25	0.70
1:B:491:SER:O	1:B:495:LYS:HG3	1.91	0.70
1:A:269:TYR:CA	1:A:270:LYS:HB2	2.08	0.70
1:B:95:SER:HB3	1:B:167:LEU:HD12	1.71	0.70
1:B:439:THR:HA	2:B:768:HOH:O	1.93	0.69
1:B:427:ARG:HH11	1:B:427:ARG:CG	2.00	0.69
1:B:185:ASN:HD21	1:B:314:THR:H	1.41	0.69
1:B:97:GLN:HE22	1:B:141:GLY:H	1.42	0.68
1:A:255:GLN:HE22	1:A:315:VAL:H	1.40	0.67
1:B:457:TYR:HB3	1:B:472:THR:HG21	1.76	0.67
1:B:49:THR:HG22	1:B:51:TYR:H	1.60	0.67
1:A:253:ASN:HB2	2:A:723:HOH:O	1.94	0.66
1:A:97:GLN:NE2	1:A:141:GLY:H	1.93	0.66
1:B:78:ARG:CA	1:B:193:MET:HE1	2.25	0.65
1:A:100:PHE:HD2	1:A:100:PHE:H	1.44	0.65
1:B:286:LYS:HD2	1:B:287:PRO:CD	2.25	0.65
1:A:262:THR:O	1:A:266:LEU:HD22	1.97	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:286:LYS:HE3	1:B:288:GLN:H	1.63	0.64
1:A:373:VAL:CG1	1:A:385:ARG:HB3	2.28	0.63
1:B:37:THR:HG22	1:B:40:GLY:O	1.99	0.63
1:B:534:GLU:CD	1:B:534:GLU:H	2.07	0.63
1:A:377:PRO:HD2	1:A:562:MET:HE1	1.80	0.63
1:B:457:TYR:CB	1:B:472:THR:HG21	2.29	0.62
1:B:535:VAL:HG13	1:B:553:MET:HG2	1.82	0.62
1:B:61:ILE:HB	1:B:114:LEU:HD23	1.82	0.62
1:A:382:ASP:OD2	1:A:390:THR:HG21	1.99	0.62
1:B:189:LYS:HE3	1:B:322:ASP:O	2.00	0.62
1:A:457:TYR:HB3	1:A:472:THR:HG21	1.82	0.62
1:A:65:GLN:HE21	1:A:79:PRO:HA	1.61	0.62
1:B:337:TRP:O	1:B:340:SER:OG	2.18	0.61
1:A:457:TYR:HB2	1:A:472:THR:HG21	1.82	0.61
1:B:96:VAL:CG1	1:B:291:ILE:HG13	2.31	0.61
1:B:212:ILE:H	1:B:238:HIS:CD2	2.17	0.61
1:B:272:GLU:HB3	2:B:656:HOH:O	2.00	0.61
1:A:212:ILE:H	1:A:238:HIS:HD2	1.46	0.61
1:B:74:LYS:NZ	1:B:283:MET:HE1	2.16	0.60
1:B:418:HIS:HE1	2:B:691:HOH:O	1.84	0.60
1:B:471:HIS:O	1:B:472:THR:CB	2.48	0.60
1:A:492:ARG:HD3	2:A:626:HOH:O	2.01	0.60
1:B:440:PRO:HB2	1:B:528:PRO:HB2	1.84	0.59
1:A:425:MET:O	1:A:429:LEU:HD22	2.03	0.59
1:B:155:LYS:HE3	2:B:771:HOH:O	2.03	0.59
1:B:253:ASN:ND2	1:B:329:PRO:HG2	2.18	0.59
1:B:49:THR:HG21	2:B:638:HOH:O	2.03	0.58
1:A:283:MET:CG	2:A:632:HOH:O	2.49	0.58
1:B:537:LYS:HA	1:B:549:ASP:HA	1.84	0.58
1:B:269:TYR:HA	1:B:270:LYS:HB2	1.84	0.58
1:A:472:THR:HG22	2:A:614:HOH:O	2.03	0.58
1:B:286:LYS:CE	1:B:287:PRO:HD2	2.34	0.57
1:A:286:LYS:HD2	1:A:287:PRO:HD2	1.86	0.56
1:A:143:ASN:H	1:A:164:GLN:HE21	1.54	0.56
1:A:225:HIS:HD2	1:A:247:ALA:H	1.53	0.56
1:A:185:ASN:ND2	1:A:313:PRO:HA	2.20	0.56
1:B:191:GLN:HE22	1:B:223:SER:HB3	1.71	0.56
1:B:49:THR:HG22	1:B:50:VAL:N	2.21	0.56
1:B:433:PHE:HB3	1:B:531:LYS:HG3	1.88	0.55
1:B:47:ARG:HD3	2:B:697:HOH:O	2.06	0.55
1:B:255:GLN:HE22	1:B:315:VAL:H	1.54	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:261:PHE:CE2	1:B:265:LYS:HD2	2.41	0.55
1:A:49:THR:HG22	1:A:51:TYR:H	1.73	0.54
1:A:185:ASN:HD21	1:A:314:THR:N	1.98	0.54
1:A:225:HIS:CD2	1:A:247:ALA:H	2.25	0.54
1:A:195:LEU:HB3	1:A:236:LEU:HD23	1.90	0.54
1:A:344:MET:CE	1:A:526:TRP:HH2	2.20	0.53
1:A:471:HIS:O	1:A:472:THR:CB	2.45	0.53
1:B:78:ARG:HA	1:B:193:MET:CE	2.31	0.53
1:B:74:LYS:HZ1	1:B:283:MET:HE1	1.73	0.53
1:B:352:GLY:HA3	1:B:413:MET:O	2.07	0.53
1:B:500:MET:HE2	1:B:500:MET:HA	1.89	0.53
1:A:77:GLN:C	1:A:193:MET:HE1	2.33	0.52
1:B:105:VAL:HG21	1:B:292:LYS:HG3	1.91	0.52
1:B:185:ASN:HD21	1:B:314:THR:N	2.06	0.52
1:A:445:ARG:HH11	1:A:540:ASN:HD21	1.56	0.52
1:A:132:TYR:HA	1:A:215:PHE:O	2.09	0.52
1:A:95:SER:O	1:A:97:GLN:HB2	2.10	0.51
1:B:228:MET:HE3	1:B:341:ILE:HD11	1.92	0.51
1:A:100:PHE:N	1:A:100:PHE:CD2	2.78	0.51
1:B:48:LEU:HD22	2:B:764:HOH:O	2.10	0.50
1:B:440:PRO:HB3	1:B:529:ILE:O	2.10	0.50
1:B:423:PHE:CE2	1:B:427:ARG:HD2	2.46	0.50
1:A:371:THR:HB	1:A:373:VAL:HG23	1.93	0.50
1:B:280:GLU:O	1:B:284:LYS:HG3	2.12	0.50
1:A:100:PHE:HD2	1:A:100:PHE:N	2.09	0.50
1:B:251:TRP:HD1	1:B:420:TYR:HD1	1.58	0.49
1:B:185:ASN:ND2	1:B:313:PRO:HA	2.27	0.49
1:A:241:ILE:HG23	1:A:344:MET:HB3	1.93	0.49
1:B:347:ASN:HB3	1:B:351:GLU:OE1	2.13	0.49
1:A:373:VAL:HG11	1:A:385:ARG:O	2.12	0.48
1:B:6:SER:HB3	1:B:9:GLU:CG	2.44	0.48
1:A:376:VAL:O	1:A:385:ARG:NH2	2.47	0.48
1:A:445:ARG:HH11	1:A:540:ASN:ND2	2.10	0.47
1:B:493:GLU:H	1:B:493:GLU:CD	2.21	0.47
1:A:225:HIS:CD2	1:A:247:ALA:HB3	2.50	0.47
1:B:100:PHE:HD2	1:B:100:PHE:H	1.61	0.47
1:B:301:GLU:CD	1:B:301:GLU:H	2.22	0.47
1:A:226:TYR:OH	1:A:253:ASN:HB3	2.14	0.47
1:B:225:HIS:O	1:B:229:LEU:HG	2.14	0.47
1:B:526:TRP:CZ3	1:B:539:LEU:HD22	2.49	0.47
1:A:557:LYS:HE3	1:A:560:GLU:OE1	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:49:THR:CG2	1:A:51:TYR:H	2.29	0.46
1:A:423:PHE:HB3	1:A:424:PRO:HD3	1.98	0.46
1:B:423:PHE:HB3	1:B:424:PRO:HD3	1.97	0.46
1:A:279:LEU:HG	1:A:283:MET:HE3	1.98	0.46
1:A:497:ILE:O	1:A:501:THR:HB	2.16	0.46
1:B:335:THR:HG23	2:B:754:HOH:O	2.16	0.46
1:A:351:GLU:HG2	1:A:421:PHE:CE2	2.51	0.46
1:B:269:TYR:CA	1:B:270:LYS:HB2	2.45	0.46
1:B:100:PHE:HD2	1:B:100:PHE:N	2.14	0.45
1:B:268:GLY:O	1:B:270:LYS:HB2	2.16	0.45
1:B:445:ARG:HH11	1:B:540:ASN:HD21	1.63	0.45
1:B:100:PHE:N	1:B:100:PHE:CD2	2.84	0.45
1:B:269:TYR:HA	1:B:270:LYS:CB	2.46	0.45
1:B:534:GLU:CD	1:B:534:GLU:N	2.73	0.45
1:B:428:LEU:O	1:B:432:ARG:HB2	2.17	0.45
1:B:177:SER:HB3	1:B:180:LEU:HD22	1.98	0.44
1:A:351:GLU:HG2	1:A:421:PHE:CD2	2.52	0.44
1:A:197:TRP:O	1:A:201:ASN:HB2	2.18	0.44
1:A:29:THR:HG22	1:A:30:ASN:N	2.32	0.44
1:B:269:TYR:CZ	1:B:271:GLY:HA3	2.53	0.44
1:B:234:ARG:NH1	1:B:337:TRP:HB3	2.33	0.43
1:A:535:VAL:HG13	1:A:553:MET:HG2	2.00	0.43
1:A:236:LEU:HD11	2:A:677:HOH:O	2.18	0.43
1:A:344:MET:HE2	1:A:444:TYR:CD2	2.54	0.43
1:A:37:THR:HG22	1:A:39:TYR:H	1.83	0.43
1:A:423:PHE:O	1:A:427:ARG:HG3	2.19	0.43
1:A:49:THR:HB	1:A:53:ASP:O	2.19	0.43
1:A:217:GLU:OE2	1:A:218:SER:HB2	2.18	0.43
1:A:218:SER:HA	1:A:244:SER:O	2.18	0.43
1:A:380:LEU:HB3	1:A:393:MET:SD	2.59	0.43
1:B:132:TYR:CE1	1:B:216:GLY:HA2	2.53	0.43
1:B:299:THR:O	1:B:303:ARG:HG3	2.18	0.43
1:A:92:LYS:HE2	1:A:113:TYR:CZ	2.53	0.42
1:A:227:MET:O	1:A:233:THR:OG1	2.37	0.42
1:B:105:VAL:HG11	1:B:288:GLN:HB2	2.02	0.42
1:A:259:ARG:NH2	2:A:605:HOH:O	2.45	0.42
1:A:440:PRO:CB	1:A:528:PRO:HB2	2.49	0.42
1:B:3:PHE:CZ	1:B:358:ILE:HD11	2.54	0.42
1:B:445:ARG:HH11	1:B:540:ASN:ND2	2.17	0.42
1:A:185:ASN:HD21	1:A:313:PRO:HA	1.82	0.42
1:A:217:GLU:HA	1:A:243:MET:HB2	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:269:TYR:CA	1:A:270:LYS:CB	2.83	0.42
1:A:5:VAL:HG21	1:A:101:ILE:HA	2.02	0.42
1:B:418:HIS:HA	1:B:422:TRP:HB2	2.01	0.42
1:A:112:LEU:HD21	1:A:166:ARG:HG3	2.02	0.41
1:A:49:THR:HG22	1:A:51:TYR:N	2.33	0.41
1:A:457:TYR:CG	1:A:472:THR:HG21	2.54	0.41
1:B:97:GLN:NE2	1:B:141:GLY:H	2.12	0.41
1:B:242:LEU:HB2	1:B:345:MET:HG2	2.03	0.41
1:B:258:HIS:HD2	2:B:716:HOH:O	2.03	0.41
1:B:261:PHE:CZ	1:B:265:LYS:HD2	2.55	0.41
1:B:368:GLU:O	1:B:371:THR:O	2.38	0.41
1:A:72:ARG:NH2	1:A:282:LEU:O	2.41	0.41
1:A:242:LEU:HB2	1:A:345:MET:HG2	2.02	0.41
1:B:226:TYR:O	1:B:230:THR:HG22	2.21	0.41
1:A:183:PRO:HG2	1:A:189:LYS:NZ	2.35	0.41
1:A:200:ASN:HB3	2:A:693:HOH:O	2.19	0.41
1:A:423:PHE:CE2	1:A:427:ARG:HD2	2.56	0.41
1:B:80:THR:HA	1:B:81:PRO:HD3	1.86	0.40
1:B:142:GLU:CD	1:B:142:GLU:N	2.79	0.40
1:A:29:THR:HG22	2:A:746:HOH:O	2.21	0.40
1:B:105:VAL:CG2	1:B:292:LYS:HG3	2.51	0.40
1:B:225:HIS:CD2	1:B:247:ALA:H	2.26	0.40
1:A:218:SER:OG	1:A:471:HIS:HE1	2.04	0.40
1:B:428:LEU:HD11	2:B:759:HOH:O	2.21	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	567/577 (98%)	537 (95%)	28 (5%)	2 (0%)	30 34

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	567/577 (98%)	545 (96%)	20 (4%)	2 (0%)	30	34
All	All	1134/1154 (98%)	1082 (95%)	48 (4%)	4 (0%)	30	34

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	270	LYS
1	B	270	LYS
1	A	472	THR
1	B	472	THR

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	501/511 (98%)	475 (95%)	26 (5%)	21	26
1	B	498/511 (98%)	469 (94%)	29 (6%)	18	22
All	All	999/1022 (98%)	944 (94%)	55 (6%)	19	24

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	THR
1	A	100	PHE
1	A	121	LEU
1	A	124	GLU
1	A	142	GLU
1	A	170	LEU
1	A	180	LEU
1	A	195	LEU
1	A	266	LEU
1	A	280	GLU
1	A	290	LEU
1	A	292	LYS

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Mol	Chain	Res	Type
1	A	314	THR
1	A	351	GLU
1	A	356	THR
1	A	360	LYS
1	A	370	GLU
1	A	378	SER
1	A	390	THR
1	A	406	THR
1	A	429	LEU
1	A	472	THR
1	A	490	GLU
1	A	501	THR
1	A	517	GLU
1	A	534	GLU
1	B	37	THR
1	B	72	ARG
1	B	80	THR
1	B	100	PHE
1	B	114	LEU
1	B	125	THR
1	B	126	LYS
1	B	173	LEU
1	B	180	LEU
1	B	188	LEU
1	B	195	LEU
1	B	242	LEU
1	B	266	LEU
1	B	290	LEU
1	B	301	GLU
1	B	306	LYS
1	B	351	GLU
1	B	389	GLU
1	B	399	LYS
1	B	403	THR
1	B	425	MET
1	B	427	ARG
1	B	429	LEU
1	B	472	THR
1	B	485	LYS
1	B	519	GLU
1	B	534	GLU
1	B	537	LYS

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Mol	Chain	Res	Type
1	B	565	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	20	ASN
1	A	65	GLN
1	A	97	GLN
1	A	162	ASN
1	A	164	GLN
1	A	185	ASN
1	A	225	HIS
1	A	255	GLN
1	A	274	ASN
1	A	288	GLN
1	A	328	HIS
1	A	471	HIS
1	A	516	ASN
1	A	540	ASN
1	B	77	GLN
1	B	97	GLN
1	B	164	GLN
1	B	185	ASN
1	B	191	GLN
1	B	204	ASN
1	B	208	ASN
1	B	225	HIS
1	B	238	HIS
1	B	255	GLN
1	B	257	GLN
1	B	258	HIS
1	B	288	GLN
1	B	411	ASN
1	B	418	HIS
1	B	435	HIS
1	B	540	ASN

5.3.3 RNA ⓘ

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 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 [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	569/577 (98%)	-1.31	0 100 100	22, 36, 52, 74	0
1	B	569/577 (98%)	-1.32	0 100 100	22, 36, 60, 73	0
All	All	1138/1154 (98%)	-1.31	0 100 100	22, 36, 56, 74	0

There are no RSRZ outliers to report.

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.