



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 13, 2018 – 02:48 PM EDT

PDB ID : 6AUN  
Title : calcium-independent phospholipase A2 beta  
Authors : Malley, K.; Koroleva, O.; Miller, I.; Sanishvili, R.; Jenkins, C.M.; Gross, R.W.;  
Korolev, S.  
Deposited on : 2017-09-01  
Resolution : 3.95 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](mailto: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.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20030736  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
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) : rb-20030736

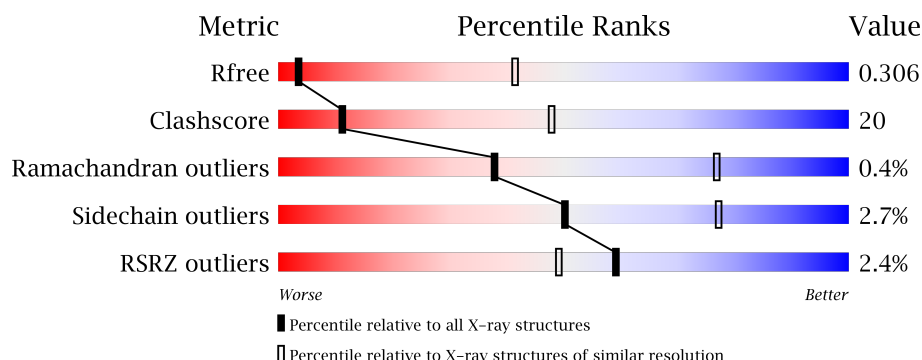
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.95 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))
$R_{free}$	100719	1057 (4.32-3.60)
Clashscore	112137	1154 (4.32-3.60)
Ramachandran outliers	110173	1111 (4.32-3.60)
Sidechain outliers	110143	1101 (4.32-3.60)
RSRZ outliers	101464	1069 (4.32-3.60)

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  $\geq 3$ , 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	752	<div> <div> <div>0.1%</div> <div>49%</div> <div>32%</div> <div>18%</div> </div> </div>
1	B	752	<div> <div>3%</div> <div>49%</div> <div>32%</div> <div>18%</div> </div>

## 2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 9396 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.

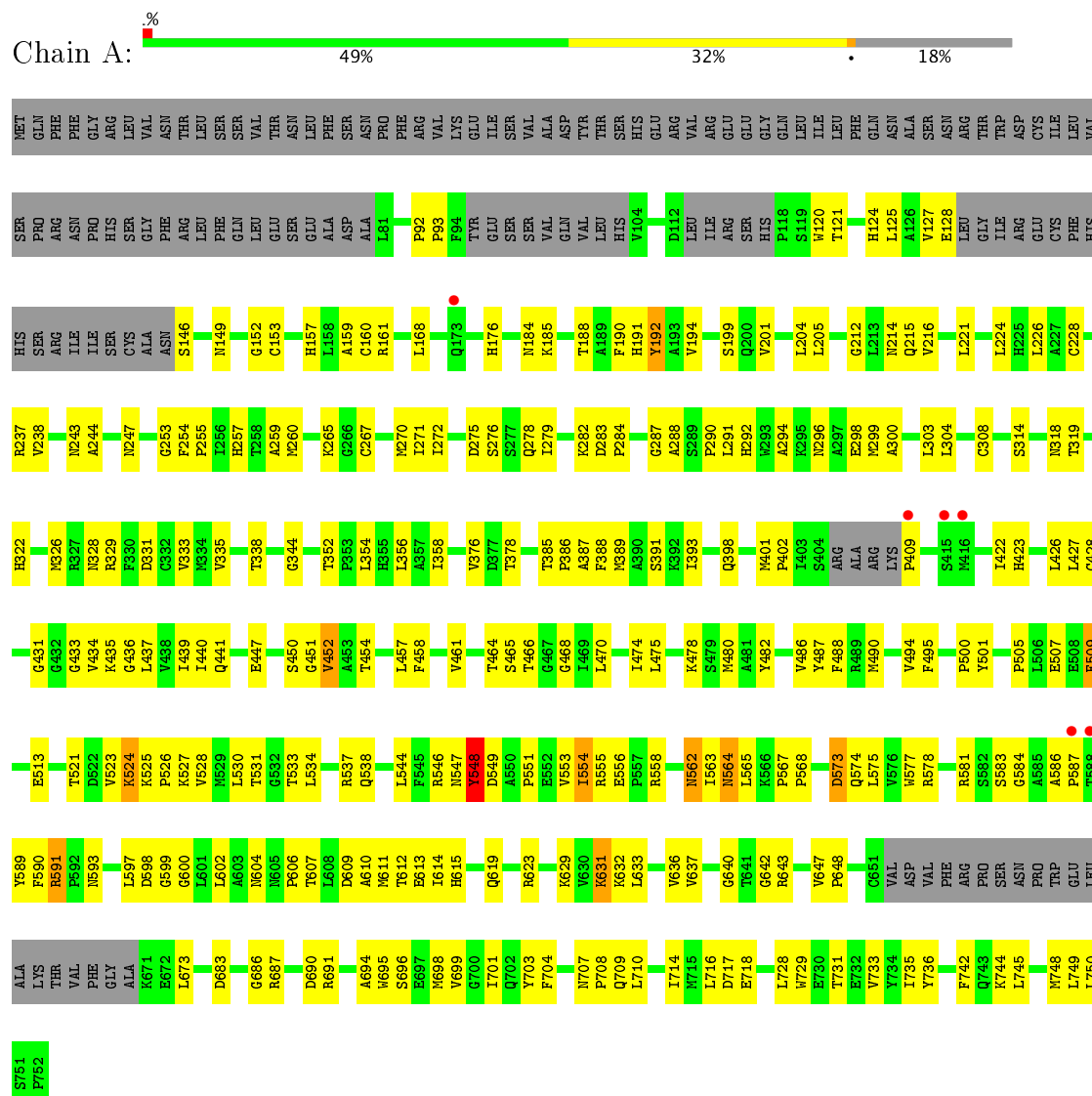
- Molecule 1 is a protein called PLA2G6, iPLA2beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	618	Total	C	N	O	S	0	0	0
			4698	2942	835	882	39			
1	B	618	Total	C	N	O	S	0	0	0
			4698	2942	835	882	39			

### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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: PLA2G6, iPLA2beta



#### • Molecule 1: PLA2G6, iPLA2beta





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 62	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	266.06 Å   266.06 Å   79.38 Å 90.00°   90.00°   120.00°	Depositor
Resolution (Å)	39.11 – 3.95 39.11 – 3.95	Depositor EDS
% Data completeness (in resolution range)	86.4 (39.11-3.95) 80.6 (39.11-3.95)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	5.14 (at 4.00 Å)	Xtriage
Refinement program	PHENIX (1.12_2829: ???)	Depositor
R, $R_{free}$	0.257 , 0.310 0.248 , 0.306	Depositor DCC
$R_{free}$ test set	2290 reflections (9.96%)	DCC
Wilson B-factor (Å <sup>2</sup> )	33.6	Xtriage
Anisotropy	0.417	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 5.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.44$ , $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	0.058 for h,-h-k,-l	Xtriage
$F_o, F_c$ correlation	0.79	EDS
Total number of atoms	9396	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.78% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	0.30	0/4782	0.60	3/6465 (0.0%)
1	B	0.32	0/4782	0.60	3/6465 (0.0%)
All	All	0.31	0/9564	0.60	6/12930 (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	6
1	B	0	1
All	All	0	7

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	409	PRO	N-CA-CB	5.93	110.42	103.30
1	B	409	PRO	N-CA-CB	5.91	110.39	103.30
1	A	93	PRO	N-CA-CB	5.88	110.36	103.30
1	B	93	PRO	N-CA-CB	5.71	110.15	103.30
1	A	92	PRO	N-CA-CB	5.70	110.13	103.30
1	B	92	PRO	N-CA-CB	5.47	109.87	103.30

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	401	MET	Peptide
1	A	452	VAL	Peptide
1	A	548	TYR	Peptide

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Mol	Chain	Res	Type	Group
1	A	562	ASN	Peptide
1	A	591	ARG	Peptide
1	A	631	LYS	Peptide
1	B	564	ASN	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	4698	0	4595	192	0
1	B	4698	0	4595	197	0
All	All	9396	0	9190	370	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 20.

All (370) 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:464:THR:HG21	1:A:604:ASN:HA	1.53	0.90
1:B:464:THR:HG21	1:B:604:ASN:HA	1.56	0.88
1:A:562:ASN:O	1:A:564:ASN:N	2.07	0.86
1:A:288:ALA:HB1	1:A:292:HIS:HB2	1.58	0.84
1:B:573:ASP:O	1:B:574:GLN:NE2	2.13	0.81
1:B:260:MET:HB3	1:B:294:ALA:HB1	1.62	0.81
1:B:295:LYS:HE3	1:B:327:ARG:HG2	1.60	0.81
1:B:575:LEU:HD13	1:B:578:ARG:HE	1.43	0.81
1:A:260:MET:HB3	1:A:294:ALA:HB1	1.62	0.80
1:A:691:ARG:HD3	1:B:608:LEU:HD12	1.63	0.80
1:A:642:GLY:HA2	1:A:708:PRO:HD2	1.62	0.79
1:B:587:PRO:HD3	1:B:598:ASP:HB2	1.62	0.79
1:B:265:LYS:NZ	1:B:298:GLU:OE1	2.17	0.77
1:B:168:LEU:HD22	1:B:201:VAL:HG22	1.67	0.77
1:B:588:THR:HG21	1:B:678:VAL:HG11	1.67	0.77
1:A:548:TYR:OH	1:A:613:GLU:OE2	2.04	0.74
1:B:643:ARG:NH1	1:B:709:GLN:OE1	2.20	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:604:ASN:ND2	1:B:682:THR:O	2.21	0.72
1:A:586:ALA:HA	1:A:598:ASP:HB2	1.72	0.72
1:A:607:THR:O	1:A:611:MET:HG3	1.91	0.70
1:A:607:THR:HG21	1:A:703:TYR:HE1	1.55	0.70
1:A:619:GLN:O	1:A:623:ARG:HB2	1.91	0.70
1:B:614:ILE:HG13	1:B:633:LEU:HD11	1.73	0.69
1:B:554:ILE:HG23	1:B:555:ARG:H	1.57	0.69
1:B:433:GLY:HA2	1:B:716:LEU:HG	1.74	0.69
1:A:643:ARG:NH1	1:A:709:GLN:OE1	2.26	0.69
1:A:600:GLY:HA2	1:A:604:ASN:HB2	1.72	0.69
1:A:587:PRO:HG3	1:A:597:LEU:HB3	1.76	0.68
1:B:120:TRP:HB3	1:B:125:LEU:HB2	1.74	0.68
1:B:600:GLY:HA2	1:B:604:ASN:HB2	1.75	0.68
1:B:433:GLY:CA	1:B:716:LEU:HG	2.24	0.68
1:B:548:TYR:HD1	1:B:549:ASP:H	1.40	0.67
1:A:146:SER:OG	1:A:176:HIS:O	2.10	0.67
1:A:615:HIS:HD2	1:B:694:ALA:HB1	1.59	0.67
1:B:437:LEU:HD21	1:B:487:TYR:HB2	1.76	0.67
1:A:647:VAL:HB	1:B:538:GLN:HB3	1.77	0.66
1:A:314:SER:HB3	1:A:318:ASN:H	1.60	0.66
1:A:427:LEU:HB3	1:A:461:VAL:HG22	1.75	0.66
1:B:198:ASN:OD1	1:B:201:VAL:N	2.18	0.66
1:A:464:THR:O	1:A:583:SER:OG	2.13	0.66
1:A:276:SER:O	1:A:279:ILE:HG12	1.96	0.66
1:B:127:VAL:HG13	1:B:167:ILE:HD13	1.76	0.66
1:B:296:ASN:HB3	1:B:299:MET:HG3	1.77	0.66
1:B:587:PRO:HD3	1:B:598:ASP:CB	2.26	0.66
1:A:696:SER:HB3	1:A:701:ILE:HB	1.79	0.65
1:A:282:LYS:HA	1:A:288:ALA:O	1.97	0.65
1:A:436:GLY:HA3	1:A:466:THR:HG21	1.79	0.65
1:B:282:LYS:HA	1:B:288:ALA:O	1.97	0.64
1:A:609:ASP:OD1	1:B:687:ARG:NE	2.28	0.63
1:A:149:ASN:HB3	1:A:153:CYS:H	1.64	0.63
1:A:300:ALA:O	1:A:304:LEU:HB2	1.99	0.63
1:A:687:ARG:NH2	1:B:609:ASP:OD1	2.32	0.62
1:A:501:TYR:OH	1:A:717:ASP:OD2	2.17	0.62
1:B:163:GLY:HA2	1:B:198:ASN:HD22	1.65	0.62
1:A:744:LYS:O	1:A:748:MET:HG3	2.00	0.62
1:B:289:SER:HB2	1:B:292:HIS:ND1	2.15	0.62
1:B:642:GLY:HA2	1:B:708:PRO:HD2	1.81	0.62
1:A:121:THR:H	1:A:124:HIS:HB3	1.63	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:322:HIS:HD2	1:B:356:LEU:HD12	1.63	0.62
1:B:524:LYS:HA	1:B:547:ASN:HD21	1.64	0.62
1:A:468:GLY:HA2	1:A:530:LEU:HD22	1.82	0.61
1:B:738:HIS:O	1:B:740:GLU:N	2.34	0.61
1:B:291:LEU:HD22	1:B:308:CYS:SG	2.40	0.61
1:A:212:GLY:HA2	1:A:215:GLN:HB2	1.83	0.61
1:B:542:LEU:HD13	1:B:602:LEU:HD11	1.83	0.61
1:A:565:LEU:HD11	1:B:648:PRO:HD3	1.83	0.60
1:B:633:LEU:O	1:B:701:ILE:HD13	2.02	0.60
1:A:161:ARG:HH11	1:A:192:TYR:HE1	1.47	0.60
1:A:587:PRO:HD3	1:A:598:ASP:HB2	1.81	0.60
1:A:615:HIS:CD2	1:B:694:ALA:HB1	2.36	0.60
1:A:548:TYR:HD1	1:A:549:ASP:H	1.49	0.60
1:B:455:LYS:NZ	1:B:477:SER:OG	2.34	0.60
1:B:533:THR:OG1	1:B:598:ASP:O	2.15	0.60
1:A:283:ASP:O	1:A:287:GLY:N	2.35	0.59
1:B:220:GLY:O	1:B:251:PRO:HD3	2.01	0.59
1:B:566:LYS:C	1:B:568:PRO:HD3	2.22	0.59
1:B:587:PRO:HG3	1:B:597:LEU:HB3	1.82	0.59
1:A:447:GLU:OE1	1:A:454:THR:OG1	2.10	0.59
1:A:648:PRO:HG2	1:B:565:LEU:HD23	1.85	0.59
1:B:696:SER:HB3	1:B:701:ILE:HB	1.83	0.59
1:B:163:GLY:HA2	1:B:198:ASN:HB2	1.85	0.58
1:B:319:THR:OG1	1:B:322:HIS:ND1	2.35	0.58
1:A:553:VAL:C	1:A:555:ARG:H	2.07	0.58
1:A:731:THR:O	1:A:735:ILE:HG12	2.03	0.58
1:A:120:TRP:HB3	1:A:125:LEU:HB2	1.86	0.57
1:B:744:LYS:O	1:B:748:MET:HG2	2.04	0.57
1:B:157:HIS:HB3	1:B:192:TYR:CD2	2.40	0.57
1:B:240:LEU:HB3	1:B:274:MET:HG2	1.86	0.57
1:B:527:LYS:HB3	1:B:548:TYR:HD2	1.70	0.57
1:A:422:ILE:O	1:A:631:LYS:HB2	2.05	0.56
1:B:468:GLY:HA2	1:B:530:LEU:HD22	1.87	0.56
1:A:214:ASN:O	1:A:216:VAL:HG12	2.05	0.56
1:A:581:ARG:NH2	1:A:593:ASN:OD1	2.36	0.56
1:B:521:THR:HG22	1:B:547:ASN:HA	1.86	0.56
1:B:527:LYS:HD2	1:B:548:TYR:HB3	1.87	0.56
1:A:282:LYS:HB3	1:A:287:GLY:HA2	1.86	0.56
1:A:695:TRP:HB3	1:B:695:TRP:CZ3	2.41	0.55
1:B:313:THR:HG23	1:B:314:SER:H	1.69	0.55
1:B:694:ALA:O	1:B:698:MET:HG2	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:427:LEU:HB2	1:B:458:PHE:CD2	2.41	0.55
1:A:435:LYS:HD3	1:A:714:ILE:HB	1.88	0.55
1:A:610:ALA:O	1:A:614:ILE:HG12	2.07	0.55
1:A:607:THR:HG21	1:A:703:TYR:CE1	2.40	0.55
1:A:525:LYS:HB2	1:A:526:PRO:CD	2.36	0.55
1:A:427:LEU:HB2	1:A:458:PHE:CD2	2.42	0.55
1:A:694:ALA:O	1:A:698:MET:HG2	2.06	0.55
1:B:300:ALA:O	1:B:304:LEU:HB2	2.07	0.54
1:B:436:GLY:O	1:B:439:ILE:HG12	2.06	0.54
1:B:738:HIS:C	1:B:740:GLU:H	2.09	0.54
1:B:436:GLY:HA3	1:B:466:THR:HG21	1.88	0.54
1:A:475:LEU:HD22	1:A:528:VAL:HG21	1.88	0.54
1:A:440:ILE:HD11	1:A:487:TYR:CE2	2.43	0.54
1:A:573:ASP:O	1:A:574:GLN:NE2	2.40	0.54
1:A:391:SER:HA	1:A:398:GLN:HB2	1.89	0.54
1:B:523:VAL:HG12	1:B:525:LYS:HG3	1.89	0.54
1:B:531:THR:OG1	1:B:599:GLY:HA3	2.07	0.54
1:A:558:ARG:NH2	1:B:689:VAL:HG11	2.23	0.54
1:B:291:LEU:H	1:B:291:LEU:HD23	1.72	0.53
1:A:523:VAL:O	1:A:524:LYS:HG2	2.08	0.53
1:B:704:PHE:CD2	1:B:745:LEU:HD13	2.43	0.53
1:B:222:THR:HG22	1:B:223:PRO:HD2	1.90	0.53
1:A:470:LEU:O	1:A:474:ILE:HG23	2.08	0.53
1:A:478:LYS:HE3	1:A:513:GLU:O	2.09	0.53
1:B:289:SER:HB3	1:B:291:LEU:HD23	1.91	0.53
1:A:558:ARG:HH22	1:B:689:VAL:HG11	1.74	0.53
1:A:612:THR:OG1	1:B:691:ARG:HA	2.09	0.53
1:B:154:THR:HB	1:B:157:HIS:CG	2.43	0.53
1:B:642:GLY:HA3	1:B:710:LEU:HD11	1.91	0.53
1:A:426:LEU:HB3	1:A:636:VAL:HG22	1.91	0.52
1:A:125:LEU:HD12	1:A:128:GLU:HB3	1.92	0.52
1:B:260:MET:HB3	1:B:294:ALA:CB	2.36	0.52
1:B:476:HIS:CE1	1:B:523:VAL:HG21	2.43	0.52
1:A:188:THR:HG22	1:A:191:HIS:ND1	2.25	0.52
1:B:429:LEU:HD12	1:B:471:ALA:HB2	1.91	0.52
1:A:127:VAL:HG11	1:A:159:ALA:HA	1.92	0.52
1:B:524:LYS:HA	1:B:547:ASN:ND2	2.25	0.52
1:A:260:MET:HB3	1:A:294:ALA:CB	2.36	0.52
1:B:639:LEU:HD23	1:B:706:LEU:HB2	1.91	0.52
1:A:296:ASN:HB3	1:A:299:MET:HG3	1.92	0.52
1:A:695:TRP:CZ3	1:B:695:TRP:HB3	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:482:TYR:O	1:A:486:VAL:HG23	2.10	0.51
1:B:295:LYS:HD2	1:B:329:ARG:HD3	1.92	0.51
1:A:647:VAL:HB	1:B:538:GLN:CB	2.39	0.51
1:A:157:HIS:HB3	1:A:192:TYR:CD2	2.45	0.51
1:A:699:VAL:HG21	1:B:698:MET:HE2	1.93	0.51
1:B:629:LYS:HZ1	1:B:631:LYS:HE3	1.74	0.51
1:B:464:THR:O	1:B:583:SER:OG	2.24	0.51
1:A:637:VAL:HG22	1:A:704:PHE:HB2	1.93	0.51
1:A:565:LEU:HB3	1:A:567:PRO:HD2	1.93	0.51
1:B:605:ASN:O	1:B:605:ASN:ND2	2.44	0.51
1:B:287:GLY:O	1:B:314:SER:HA	2.11	0.50
1:B:438:VAL:HG13	1:B:731:THR:OG1	2.12	0.50
1:B:704:PHE:HD2	1:B:745:LEU:HD13	1.75	0.50
1:A:224:LEU:HD21	1:A:244:ALA:CB	2.42	0.50
1:B:619:GLN:O	1:B:623:ARG:HB2	2.11	0.50
1:A:267:CYS:O	1:A:271:ILE:HG13	2.12	0.50
1:A:507:GLU:HG3	1:A:577:TRP:CZ3	2.46	0.50
1:B:119:SER:O	1:B:149:ASN:ND2	2.45	0.50
1:A:201:VAL:HA	1:A:204:LEU:HD12	1.94	0.50
1:A:553:VAL:O	1:A:555:ARG:N	2.45	0.50
1:B:152:GLY:HA3	1:B:185:LYS:HE2	1.94	0.50
1:A:161:ARG:NH1	1:A:192:TYR:HE1	2.10	0.49
1:A:614:ILE:HG21	1:A:633:LEU:HD13	1.94	0.49
1:B:525:LYS:HB2	1:B:526:PRO:CD	2.42	0.49
1:A:290:PRO:O	1:A:294:ALA:HB3	2.13	0.49
1:B:602:LEU:HD12	1:B:603:ALA:H	1.78	0.49
1:A:495:PHE:CE2	1:A:717:ASP:HA	2.48	0.49
1:B:216:VAL:HG21	1:B:222:THR:N	2.27	0.49
1:A:433:GLY:HA2	1:A:716:LEU:HG	1.93	0.49
1:A:304:LEU:HA	1:A:308:CYS:HB3	1.95	0.49
1:B:428:CYS:SG	1:B:607:THR:HA	2.53	0.49
1:B:505:PRO:O	1:B:509:PHE:HB3	2.13	0.49
1:A:436:GLY:O	1:A:439:ILE:HG12	2.11	0.49
1:B:122:VAL:O	1:B:126:ALA:HB2	2.13	0.48
1:B:457:LEU:HD11	1:B:749:LEU:HB3	1.95	0.48
1:A:694:ALA:HB1	1:B:615:HIS:CD2	2.48	0.48
1:B:495:PHE:CE2	1:B:717:ASP:HA	2.49	0.48
1:B:287:GLY:H	1:B:315:ALA:HB2	1.79	0.48
1:B:314:SER:HB3	1:B:317:GLY:H	1.77	0.48
1:B:475:LEU:HB2	1:B:528:VAL:HG21	1.95	0.48
1:A:575:LEU:HD13	1:A:578:ARG:HE	1.79	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:554:ILE:O	1:B:555:ARG:HG2	2.13	0.48
1:A:694:ALA:HB1	1:B:615:HIS:HD2	1.78	0.48
1:A:120:TRP:CG	1:A:125:LEU:HD13	2.49	0.48
1:A:441:GLN:OE1	1:A:728:LEU:HD13	2.14	0.48
1:B:188:THR:HG22	1:B:191:HIS:ND1	2.29	0.48
1:A:387:ALA:HB1	1:A:402:PRO:HG3	1.95	0.48
1:B:188:THR:HG22	1:B:191:HIS:CE1	2.49	0.48
1:A:216:VAL:HG23	1:A:221:LEU:C	2.34	0.47
1:A:389:MET:O	1:A:393:ILE:HG12	2.14	0.47
1:A:544:LEU:HB2	1:A:546:ARG:NH2	2.29	0.47
1:B:260:MET:HE1	1:B:290:PRO:HB3	1.96	0.47
1:B:321:LEU:CD2	1:B:341:ALA:HB1	2.44	0.47
1:B:402:PRO:HB2	1:B:736:TYR:OH	2.14	0.47
1:A:237:ARG:HG3	1:A:270:MET:SD	2.54	0.47
1:A:427:LEU:HD13	1:A:458:PHE:CE2	2.49	0.47
1:B:440:ILE:HG12	1:B:470:LEU:HD22	1.97	0.47
1:A:534:LEU:HD21	1:A:537:ARG:NE	2.29	0.47
1:B:529:MET:HA	1:B:545:PHE:O	2.14	0.47
1:B:643:ARG:HD2	1:B:709:GLN:OE1	2.14	0.47
1:B:232:LYS:HB3	1:B:235:MET:HG2	1.97	0.47
1:A:216:VAL:HG23	1:A:221:LEU:N	2.30	0.47
1:B:523:VAL:O	1:B:524:LYS:HG2	2.14	0.47
1:B:331:ASP:O	1:B:335:VAL:HG23	2.15	0.47
1:A:326:MET:HG3	1:A:356:LEU:HD22	1.97	0.47
1:B:163:GLY:CA	1:B:198:ASN:HD22	2.28	0.47
1:B:256:ILE:HD12	1:B:278:GLN:HB3	1.95	0.47
1:A:160:CYS:SG	1:A:205:LEU:HD11	2.55	0.47
1:A:447:GLU:O	1:A:451:GLY:N	2.45	0.47
1:A:554:ILE:HG23	1:A:556:GLU:HG2	1.97	0.47
1:B:470:LEU:O	1:B:474:ILE:HG23	2.14	0.47
1:A:322:HIS:HD2	1:A:356:LEU:HD12	1.80	0.47
1:A:329:ARG:HE	1:A:329:ARG:HB3	1.58	0.47
1:A:527:LYS:H	1:A:527:LYS:HG2	1.57	0.47
1:A:500:PRO:HA	1:A:589:TYR:O	2.15	0.47
1:A:631:LYS:HG3	1:A:632:LYS:HG2	1.97	0.47
1:B:607:THR:O	1:B:611:MET:HG3	2.15	0.47
1:B:434:VAL:HG11	1:B:718:GLU:O	2.14	0.47
1:A:291:LEU:HD23	1:A:303:LEU:HD23	1.98	0.46
1:B:447:GLU:OE1	1:B:454:THR:HG22	2.15	0.46
1:B:465:SER:HA	1:B:584:GLY:HA2	1.97	0.46
1:A:319:THR:H	1:A:319:THR:HG1	1.52	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:463:GLY:O	1:B:531:THR:HG23	2.16	0.46
1:A:507:GLU:HG3	1:A:577:TRP:HZ3	1.80	0.46
1:B:631:LYS:HG2	1:B:632:LYS:H	1.80	0.46
1:A:191:HIS:HB2	1:A:226:LEU:HD23	1.96	0.46
1:A:426:LEU:HD12	1:A:427:LEU:H	1.80	0.46
1:B:121:THR:H	1:B:124:HIS:HB3	1.79	0.46
1:B:465:SER:O	1:B:584:GLY:HA2	2.16	0.46
1:A:533:THR:HG23	1:A:602:LEU:HD23	1.97	0.46
1:A:544:LEU:HD23	1:A:544:LEU:HA	1.84	0.46
1:A:575:LEU:HA	1:A:575:LEU:HD23	1.73	0.46
1:A:704:PHE:CD2	1:A:745:LEU:HD13	2.51	0.46
1:B:228:CYS:SG	1:B:259:ALA:HB2	2.56	0.46
1:A:254:PHE:HB3	1:A:255:PRO:HD2	1.98	0.46
1:A:606:PRO:O	1:A:610:ALA:N	2.33	0.45
1:A:254:PHE:HB2	1:A:257:HIS:CE1	2.51	0.45
1:A:521:THR:HG22	1:A:547:ASN:HA	1.98	0.45
1:A:553:VAL:HG13	1:A:555:ARG:H	1.80	0.45
1:A:590:PHE:O	1:A:591:ARG:HG2	2.16	0.45
1:A:642:GLY:HA3	1:A:710:LEU:HD11	1.98	0.45
1:B:322:HIS:CD2	1:B:356:LEU:HD12	2.49	0.45
1:B:461:VAL:HG12	1:B:471:ALA:HB1	1.97	0.45
1:A:329:ARG:O	1:A:333:VAL:HG23	2.16	0.45
1:B:120:TRP:CG	1:B:125:LEU:HD13	2.51	0.45
1:B:317:GLY:O	1:B:347:GLY:HA2	2.16	0.45
1:B:708:PRO:HB3	1:B:734:TYR:CB	2.46	0.45
1:B:380:ASN:HD21	1:B:384:GLU:HB2	1.81	0.45
1:A:437:LEU:HD21	1:A:487:TYR:HB2	1.98	0.45
1:A:338:THR:HG23	1:A:488:PHE:HB3	1.97	0.45
1:A:433:GLY:CA	1:A:716:LEU:HG	2.47	0.45
1:B:363:MET:HA	1:B:366:ILE:HD12	1.99	0.45
1:A:168:LEU:HD22	1:A:201:VAL:HG13	1.99	0.45
1:A:440:ILE:HG23	1:A:480:MET:HE3	1.98	0.45
1:A:640:GLY:O	1:A:707:ASN:HB2	2.17	0.45
1:B:437:LEU:HD23	1:B:440:ILE:HD12	1.99	0.45
1:A:450:SER:O	1:A:452:VAL:HG23	2.17	0.45
1:B:188:THR:HG23	1:B:190:PHE:H	1.81	0.45
1:B:427:LEU:HD13	1:B:458:PHE:CE2	2.51	0.45
1:B:482:TYR:O	1:B:486:VAL:HG23	2.17	0.45
1:B:476:HIS:ND1	1:B:523:VAL:HG21	2.32	0.45
1:B:629:LYS:NZ	1:B:631:LYS:HG3	2.32	0.45
1:B:589:TYR:CD2	1:B:716:LEU:HD22	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:523:VAL:HG23	1:A:524:LYS:H	1.82	0.45
1:A:733:VAL:O	1:A:736:TYR:HB2	2.17	0.45
1:A:199:SER:HB2	1:A:238:VAL:HG21	1.99	0.45
1:A:683:ASP:OD2	1:A:686:GLY:HA3	2.17	0.45
1:B:169:VAL:O	1:B:173:GLN:HB3	2.17	0.45
1:A:188:THR:HG23	1:A:190:PHE:H	1.82	0.44
1:A:505:PRO:O	1:A:509:PHE:HB3	2.17	0.44
1:A:600:GLY:HA2	1:A:604:ASN:CB	2.42	0.44
1:A:629:LYS:HE2	1:A:631:LYS:HD2	1.99	0.44
1:B:325:VAL:HG22	1:B:365:MET:SD	2.58	0.44
1:B:295:LYS:CE	1:B:327:ARG:HG2	2.39	0.44
1:A:494:VAL:HA	1:A:505:PRO:HB2	1.99	0.44
1:B:354:LEU:O	1:B:358:ILE:HG13	2.17	0.44
1:B:458:PHE:HE1	1:B:749:LEU:HD21	1.83	0.44
1:A:385:THR:O	1:A:388:PHE:HB3	2.18	0.44
1:B:696:SER:CB	1:B:701:ILE:HB	2.47	0.44
1:B:729:TRP:O	1:B:733:VAL:HG23	2.18	0.44
1:B:216:VAL:HG21	1:B:221:LEU:C	2.38	0.44
1:B:565:LEU:HG	1:B:567:PRO:HD3	1.99	0.44
1:B:731:THR:O	1:B:735:ILE:HG13	2.17	0.44
1:B:224:LEU:HD21	1:B:244:ALA:HB3	2.00	0.44
1:B:276:SER:O	1:B:279:ILE:HG12	2.18	0.44
1:A:191:HIS:HA	1:A:194:VAL:HG22	1.99	0.44
1:A:458:PHE:HE1	1:A:749:LEU:HD21	1.82	0.44
1:A:494:VAL:HG22	1:A:505:PRO:HB2	2.00	0.44
1:A:452:VAL:HG21	1:A:750:LEU:HD21	2.00	0.44
1:B:585:ALA:O	1:B:598:ASP:HB2	2.17	0.44
1:A:344:GLY:HA2	1:A:352:THR:HB	2.00	0.43
1:A:567:PRO:N	1:A:568:PRO:CD	2.81	0.43
1:A:673:LEU:HD13	1:B:592:PRO:HG3	2.00	0.43
1:A:378:THR:O	1:A:386:PRO:HD3	2.18	0.43
1:A:457:LEU:HD11	1:A:749:LEU:HB3	1.99	0.43
1:A:376:VAL:HA	1:A:386:PRO:HG2	2.00	0.43
1:B:640:GLY:O	1:B:707:ASN:HB2	2.19	0.43
1:A:296:ASN:HA	1:A:329:ARG:NH2	2.34	0.43
1:A:553:VAL:C	1:A:555:ARG:N	2.72	0.43
1:B:602:LEU:HD12	1:B:603:ALA:N	2.33	0.43
1:A:548:TYR:HD1	1:A:549:ASP:N	2.14	0.43
1:B:151:GLU:O	1:B:185:LYS:HE3	2.18	0.43
1:B:502:GLU:HG2	1:B:591:ARG:NH1	2.33	0.43
1:A:465:SER:HA	1:A:584:GLY:HA2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:166:GLU:O	1:B:170:GLU:HG3	2.19	0.43
1:B:525:LYS:HB2	1:B:526:PRO:HD3	1.99	0.43
1:A:185:LYS:HA	1:A:185:LYS:HD3	1.87	0.43
1:B:218:LYS:HA	1:B:218:LYS:HD2	1.88	0.43
1:A:331:ASP:O	1:A:335:VAL:HG23	2.19	0.43
1:A:735:ILE:HD13	1:A:742:PHE:CZ	2.54	0.43
1:B:314:SER:OG	1:B:315:ALA:N	2.51	0.43
1:B:429:LEU:O	1:B:463:GLY:HA2	2.19	0.43
1:B:267:CYS:O	1:B:271:ILE:HG13	2.18	0.43
1:B:475:LEU:HD12	1:B:526:PRO:HD2	2.01	0.43
1:A:265:LYS:NZ	1:A:298:GLU:HB3	2.33	0.42
1:A:704:PHE:HZ	1:A:744:LYS:HE2	1.85	0.42
1:B:313:THR:OG1	1:B:317:GLY:HA2	2.19	0.42
1:A:201:VAL:O	1:A:205:LEU:HG	2.19	0.42
1:A:422:ILE:HG13	1:A:423:HIS:H	1.82	0.42
1:B:218:LYS:O	1:B:218:LYS:NZ	2.43	0.42
1:A:551:PRO:HG3	1:A:613:GLU:OE2	2.19	0.42
1:B:508:GLU:HA	1:B:511:LYS:HD2	2.02	0.42
1:A:428:CYS:SG	1:A:607:THR:HA	2.59	0.42
1:A:531:THR:OG1	1:A:599:GLY:HA3	2.20	0.42
1:B:162:LYS:HD3	1:B:162:LYS:HA	1.87	0.42
1:B:598:ASP:C	1:B:600:GLY:H	2.22	0.42
1:B:693:ARG:HB2	1:B:703:TYR:CD2	2.55	0.42
1:A:354:LEU:O	1:A:358:ILE:HG13	2.20	0.41
1:A:439:ILE:HD11	1:A:470:LEU:HD23	2.01	0.41
1:B:355:HIS:NE2	1:B:384:GLU:O	2.53	0.41
1:B:394:SER:HB3	1:B:397:LEU:CD2	2.50	0.41
1:B:426:LEU:HD12	1:B:427:LEU:H	1.85	0.41
1:A:729:TRP:O	1:A:733:VAL:HG23	2.21	0.41
1:B:120:TRP:CD2	1:B:125:LEU:HD13	2.56	0.41
1:B:386:PRO:O	1:B:389:MET:HB2	2.20	0.41
1:A:188:THR:HG22	1:A:191:HIS:CE1	2.55	0.41
1:A:319:THR:OG1	1:A:322:HIS:ND1	2.44	0.41
1:B:450:SER:OG	1:B:452:VAL:HG22	2.20	0.41
1:A:524:LYS:HA	1:A:547:ASN:ND2	2.35	0.41
1:B:329:ARG:O	1:B:333:VAL:HG23	2.21	0.41
1:A:212:GLY:CA	1:A:215:GLN:HB2	2.49	0.41
1:A:265:LYS:HA	1:A:299:MET:HE1	2.03	0.41
1:B:534:LEU:HD21	1:B:537:ARG:CZ	2.51	0.41
1:B:553:VAL:HG22	1:B:554:ILE:HA	2.01	0.41
1:A:228:CYS:SG	1:A:259:ALA:HB2	2.60	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:152:GLY:HA2	1:A:184:ASN:HB2	2.02	0.41
1:A:296:ASN:HA	1:A:329:ARG:HH22	1.85	0.41
1:A:434:VAL:HG11	1:A:718:GLU:O	2.21	0.41
1:A:224:LEU:HD21	1:A:244:ALA:HB3	2.02	0.41
1:A:431:GLY:HA2	1:A:465:SER:H	1.86	0.41
1:A:253:GLY:HA2	1:A:284:PRO:HD2	2.03	0.41
1:A:538:GLN:HB3	1:B:647:VAL:O	2.21	0.41
1:A:749:LEU:HD23	1:A:749:LEU:HA	1.81	0.41
1:B:321:LEU:HD23	1:B:341:ALA:HB1	2.02	0.41
1:B:553:VAL:HA	1:B:554:ILE:HA	1.73	0.41
1:B:596:PHE:O	1:B:597:LEU:HD23	2.21	0.41
1:B:527:LYS:HG2	1:B:527:LYS:H	1.68	0.40
1:A:554:ILE:O	1:A:554:ILE:HG13	2.21	0.40
1:B:232:LYS:HB3	1:B:235:MET:CG	2.51	0.40
1:B:544:LEU:HA	1:B:544:LEU:HD23	1.75	0.40
1:B:687:ARG:HB3	1:B:691:ARG:HE	1.86	0.40
1:B:491:LYS:HE2	1:B:721:ASP:HB3	2.03	0.40
1:A:188:THR:HG23	1:A:190:PHE:N	2.37	0.40
1:A:271:ILE:HG22	1:A:278:GLN:HG3	2.04	0.40
1:A:272:ILE:HA	1:A:275:ASP:O	2.21	0.40
1:A:524:LYS:HA	1:A:547:ASN:HD21	1.86	0.40
1:B:236:VAL:O	1:B:240:LEU:HG	2.22	0.40
1:B:430:ASP:N	1:B:430:ASP:OD1	2.52	0.40
1:B:565:LEU:HD12	1:B:565:LEU:HA	1.94	0.40
1:B:607:THR:CG2	1:B:703:TYR:HE1	2.35	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	606/752 (81%)	509 (84%)	94 (16%)	3 (0%)	32 73

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	606/752 (81%)	500 (82%)	104 (17%)	2 (0%)	44	80
All	All	1212/1504 (81%)	1009 (83%)	198 (16%)	5 (0%)	38	77

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	563	ILE
1	A	554	ILE
1	B	739	ARG
1	A	564	ASN
1	B	572	ALA

### 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	495/646 (77%)	485 (98%)	10 (2%)	60	83
1	B	495/646 (77%)	478 (97%)	17 (3%)	42	73
All	All	990/1292 (77%)	963 (97%)	27 (3%)	50	77

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	192	TYR
1	A	243	ASN
1	A	247	ASN
1	A	328	ASN
1	A	490	MET
1	A	509	PHE
1	A	524	LYS
1	A	548	TYR
1	A	573	ASP
1	A	690	ASP
1	B	192	TYR

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Mol	Chain	Res	Type
1	B	247	ASN
1	B	249	MET
1	B	286	TYR
1	B	291	LEU
1	B	328	ASN
1	B	424	ASP
1	B	490	MET
1	B	509	PHE
1	B	546	ARG
1	B	548	TYR
1	B	573	ASP
1	B	609	ASP
1	B	629	LYS
1	B	690	ASP
1	B	707	ASN
1	B	736	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	615	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 carbohydrates 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<sup>th</sup> 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 <sup>2</sup> )	Q < 0.9
1	A	618/752 (82%)	0.02	6 (0%) 82 74	22, 53, 106, 135	0
1	B	618/752 (82%)	0.25	24 (3%) 40 31	19, 68, 148, 182	0
All	All	1236/1504 (82%)	0.14	30 (2%) 59 49	19, 60, 142, 182	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	413	LEU	5.9
1	B	126	ALA	5.4
1	B	159	ALA	5.2
1	B	416	MET	4.9
1	B	415	SER	4.8
1	B	564	ASN	3.9
1	B	175	CYS	3.8
1	B	414	SER	3.7
1	A	409	PRO	3.6
1	B	164	ASP	3.4
1	B	246	CYS	3.2
1	B	189	ALA	2.9
1	A	587	PRO	2.9
1	B	127	VAL	2.8
1	B	563	ILE	2.8
1	B	412	ILE	2.8
1	B	166	GLU	2.7
1	B	165	SER	2.6
1	A	415	SER	2.6
1	B	168	LEU	2.5
1	B	176	HIS	2.5
1	B	174	TYR	2.4
1	B	409	PRO	2.4
1	A	588	THR	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	163	GLY	2.3
1	A	416	MET	2.3
1	A	173	GLN	2.3
1	B	221	LEU	2.1
1	B	167	ILE	2.1
1	B	128	GLU	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 carbohydrates 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.