



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 2, 2026 – 02:24 AM UTC

PDB ID : 5Z70 / pdb_00005z70
Title : Crystal structure of oleate hydratase from *Stenotrophomonas* sp. KCTC 12332
Authors : Park, A.K.
Deposited on : 2018-01-26
Resolution : 2.91 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

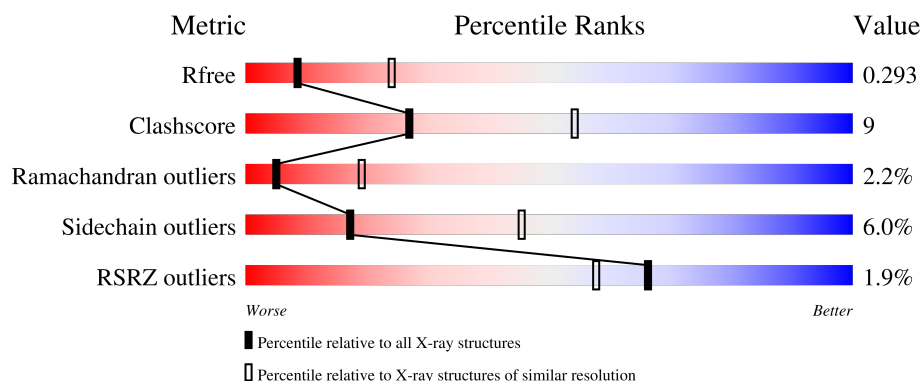
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

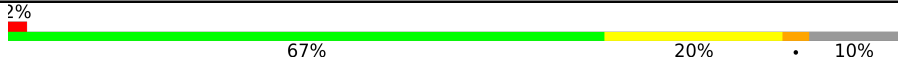
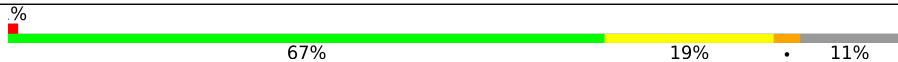
The reported resolution of this entry is 2.91 Å.

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	2995 (2.94-2.90)
Clashscore	190562	3213 (2.94-2.90)
Ramachandran outliers	187476	3128 (2.94-2.90)
Sidechain outliers	187428	3130 (2.94-2.90)
RSRZ outliers	180081	2995 (2.94-2.90)

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	605	
1	B	605	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8606 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 Oleate hydratase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	543	Total	C	N	O	S	0	0	0
			4316	2754	732	809	21			
1	B	538	Total	C	N	O	S	0	0	0
			4283	2733	727	802	21			

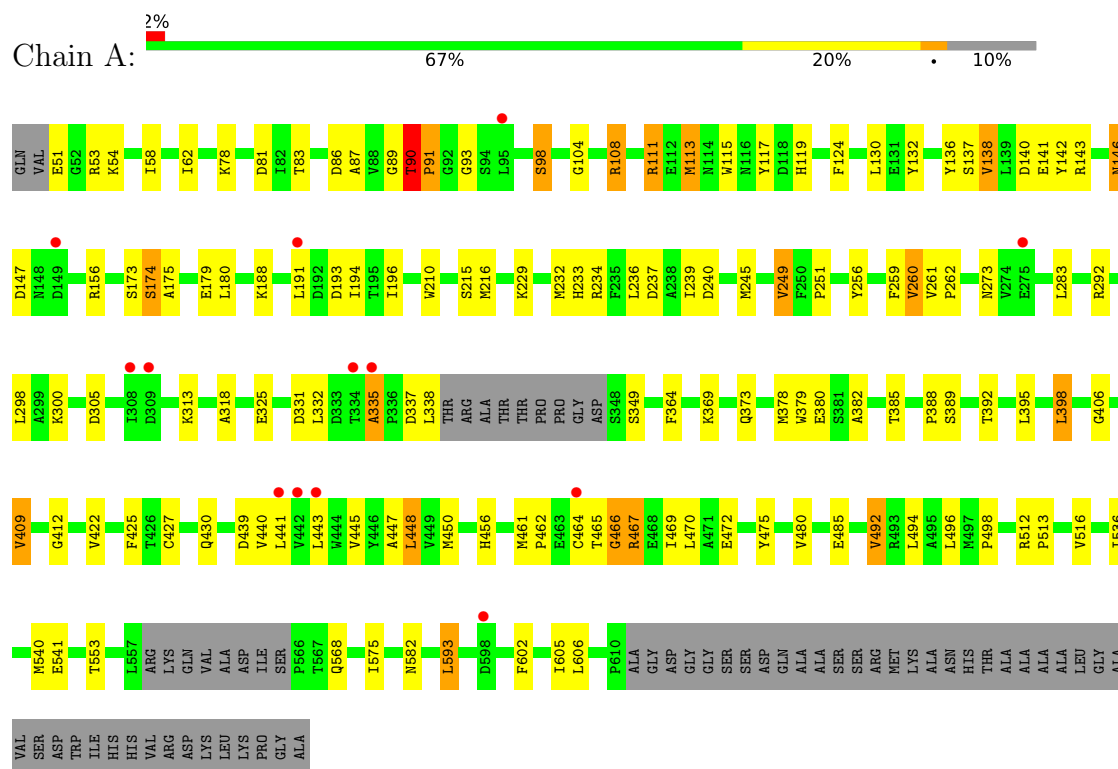
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	3	Total	O	0	0
			3	3		
2	B	4	Total	O	0	0
			4	4		

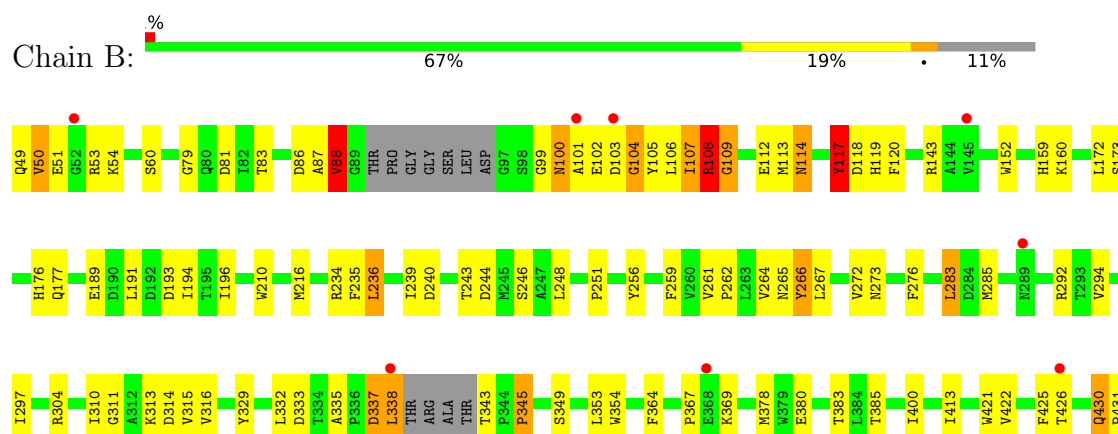
3 Residue-property plots

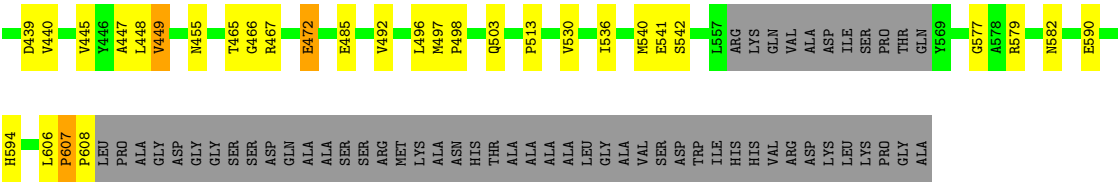
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Oleate hydratase



• Molecule 1: Oleate hydratase





4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	62.06Å 235.25Å 176.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.83 – 2.91 35.83 – 2.91	Depositor EDS
% Data completeness (in resolution range)	99.5 (35.83-2.91) 99.5 (35.83-2.91)	Depositor EDS
R_{merge}	0.26	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.86 (at 2.91Å)	Xtriage
Refinement program	PHENIX 1.12_2829	Depositor
R, R_{free}	0.227 , 0.290 0.230 , 0.293	Depositor DCC
R_{free} test set	1514 reflections (5.23%)	wwPDB-VP
Wilson B-factor (Å ²)	74.6	Xtriage
Anisotropy	0.445	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 27.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	8606	wwPDB-VP
Average B, all atoms (Å ²)	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.97% 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	0.17	0/4421	0.45	0/6003
1	B	0.24	0/4386	0.49	4/5954 (0.1%)
All	All	0.20	0/8807	0.47	4/11957 (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 (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	109	GLY	N-CA-C	-6.22	98.44	113.18
1	B	108	ARG	N-CA-C	5.72	122.98	110.80
1	B	107	ILE	CB-CA-C	-5.39	105.41	111.23
1	B	117	TYR	CA-CB-CG	5.02	122.93	113.90

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	51	GLU	Peptide
1	B	430	GLN	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	4316	0	4226	77	0
1	B	4283	0	4189	79	0
2	A	3	0	0	0	0
2	B	4	0	0	0	0
All	All	8606	0	8415	156	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 9.

All (156) 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:234:ARG:NH2	1:A:541:GLU:OE2	2.08	0.87
1:A:90:THR:HB	1:A:91:PRO:HD3	1.56	0.87
1:A:300:LYS:NZ	1:A:305:ASP:OD1	2.11	0.84
1:B:607:PRO:HB2	1:B:608:PRO:HD2	1.67	0.76
1:A:332:LEU:HA	1:A:466:GLY:HA3	1.68	0.75
1:B:310:ILE:HG23	1:B:314:ASP:HB2	1.68	0.75
1:B:88:VAL:HA	1:B:276:PHE:HB3	1.69	0.75
1:B:455:ASN:N	1:B:472:GLU:OE2	2.20	0.73
1:B:99:GLY:HA2	1:B:106:LEU:H	1.56	0.71
1:A:516:VAL:HG21	1:A:553:THR:HG21	1.73	0.70
1:A:156:ARG:HH22	1:A:215:SER:HB2	1.57	0.70
1:B:579:ARG:NH1	1:B:607:PRO:HG2	2.07	0.69
1:B:378:MET:HB2	1:B:498:PRO:HA	1.75	0.69
1:A:470:LEU:HD21	1:A:492:VAL:HG11	1.76	0.67
1:A:62:ILE:HG21	1:A:540:MET:HG3	1.78	0.66
1:B:345:PRO:HD2	1:B:349:SER:HB2	1.78	0.65
1:B:234:ARG:NH2	1:B:541:GLU:OE2	2.30	0.65
1:B:579:ARG:NH1	1:B:590:GLU:OE2	2.31	0.63
1:B:117:TYR:HD2	1:B:117:TYR:H	1.46	0.63
1:B:193:ASP:OD1	1:B:369:LYS:NZ	2.30	0.63
1:B:332:LEU:HA	1:B:466:GLY:HA3	1.81	0.63
1:B:117:TYR:HB3	1:B:120:PHE:HB3	1.81	0.62
1:B:364:PHE:HA	1:B:513:PRO:HB3	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:335:ALA:HA	1:A:465:THR:HG21	1.82	0.61
1:A:380:GLU:OE2	1:A:494:LEU:HD23	2.01	0.61
1:B:261:VAL:HG22	1:B:262:PRO:HD3	1.82	0.61
1:A:188:LYS:HE3	1:A:229:LYS:HD3	1.81	0.60
1:B:191:LEU:HA	1:B:194:ILE:HD12	1.82	0.59
1:A:191:LEU:HA	1:A:194:ILE:HD12	1.85	0.59
1:A:216:MET:HE1	1:A:245:MET:SD	2.43	0.58
1:B:53:ARG:NH1	1:B:313:LYS:O	2.36	0.58
1:B:607:PRO:HB2	1:B:608:PRO:CD	2.33	0.58
1:B:53:ARG:NH1	1:B:315:VAL:HG13	2.19	0.58
1:A:378:MET:HB2	1:A:498:PRO:HA	1.86	0.57
1:A:395:LEU:HA	1:A:398:LEU:HD12	1.86	0.57
1:B:87:ALA:O	1:B:88:VAL:HG12	2.03	0.57
1:A:54:LYS:HA	1:A:81:ASP:HB3	1.86	0.56
1:A:475:TYR:HA	1:A:480:VAL:HB	1.86	0.56
1:A:193:ASP:OD1	1:A:369:LYS:NZ	2.38	0.56
1:B:60:SER:HB3	1:B:86:ASP:OD2	2.06	0.56
1:B:503:GLN:HB2	1:B:536:ILE:HD12	1.87	0.56
1:B:49:GLN:HG3	1:B:50:VAL:HG23	1.89	0.55
1:A:575:ILE:HD13	1:A:606:LEU:HD23	1.88	0.55
1:A:124:PHE:HB3	1:A:138:VAL:HG12	1.88	0.55
1:A:389:SER:H	1:A:392:THR:HG22	1.70	0.55
1:A:115:TRP:CZ2	1:A:117:TYR:HB3	2.42	0.54
1:A:331:ASP:HA	1:A:494:LEU:HD13	1.90	0.54
1:B:422:VAL:HB	1:B:448:LEU:HB2	1.89	0.54
1:B:172:LEU:HD22	1:B:176:HIS:HB3	1.88	0.54
1:A:450:MET:HE1	1:A:461:MET:HE2	1.90	0.54
1:B:594:HIS:CD2	1:B:606:LEU:HD13	2.43	0.53
1:B:264:VAL:O	1:B:266:TYR:N	2.36	0.53
1:B:177:GLN:NE2	1:B:582:ASN:HA	2.22	0.53
1:B:216:MET:HE2	1:B:248:LEU:HD11	1.91	0.53
1:A:466:GLY:HA2	1:A:492:VAL:HG21	1.91	0.53
1:A:196:ILE:HG21	1:A:210:TRP:CZ2	2.43	0.52
1:A:325:GLU:OE1	1:A:349:SER:HB2	2.09	0.52
1:B:104:GLY:O	1:B:385:THR:OG1	2.16	0.52
1:A:256:TYR:HA	1:A:260:VAL:HG13	1.92	0.52
1:A:512:ARG:HH12	1:A:536:ILE:HG12	1.74	0.52
1:B:354:TRP:CD2	1:B:367:PRO:HB3	2.46	0.51
1:B:112:GLU:N	1:B:112:GLU:OE1	2.44	0.51
1:A:406:GLY:HA2	1:A:430:GLN:O	2.11	0.51
1:A:83:THR:HG22	1:A:273:ASN:HB2	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:98:SER:OG	1:A:108:ARG:NH2	2.44	0.50
1:A:175:ALA:O	1:A:179:GLU:HG3	2.12	0.50
1:A:173:SER:OG	1:A:174:SER:N	2.44	0.50
1:A:180:LEU:HB3	1:A:232:MET:HE2	1.94	0.50
1:A:233:HIS:O	1:A:568:GLN:HB3	2.12	0.50
1:B:196:ILE:HG21	1:B:210:TRP:CZ2	2.47	0.49
1:A:382:ALA:HB3	1:A:445:VAL:HG12	1.95	0.49
1:B:425:PHE:HB2	1:B:445:VAL:HG22	1.94	0.49
1:A:143:ARG:O	1:A:147:ASP:HB2	2.12	0.49
1:B:106:LEU:HD21	1:B:329:TYR:CE1	2.47	0.49
1:A:245:MET:HE3	1:A:412:GLY:HA3	1.95	0.48
1:B:311:GLY:O	1:B:313:LYS:N	2.40	0.48
1:A:232:MET:O	1:A:236:LEU:HG	2.13	0.48
1:B:114:ASN:HD21	1:B:540:MET:HE2	1.77	0.48
1:B:335:ALA:HB2	1:B:465:THR:HG22	1.94	0.48
1:B:236:LEU:HD22	1:B:577:GLY:HA3	1.95	0.48
1:A:335:ALA:CA	1:A:465:THR:HG21	2.43	0.48
1:A:409:VAL:HG13	1:A:427:CYS:O	2.14	0.48
1:A:86:ASP:OD1	1:A:87:ALA:N	2.46	0.47
1:A:380:GLU:HB2	1:A:496:LEU:HD13	1.95	0.47
1:B:117:TYR:O	1:B:119:HIS:N	2.44	0.47
1:B:189:GLU:OE2	1:B:189:GLU:N	2.41	0.47
1:A:58:ILE:HB	1:A:318:ALA:HA	1.96	0.47
1:B:117:TYR:N	1:B:117:TYR:CD2	2.82	0.47
1:B:117:TYR:HB2	1:B:118:ASP:H	1.53	0.47
1:B:235:PHE:O	1:B:239:ILE:HG23	2.14	0.47
1:A:251:PRO:HB3	1:A:259:PHE:CZ	2.49	0.47
1:A:90:THR:HB	1:A:91:PRO:CD	2.36	0.47
1:B:54:LYS:HA	1:B:81:ASP:HB3	1.96	0.47
1:B:283:LEU:HD23	1:B:297:ILE:HG12	1.96	0.47
1:A:380:GLU:HB3	1:A:461:MET:HE3	1.96	0.47
1:A:146:ASN:CG	1:A:249:VAL:HG23	2.40	0.46
1:B:378:MET:CB	1:B:498:PRO:HA	2.44	0.46
1:A:425:PHE:HB3	1:A:445:VAL:HG23	1.98	0.46
1:B:259:PHE:C	1:B:262:PRO:HD2	2.41	0.45
1:B:294:VAL:HG11	1:B:316:VAL:HG21	1.98	0.45
1:B:79:GLY:O	1:B:272:VAL:HA	2.16	0.45
1:B:311:GLY:C	1:B:313:LYS:H	2.24	0.45
1:B:60:SER:HB3	1:B:86:ASP:CG	2.42	0.45
1:A:422:VAL:HB	1:A:448:LEU:HB2	1.98	0.45
1:A:388:PRO:HD3	1:A:439:ASP:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:461:MET:N	1:A:462:PRO:HD2	2.31	0.45
1:B:256:TYR:O	1:B:261:VAL:HG13	2.17	0.45
1:B:251:PRO:HB3	1:B:259:PHE:CZ	2.52	0.44
1:A:132:TYR:HB3	1:A:136:TYR:HD2	1.83	0.44
1:A:364:PHE:HA	1:A:513:PRO:HB3	1.99	0.44
1:B:337:ASP:OD1	1:B:338:LEU:N	2.50	0.44
1:B:421:TRP:HA	1:B:449:VAL:HG22	1.99	0.44
1:B:579:ARG:HH12	1:B:607:PRO:HG2	1.77	0.44
1:A:379:TRP:HB2	1:A:447:ALA:O	2.17	0.44
1:A:602:PHE:O	1:A:605:ILE:HG12	2.17	0.44
1:B:244:ASP:OD1	1:B:246:SER:HB3	2.17	0.44
1:B:413:ILE:HG22	1:B:426:THR:OG1	2.18	0.44
1:B:421:TRP:HB3	1:B:447:ALA:HB1	2.00	0.44
1:A:237:ASP:OD1	1:A:237:ASP:N	2.51	0.43
1:B:285:MET:HE2	1:B:292:ARG:HB3	1.98	0.43
1:A:113:MET:H	1:A:113:MET:HE2	1.83	0.43
1:B:582:ASN:O	1:B:582:ASN:ND2	2.51	0.43
1:A:117:TYR:CD1	1:A:119:HIS:HB2	2.54	0.43
1:A:369:LYS:HE2	1:A:369:LYS:HB3	1.82	0.43
1:A:485:GLU:OE1	1:A:485:GLU:N	2.48	0.43
1:B:267:LEU:HB3	1:B:272:VAL:HG21	2.01	0.43
1:A:332:LEU:HD12	1:A:467:ARG:N	2.33	0.43
1:B:234:ARG:NH1	1:B:235:PHE:CZ	2.87	0.43
1:A:138:VAL:HG21	1:A:262:PRO:HG2	2.00	0.42
1:A:130:LEU:H	1:A:141:GLU:CD	2.28	0.42
1:A:137:SER:OG	1:A:140:ASP:OD1	2.31	0.42
1:B:54:LYS:HE2	1:B:54:LYS:HB3	1.90	0.42
1:A:430:GLN:HG3	1:A:441:LEU:HD22	2.01	0.42
1:B:173:SER:O	1:B:177:GLN:HG3	2.20	0.42
1:B:369:LYS:HE2	1:B:369:LYS:HB3	1.65	0.42
1:B:109:GLY:HA3	1:B:497:MET:SD	2.60	0.42
1:B:152:TRP:CZ2	1:B:400:ILE:HD12	2.55	0.42
1:B:107:ILE:HG13	1:B:383:THR:HG21	2.01	0.42
1:A:335:ALA:CB	1:A:465:THR:HG21	2.49	0.41
1:A:337:ASP:OD1	1:A:338:LEU:N	2.54	0.41
1:A:382:ALA:HB3	1:A:445:VAL:CG1	2.51	0.41
1:A:593:LEU:HD13	1:A:593:LEU:HA	1.90	0.41
1:B:332:LEU:HD21	1:B:467:ARG:NH2	2.36	0.41
1:A:111:ARG:NH2	1:A:216:MET:O	2.54	0.41
1:A:380:GLU:HB3	1:A:461:MET:CE	2.50	0.41
1:B:83:THR:HG23	1:B:273:ASN:HB2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:142:TYR:CE1	1:A:259:PHE:HZ	2.39	0.40
1:A:456:HIS:HB2	1:A:475:TYR:CG	2.56	0.40
1:B:152:TRP:CE2	1:B:400:ILE:HD12	2.56	0.40
1:B:234:ARG:NH1	1:B:235:PHE:CE1	2.89	0.40
1:B:264:VAL:C	1:B:266:TYR:H	2.23	0.40
1:B:332:LEU:HD23	1:B:333:ASP:CG	2.46	0.40
1:B:380:GLU:HB2	1:B:496:LEU:HD13	2.03	0.40
1:A:216:MET:HB2	1:A:216:MET:HE2	1.64	0.40
1:A:53:ARG:HD3	1:A:313:LYS:O	2.22	0.40
1:B:261:VAL:CG2	1:B:262:PRO:HD3	2.48	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	537/605 (89%)	493 (92%)	34 (6%)	10 (2%)	6	22
1	B	530/605 (88%)	484 (91%)	32 (6%)	14 (3%)	4	15
All	All	1067/1210 (88%)	977 (92%)	66 (6%)	24 (2%)	5	19

All (24) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	91	PRO
1	A	108	ARG
1	A	335	ALA
1	B	50	VAL
1	B	88	VAL
1	B	108	ARG
1	B	265	ASN
1	B	430	GLN

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Mol	Chain	Res	Type
1	A	466	GLY
1	B	431	PRO
1	A	90	THR
1	B	101	ALA
1	B	103	ASP
1	B	104	GLY
1	B	266	TYR
1	B	345	PRO
1	A	111	ARG
1	B	100	ASN
1	B	105	TYR
1	B	607	PRO
1	A	98	SER
1	A	104	GLY
1	A	93	GLY
1	A	89	GLY

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	469/513 (91%)	441 (94%)	28 (6%)	17	45
1	B	465/513 (91%)	437 (94%)	28 (6%)	17	45
All	All	934/1026 (91%)	878 (94%)	56 (6%)	17	45

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	78	LYS
1	A	90	THR
1	A	113	MET
1	A	138	VAL
1	A	146	ASN
1	A	174	SER
1	A	239	ILE

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Mol	Chain	Res	Type
1	A	240	ASP
1	A	249	VAL
1	A	260	VAL
1	A	261	VAL
1	A	283	LEU
1	A	292	ARG
1	A	298	LEU
1	A	373	GLN
1	A	385	THR
1	A	398	LEU
1	A	409	VAL
1	A	440	VAL
1	A	443	LEU
1	A	448	LEU
1	A	464	CYS
1	A	467	ARG
1	A	469	ILE
1	A	472	GLU
1	A	492	VAL
1	A	582	ASN
1	A	593	LEU
1	B	51	GLU
1	B	88	VAL
1	B	100	ASN
1	B	102	GLU
1	B	108	ARG
1	B	113	MET
1	B	114	ASN
1	B	117	TYR
1	B	143	ARG
1	B	159	HIS
1	B	160	LYS
1	B	236	LEU
1	B	240	ASP
1	B	243	THR
1	B	283	LEU
1	B	304	ARG
1	B	337	ASP
1	B	338	LEU
1	B	343	THR
1	B	353	LEU
1	B	439	ASP

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Mol	Chain	Res	Type
1	B	440	VAL
1	B	449	VAL
1	B	472	GLU
1	B	485	GLU
1	B	492	VAL
1	B	530	VAL
1	B	542	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	119	HIS
1	A	163	GLN
1	A	177	GLN
1	A	428	ASN
1	B	114	ASN
1	B	177	GLN
1	B	482	GLN
1	B	584	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 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

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	543/605 (89%)	0.19	13 (2%) 59 50	41, 78, 109, 139	0
1	B	538/605 (88%)	0.02	8 (1%) 72 64	40, 65, 92, 110	0
All	All	1081/1210 (89%)	0.10	21 (1%) 66 58	40, 70, 105, 139	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	289	ASN	4.3
1	A	464	CYS	4.2
1	A	309	ASP	3.6
1	A	443	LEU	3.1
1	B	101	ALA	3.1
1	B	338	LEU	3.0
1	A	334	THR	3.0
1	A	191	LEU	2.9
1	B	145	VAL	2.9
1	A	275	GLU	2.8
1	B	368	GLU	2.7
1	B	52	GLY	2.7
1	A	95	LEU	2.6
1	A	308	ILE	2.4
1	B	426	THR	2.2
1	A	598	ASP	2.2
1	A	441	LEU	2.2
1	B	103	ASP	2.1
1	A	335	ALA	2.1
1	A	442	VAL	2.1
1	A	149	ASP	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 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.