Full wwPDB X-ray Structure Validation Report ⓘ

Aug 3, 2023 – 07:51 AM EDT

PDB ID	:	1F6W
Title	:	STRUCTURE OF THE CATALYTIC DOMAIN OF HUMAN BILE SALT ACTIVATED LIPASE
Authors	:	Terzyan, S.; Zhang, X.
Deposited on	:	2000-06-23
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:

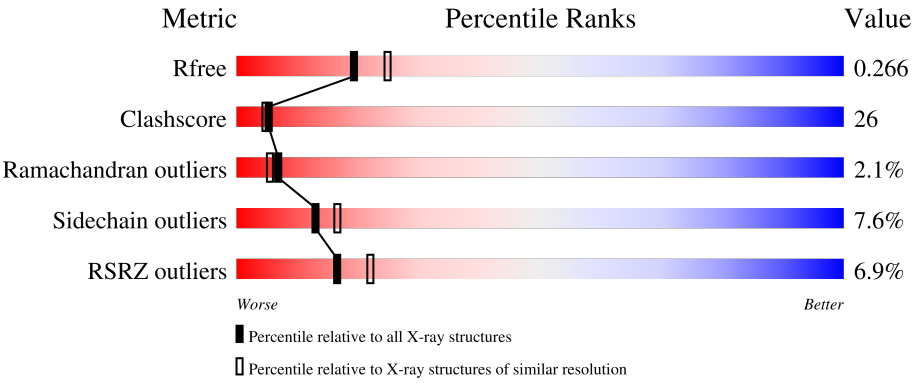
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.34
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.34

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	533	7%, 58%, 37%, 5%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 4314 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 BILE SALT ACTIVATED LIPASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	533	Total	C	N	O	S	0	0	0
			4151	2668	698	769	16			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	186	ASP	ASN	engineered mutation	UNP P19835
A	298	ASP	ALA	engineered mutation	UNP P19835

- Molecule 2 is water.

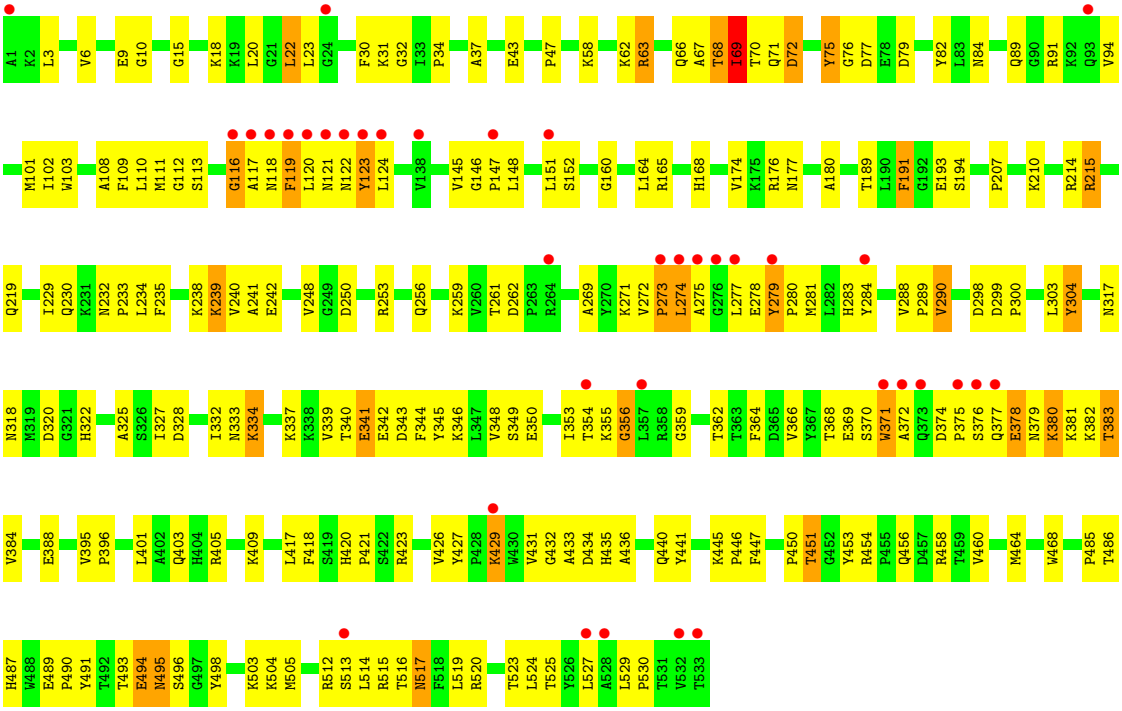
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	163	Total	O	0	0
			163	163		

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: BILE SALT ACTIVATED LIPASE

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	64.70Å 88.97Å 104.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.19 – 2.30 24.18 – 2.30	Depositor EDS
% Data completeness (in resolution range)	85.6 (24.19-2.30) 85.7 (24.18-2.30)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.65 (at 2.31Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.212 , 0.267 0.213 , 0.266	Depositor DCC
R_{free} test set	1258 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å^2)	44.4	Xtriage
Anisotropy	0.584	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.28 , 44.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	4314	wwPDB-VP
Average B, all atoms (Å^2)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.92% 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

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	0.55	0/4264	0.75	2/5801 (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

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	215	ARG	NE-CZ-NH2	-5.42	117.59	120.30
1	A	409	LYS	N-CA-C	-5.03	97.42	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	304	TYR	Sidechain

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	4151	0	4092	218	0
2	A	163	0	0	10	0
All	All	4314	0	4092	218	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 26.

All (218) 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:372:ALA:HB1	1:A:375:PRO:HG3	1.19	1.12
1:A:274:LEU:HD11	1:A:277:LEU:HG	1.29	1.07
1:A:512:ARG:HG2	1:A:513:SER:H	1.21	1.05
1:A:9:GLU:OE1	1:A:176:ARG:HD2	1.61	1.00
1:A:66:GLN:HE22	1:A:112:GLY:H	1.10	0.92
1:A:274:LEU:CD1	1:A:277:LEU:HG	1.99	0.92
1:A:117:ALA:HB3	1:A:124:LEU:HB3	1.50	0.91
1:A:372:ALA:HB1	1:A:375:PRO:CG	2.03	0.85
1:A:31:LYS:HA	1:A:84:ASN:HD22	1.42	0.84
1:A:512:ARG:HG2	1:A:513:SER:N	1.94	0.82
1:A:334:LYS:HB3	1:A:337:LYS:HB2	1.62	0.82
1:A:340:THR:HG22	1:A:342:GLU:H	1.45	0.80
1:A:372:ALA:O	1:A:375:PRO:HD3	1.82	0.80
1:A:280:PRO:HA	1:A:328:ASP:OD1	1.82	0.79
1:A:441:TYR:HB2	1:A:464:MET:HE3	1.65	0.78
1:A:519:LEU:O	1:A:523:THR:HG23	1.85	0.77
1:A:372:ALA:CB	1:A:375:PRO:HG3	2.08	0.76
1:A:70:THR:OG1	1:A:72:ASP:HB2	1.86	0.75
1:A:63:ARG:HD2	2:A:547:HOH:O	1.87	0.74
1:A:512:ARG:CG	1:A:513:SER:H	1.97	0.73
1:A:32:GLY:H	1:A:84:ASN:ND2	1.86	0.72
1:A:350:GLU:O	1:A:353:ILE:HG13	1.88	0.72
1:A:66:GLN:NE2	1:A:112:GLY:H	1.85	0.72
1:A:374:ASP:OD1	1:A:377:GLN:HB2	1.91	0.70
1:A:68:THR:HA	1:A:75:TYR:CE2	2.26	0.70
1:A:354:THR:O	1:A:355:LYS:HD2	1.91	0.70
1:A:118:ASN:O	1:A:119:PHE:CD1	2.44	0.70
1:A:451:THR:HB	2:A:655:HOH:O	1.92	0.69
1:A:487:HIS:HE1	1:A:489:GLU:OE1	1.77	0.68
1:A:426:VAL:HG23	1:A:427:TYR:HD1	1.60	0.66
1:A:240:VAL:HG13	1:A:269:ALA:CB	2.25	0.66
1:A:340:THR:HB	1:A:343:ASP:CG	2.16	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:441:TYR:HB2	1:A:464:MET:CE	2.24	0.66
1:A:317:ASN:ND2	1:A:418:PHE:HB3	2.10	0.65
1:A:22:LEU:N	1:A:22:LEU:HD12	2.11	0.65
1:A:239:LYS:HE3	1:A:239:LYS:HA	1.77	0.65
1:A:441:TYR:CB	1:A:464:MET:HE3	2.26	0.65
1:A:118:ASN:HB3	1:A:123:TYR:CD2	2.32	0.65
1:A:234:LEU:HD11	1:A:238:LYS:HE3	1.79	0.65
1:A:23:LEU:HD12	1:A:91:ARG:NH2	2.12	0.64
1:A:340:THR:HG22	1:A:342:GLU:N	2.12	0.64
1:A:119:PHE:HB2	1:A:121:ASN:OD1	1.99	0.63
1:A:9:GLU:H	1:A:177:ASN:ND2	1.97	0.63
1:A:426:VAL:HG23	1:A:427:TYR:N	2.14	0.63
1:A:273:PRO:HB3	1:A:277:LEU:CD1	2.29	0.62
1:A:493:THR:HB	1:A:494:GLU:OE1	1.99	0.62
1:A:279:TYR:CB	1:A:280:PRO:HD2	2.30	0.61
1:A:121:ASN:CG	1:A:122:ASN:H	2.04	0.61
1:A:520:ARG:HG3	1:A:524:LEU:HD12	1.83	0.61
1:A:281:MET:SD	1:A:327:ILE:HD12	2.41	0.61
1:A:274:LEU:HD13	1:A:275:ALA:H	1.66	0.60
1:A:362:THR:O	1:A:366:VAL:HG23	2.01	0.60
1:A:423:ARG:HH11	1:A:423:ARG:HG3	1.66	0.60
1:A:62:LYS:HD3	1:A:77:ASP:HB2	1.83	0.60
1:A:274:LEU:CD1	1:A:275:ALA:H	2.15	0.60
1:A:420:HIS:O	1:A:433:ALA:HB2	2.02	0.60
1:A:118:ASN:HA	1:A:123:TYR:HA	1.83	0.60
1:A:426:VAL:HG23	1:A:427:TYR:CD1	2.37	0.59
1:A:498:TYR:HE2	1:A:514:LEU:HB2	1.67	0.59
1:A:108:ALA:O	1:A:109:PHE:HB2	2.02	0.59
1:A:513:SER:HB3	1:A:516:THR:HG23	1.84	0.59
1:A:117:ALA:O	1:A:124:LEU:N	2.30	0.59
1:A:71:GLN:HE22	1:A:272:VAL:HG12	1.68	0.59
1:A:274:LEU:CD1	1:A:275:ALA:N	2.65	0.59
1:A:71:GLN:HE22	1:A:272:VAL:H	1.50	0.58
1:A:426:VAL:HG23	1:A:427:TYR:H	1.68	0.58
1:A:234:LEU:C	1:A:234:LEU:HD13	2.24	0.58
1:A:6:VAL:HG13	1:A:180:ALA:HB1	1.86	0.57
1:A:9:GLU:H	1:A:177:ASN:HD21	1.52	0.57
1:A:232:ASN:ND2	1:A:235:PHE:HB2	2.18	0.57
1:A:18:LYS:HE2	2:A:679:HOH:O	2.04	0.57
1:A:34:PRO:HB3	1:A:82:TYR:CE2	2.39	0.56
1:A:339:VAL:HG13	1:A:376:SER:HA	1.85	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:344:PHE:O	1:A:348:VAL:HG13	2.05	0.56
1:A:435:HIS:O	1:A:436:ALA:HB3	2.06	0.56
1:A:66:GLN:NE2	1:A:111:MET:HB2	2.20	0.56
1:A:111:MET:HA	1:A:145:VAL:HG21	1.86	0.56
1:A:279:TYR:HB2	1:A:280:PRO:HD2	1.87	0.56
1:A:340:THR:HG22	1:A:341:GLU:N	2.20	0.56
1:A:68:THR:HA	1:A:75:TYR:HE2	1.69	0.55
1:A:67:ALA:O	1:A:75:TYR:HE2	1.90	0.55
1:A:495:ASN:OD1	1:A:496:SER:N	2.41	0.54
1:A:32:GLY:H	1:A:84:ASN:HD22	1.55	0.54
1:A:447:PHE:O	1:A:450:PRO:HD3	2.08	0.53
1:A:71:GLN:O	1:A:271:LYS:NZ	2.41	0.53
1:A:234:LEU:HD13	1:A:234:LEU:O	2.08	0.53
1:A:239:LYS:HE3	1:A:242:GLU:OE2	2.08	0.53
1:A:494:GLU:OE1	1:A:494:GLU:N	2.36	0.53
1:A:147:PRO:HD2	2:A:550:HOH:O	2.08	0.53
1:A:148:LEU:HD23	2:A:550:HOH:O	2.09	0.53
1:A:454:ARG:NH1	1:A:456:GLN:HE21	2.07	0.53
1:A:164:LEU:HD11	1:A:289:PRO:HB3	1.91	0.53
1:A:380:LYS:HA	1:A:383:THR:HG23	1.91	0.52
1:A:370:SER:C	1:A:372:ALA:N	2.61	0.52
1:A:374:ASP:HB2	1:A:379:ASN:HB3	1.92	0.52
1:A:230:GLN:HB2	1:A:288:VAL:CG2	2.39	0.52
1:A:250:ASP:HB3	1:A:253:ARG:HB2	1.90	0.52
1:A:498:TYR:CE2	1:A:514:LEU:HB2	2.44	0.52
1:A:43:GLU:OE2	1:A:259:LYS:HD3	2.10	0.52
1:A:333:ASN:OD1	1:A:426:VAL:HG11	2.10	0.52
1:A:274:LEU:HD12	1:A:275:ALA:N	2.24	0.52
1:A:503:LYS:HG3	1:A:504:LYS:HG3	1.92	0.52
1:A:283:HIS:ND1	1:A:283:HIS:O	2.43	0.51
1:A:69:ILE:HD12	1:A:69:ILE:H	1.75	0.51
1:A:230:GLN:HB2	1:A:288:VAL:HG23	1.93	0.51
1:A:66:GLN:HE21	1:A:111:MET:HB2	1.75	0.51
1:A:71:GLN:NE2	1:A:272:VAL:HG12	2.25	0.51
1:A:248:VAL:O	1:A:248:VAL:HG23	2.12	0.50
1:A:339:VAL:CG1	1:A:376:SER:HA	2.42	0.50
1:A:454:ARG:CZ	1:A:456:GLN:NE2	2.75	0.50
1:A:374:ASP:CG	1:A:377:GLN:HB2	2.30	0.50
1:A:364:PHE:CZ	1:A:368:THR:HG21	2.47	0.49
1:A:110:LEU:HD12	1:A:146:GLY:HA2	1.94	0.49
1:A:380:LYS:O	1:A:384:VAL:HG23	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:31:LYS:CA	1:A:84:ASN:HD22	2.17	0.49
1:A:113:SER:OG	1:A:116:GLY:HA2	2.12	0.49
1:A:300:PRO:HA	1:A:303:LEU:HD12	1.95	0.49
1:A:446:PRO:O	1:A:450:PRO:HA	2.13	0.49
1:A:31:LYS:HA	1:A:84:ASN:ND2	2.20	0.49
1:A:513:SER:HB3	1:A:516:THR:CG2	2.42	0.48
1:A:273:PRO:HG2	1:A:284:TYR:HE2	1.78	0.48
1:A:273:PRO:HB3	1:A:277:LEU:HD12	1.94	0.48
1:A:372:ALA:C	1:A:374:ASP:N	2.64	0.48
1:A:420:HIS:CG	1:A:460:VAL:HG21	2.48	0.48
1:A:322:HIS:HB3	1:A:434:ASP:OD2	2.14	0.48
1:A:370:SER:O	1:A:371:TRP:C	2.50	0.48
1:A:69:ILE:HD12	1:A:69:ILE:N	2.29	0.48
1:A:238:LYS:O	1:A:242:GLU:HG3	2.14	0.47
1:A:529:LEU:HG	1:A:530:PRO:HD2	1.96	0.47
1:A:31:LYS:NZ	2:A:670:HOH:O	2.46	0.47
1:A:273:PRO:CB	1:A:277:LEU:HD12	2.44	0.47
1:A:487:HIS:CE1	1:A:489:GLU:OE1	2.63	0.47
1:A:441:TYR:CB	1:A:464:MET:CE	2.90	0.47
1:A:418:PHE:CE1	1:A:464:MET:CE	2.98	0.47
1:A:63:ARG:CB	1:A:113:SER:HB3	2.45	0.47
1:A:379:ASN:O	1:A:382:LYS:N	2.48	0.47
1:A:79:ASP:O	1:A:82:TYR:HE1	1.98	0.46
1:A:317:ASN:HD22	1:A:418:PHE:HB3	1.79	0.46
1:A:344:PHE:CE1	1:A:348:VAL:HG11	2.50	0.46
1:A:485:PRO:HG3	1:A:505:MET:HG3	1.97	0.46
1:A:261:THR:HG22	1:A:262:ASP:N	2.31	0.46
1:A:240:VAL:HG13	1:A:269:ALA:HB2	1.94	0.46
1:A:304:TYR:CD1	1:A:403:GLN:HG2	2.51	0.46
1:A:234:LEU:O	1:A:235:PHE:C	2.55	0.45
1:A:322:HIS:HD2	2:A:675:HOH:O	1.99	0.45
1:A:325:ALA:HB1	1:A:332:ILE:HD11	1.98	0.45
1:A:241:ALA:HB3	1:A:248:VAL:HG12	1.98	0.45
1:A:71:GLN:HE22	1:A:272:VAL:N	2.13	0.45
1:A:94:VAL:HG12	2:A:546:HOH:O	2.16	0.45
1:A:298:ASP:OD2	1:A:299:ASP:N	2.45	0.45
1:A:355:LYS:HB3	1:A:359:GLY:HA3	1.98	0.45
1:A:279:TYR:CB	1:A:280:PRO:CD	2.95	0.45
1:A:454:ARG:NH1	1:A:456:GLN:NE2	2.64	0.45
1:A:513:SER:CB	1:A:516:THR:HG23	2.47	0.45
1:A:378:GLU:HG3	1:A:427:TYR:CD2	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:89:GLN:HA	1:A:89:GLN:OE1	2.17	0.45
1:A:447:PHE:CE2	1:A:458:ARG:HG2	2.52	0.45
1:A:69:ILE:H	1:A:69:ILE:CD1	2.29	0.45
1:A:279:TYR:CE2	1:A:283:HIS:CE1	3.05	0.45
1:A:340:THR:CG2	1:A:341:GLU:N	2.79	0.45
1:A:273:PRO:HG2	1:A:284:TYR:CE2	2.52	0.44
1:A:273:PRO:HB3	1:A:277:LEU:HD11	1.97	0.44
1:A:355:LYS:O	1:A:356:GLY:O	2.35	0.44
1:A:426:VAL:CG2	1:A:427:TYR:N	2.80	0.44
1:A:429:LYS:C	1:A:431:VAL:H	2.21	0.44
1:A:101:MET:HA	1:A:189:THR:O	2.18	0.43
1:A:421:PRO:HA	1:A:432:GLY:O	2.18	0.43
1:A:520:ARG:HG3	1:A:520:ARG:HH11	1.82	0.43
1:A:103:TRP:HA	1:A:191:PHE:O	2.19	0.43
1:A:325:ALA:HB2	1:A:388:GLU:HG3	2.00	0.43
1:A:119:PHE:HB3	1:A:120:LEU:H	1.72	0.43
1:A:3:LEU:HG	1:A:89:GLN:NE2	2.34	0.43
1:A:207:PRO:O	1:A:210:LYS:HG3	2.19	0.43
1:A:503:LYS:HG2	2:A:544:HOH:O	2.19	0.43
1:A:445:LYS:N	1:A:446:PRO:CD	2.81	0.43
1:A:32:GLY:N	1:A:84:ASN:HD22	2.16	0.42
1:A:379:ASN:O	1:A:381:LYS:N	2.52	0.42
1:A:395:VAL:HB	1:A:396:PRO:HD3	2.00	0.42
1:A:63:ARG:HB2	1:A:113:SER:HB3	2.01	0.42
1:A:117:ALA:CB	1:A:124:LEU:HB3	2.36	0.42
1:A:121:ASN:CG	1:A:122:ASN:N	2.72	0.42
1:A:345:TYR:O	1:A:349:SER:HB2	2.19	0.42
1:A:234:LEU:HD11	1:A:238:LYS:CE	2.47	0.42
1:A:288:VAL:HB	1:A:289:PRO:HD2	2.01	0.42
1:A:370:SER:O	1:A:372:ALA:N	2.52	0.42
1:A:527:LEU:C	1:A:527:LEU:HD23	2.40	0.42
1:A:68:THR:OG1	1:A:69:ILE:HD12	2.20	0.42
1:A:102:ILE:HD11	1:A:174:VAL:HG11	2.01	0.42
1:A:152:SER:HB2	1:A:160:GLY:HA3	2.02	0.42
1:A:426:VAL:CG2	1:A:427:TYR:H	2.31	0.42
1:A:9:GLU:HG3	1:A:10:GLY:N	2.34	0.42
1:A:368:THR:O	1:A:380:LYS:HE3	2.20	0.42
1:A:117:ALA:HB3	1:A:124:LEU:CB	2.35	0.41
1:A:165:ARG:O	1:A:168:HIS:HB3	2.20	0.41
1:A:515:ARG:C	1:A:517:ASN:N	2.72	0.41
1:A:277:LEU:HB3	1:A:278:GLU:H	1.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:229:ILE:HG13	2:A:613:HOH:O	2.20	0.41
1:A:233:PRO:HG3	1:A:290:VAL:HG21	2.01	0.41
1:A:325:ALA:HB1	1:A:332:ILE:CD1	2.49	0.41
1:A:76:GLY:O	1:A:77:ASP:HB2	2.20	0.41
1:A:491:TYR:CD2	1:A:491:TYR:C	2.93	0.41
1:A:30:PHE:CD1	1:A:30:PHE:N	2.88	0.41
1:A:193:GLU:HA	1:A:219:GLN:O	2.20	0.41
1:A:374:ASP:HB2	1:A:379:ASN:CB	2.51	0.41
1:A:440:GLN:NE2	1:A:453:TYR:OH	2.53	0.41
1:A:372:ALA:HB1	1:A:375:PRO:CD	2.51	0.41
1:A:3:LEU:HD23	1:A:94:VAL:HG11	2.02	0.41
1:A:241:ALA:CB	1:A:248:VAL:HG12	2.51	0.41
1:A:37:ALA:HB2	1:A:47:PRO:O	2.21	0.40
1:A:279:TYR:HB2	1:A:280:PRO:CD	2.51	0.40
1:A:279:TYR:CD2	1:A:284:TYR:CZ	3.09	0.40
1:A:374:ASP:O	1:A:375:PRO:C	2.60	0.40
1:A:489:GLU:HA	1:A:490:PRO:HD3	1.87	0.40
1:A:6:VAL:CG1	1:A:180:ALA:HB1	2.49	0.40
1:A:15:GLY:O	1:A:58:LYS:NZ	2.55	0.40
1:A:232:ASN:HD21	1:A:235:PHE:HB2	1.84	0.40
1:A:272:VAL:O	1:A:273:PRO:O	2.39	0.40
1:A:435:HIS:O	1:A:436:ALA:CB	2.70	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	531/533 (100%)	460 (87%)	60 (11%)	11 (2%)	7 5

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	273	PRO
1	A	274	LEU
1	A	116	GLY
1	A	123	TYR
1	A	356	GLY
1	A	369	GLU
1	A	378	GLU
1	A	371	TRP
1	A	119	PHE
1	A	494	GLU
1	A	69	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	437/437 (100%)	404 (92%)	33 (8%)	13 16

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	LEU
1	A	22	LEU
1	A	63	ARG
1	A	68	THR
1	A	69	ILE
1	A	72	ASP
1	A	75	TYR
1	A	151	LEU
1	A	191	PHE
1	A	194	SER
1	A	214	ARG
1	A	215	ARG
1	A	239	LYS
1	A	256	GLN
1	A	279	TYR
1	A	290	VAL

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Mol	Chain	Res	Type
1	A	318	ASN
1	A	320	ASP
1	A	334	LYS
1	A	341	GLU
1	A	346	LYS
1	A	380	LYS
1	A	383	THR
1	A	401	LEU
1	A	405	ARG
1	A	417	LEU
1	A	429	LYS
1	A	451	THR
1	A	468	TRP
1	A	486	THR
1	A	495	ASN
1	A	517	ASN
1	A	525	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	GLN
1	A	71	GLN
1	A	84	ASN
1	A	118	ASN
1	A	177	ASN
1	A	232	ASN
1	A	306	ASN
1	A	317	ASN
1	A	318	ASN
1	A	440	GLN
1	A	487	HIS

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)

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, 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	533/533 (100%)	0.34	37 (6%) 16 22	22, 49, 104, 124	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	119	PHE	10.0
1	A	118	ASN	9.7
1	A	120	LEU	9.2
1	A	122	ASN	7.7
1	A	275	ALA	7.4
1	A	276	GLY	6.2
1	A	371	TRP	5.5
1	A	532	VAL	5.4
1	A	274	LEU	5.3
1	A	375	PRO	5.1
1	A	117	ALA	4.3
1	A	277	LEU	4.1
1	A	279	TYR	3.7
1	A	121	ASN	3.6
1	A	284	TYR	3.5
1	A	357	LEU	3.4
1	A	116	GLY	3.4
1	A	123	TYR	3.3
1	A	429	LYS	3.3
1	A	533	THR	3.1
1	A	376	SER	3.0
1	A	124	LEU	2.9
1	A	513	SER	2.9
1	A	273	PRO	2.8
1	A	24	GLY	2.7
1	A	527	LEU	2.7
1	A	373	GLN	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	138	VAL	2.6
1	A	528	ALA	2.6
1	A	377	GLN	2.4
1	A	151	LEU	2.4
1	A	264	ARG	2.4
1	A	1	ALA	2.3
1	A	354	THR	2.3
1	A	93	GLN	2.2
1	A	147	PRO	2.1
1	A	372	ALA	2.0

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)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.