



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 24, 2023 – 07:00 PM EDT

PDB ID : 3GEC
Title : Crystal structure of a tandem PAS domain fragment of Drosophila PERIOD
Authors : Yildiz, O.; Wolf, E.
Deposited on : 2009-02-25
Resolution : 4.00 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 : 2.36
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) : 2.36

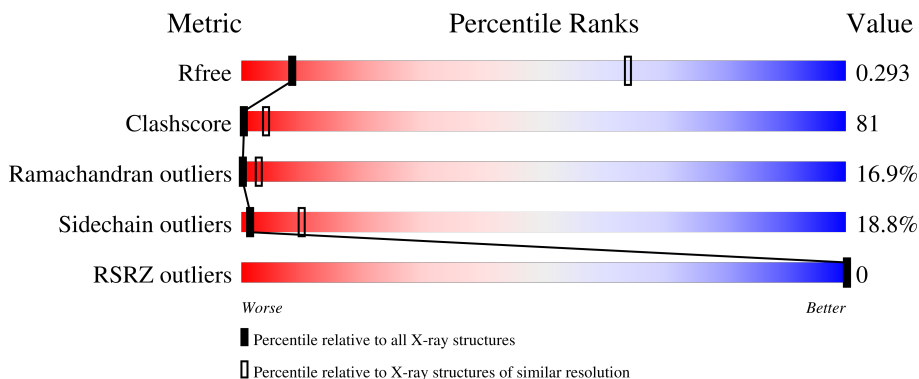
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.00 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	1087 (4.30-3.70)
Clashscore	141614	1148 (4.30-3.70)
Ramachandran outliers	138981	1108 (4.30-3.70)
Sidechain outliers	138945	1099 (4.30-3.70)
RSRZ outliers	127900	1028 (4.34-3.66)

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	312	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2054 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 Period circadian protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	257	2054	1315	351	373	15	0	0	0

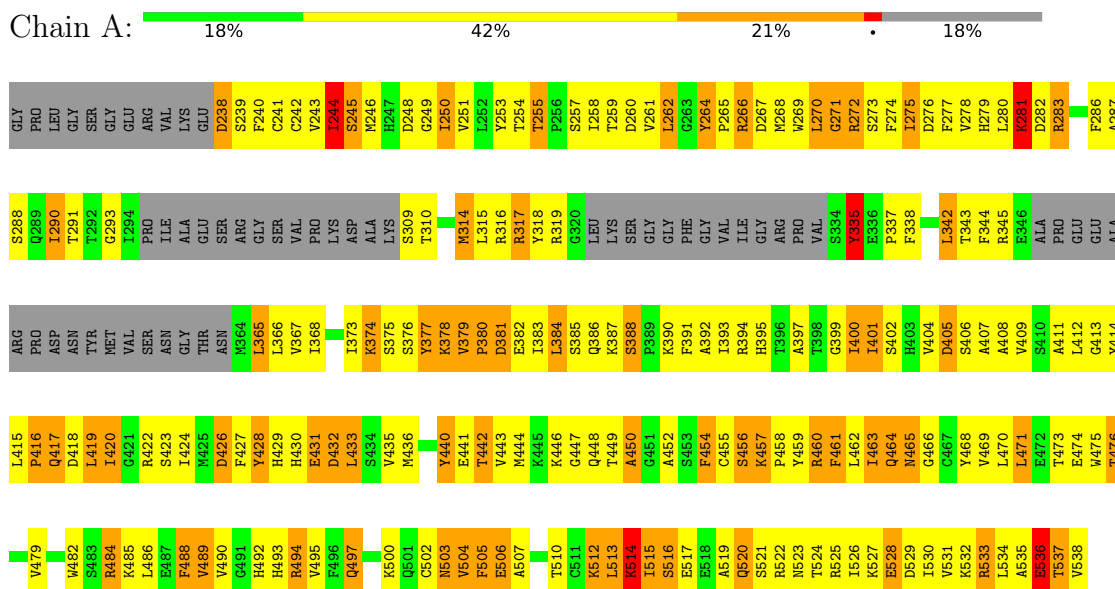
There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	227	GLY	-	expression tag	UNP P07663
A	228	PRO	-	expression tag	UNP P07663
A	229	LEU	-	expression tag	UNP P07663
A	230	GLY	-	expression tag	UNP P07663
A	231	SER	-	expression tag	UNP P07663

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: Period circadian protein



4 Data and refinement statistics i

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	114.95Å 114.95Å 85.70Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.68 – 4.00 19.68 – 4.00	Depositor EDS
% Data completeness (in resolution range)	97.7 (19.68-4.00) 97.6 (19.68-4.00)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.66 (at 4.06Å)	Xtrriage
Refinement program	CNS	Depositor
R, R_{free}	0.244 , 0.312 0.240 , 0.293	Depositor DCC
R_{free} test set	279 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	162.8	Xtrriage
Anisotropy	0.343	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 114.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtrriage
Estimated twinning fraction	0.078 for -h,-k,l	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	2054	wwPDB-VP
Average B, all atoms (Å ²)	148.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.46% 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 [i](#)

5.1 Standard geometry [i](#)

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	0.40	0/2100	0.75	1/2834 (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 (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	246	MET	N-CA-C	5.22	125.10	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	377	TYR	Sidechain

5.2 Too-close contacts [i](#)

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	2054	0	2054	333	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	2054	0	2054	333	1

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 81.

All (333) 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:506:GLU:HG3	1:A:507:ALA:H	1.17	1.09
1:A:314:MET:HE2	1:A:337:PRO:HA	1.39	1.03
1:A:243:VAL:O	1:A:244:ILE:HG23	1.59	1.03
1:A:250:ILE:HG22	1:A:251:VAL:H	1.22	1.01
1:A:440:TYR:O	1:A:443:VAL:HG12	1.59	0.99
1:A:309:SER:O	1:A:310:THR:HG22	1.67	0.94
1:A:280:LEU:HA	1:A:283:ARG:HD2	1.47	0.94
1:A:463:ILE:HG22	1:A:464:GLN:H	1.30	0.94
1:A:379:VAL:HB	1:A:380:PRO:CD	2.00	0.91
1:A:433:LEU:HA	1:A:436:MET:CB	2.02	0.90
1:A:254:THR:HG22	1:A:258:ILE:HG21	1.54	0.89
1:A:433:LEU:HA	1:A:436:MET:HB2	1.55	0.88
1:A:342:LEU:HB3	1:A:368:ILE:HD13	1.55	0.86
1:A:386:GLN:HA	1:A:494:ARG:NH1	1.93	0.83
1:A:264:TYR:HB2	1:A:268:MET:HG2	1.59	0.83
1:A:484:ARG:HG2	1:A:538:VAL:HG13	1.60	0.83
1:A:265:PRO:O	1:A:268:MET:HB3	1.77	0.83
1:A:317:ARG:HD3	1:A:319:ARG:HH22	1.44	0.83
1:A:513:LEU:O	1:A:514:LYS:HB3	1.79	0.81
1:A:519:ALA:HA	1:A:522:ARG:HB3	1.61	0.81
1:A:526:ILE:O	1:A:530:ILE:HG13	1.81	0.81
1:A:379:VAL:HB	1:A:380:PRO:HD2	1.62	0.81
1:A:259:THR:OG1	1:A:265:PRO:HA	1.81	0.81
1:A:430:HIS:CE1	1:A:431:GLU:HG3	2.16	0.81
1:A:267:ASP:HA	1:A:270:LEU:HD23	1.61	0.80
1:A:393:ILE:HG22	1:A:404:VAL:HG23	1.64	0.79
1:A:266:ARG:HG3	1:A:267:ASP:H	1.46	0.78
1:A:279:HIS:HA	1:A:335:TYR:HE2	1.47	0.78
1:A:413:GLY:N	1:A:504:VAL:HG11	1.99	0.77
1:A:342:LEU:HD13	1:A:342:LEU:H	1.49	0.77
1:A:337:PRO:HB2	1:A:373:ILE:CG2	2.15	0.76
1:A:502:CYS:SG	1:A:503:ASN:N	2.59	0.75
1:A:342:LEU:HB3	1:A:368:ILE:CD1	2.16	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:516:SER:HA	1:A:520:GLN:HB2	1.69	0.75
1:A:402:SER:O	1:A:420:ILE:HG13	1.86	0.74
1:A:494:ARG:HH21	1:A:494:ARG:HG3	1.51	0.74
1:A:342:LEU:HD13	1:A:342:LEU:N	2.02	0.74
1:A:272:ARG:H	1:A:272:ARG:HD3	1.51	0.74
1:A:265:PRO:HB2	1:A:268:MET:HB2	1.69	0.74
1:A:244:ILE:HG21	1:A:250:ILE:O	1.88	0.73
1:A:430:HIS:C	1:A:432:ASP:H	1.90	0.73
1:A:432:ASP:O	1:A:435:VAL:HG22	1.89	0.73
1:A:290:ILE:HD13	1:A:291:THR:N	2.04	0.72
1:A:378:LYS:HD3	1:A:379:VAL:HG22	1.71	0.72
1:A:250:ILE:HG22	1:A:251:VAL:N	2.01	0.72
1:A:414:TYR:CG	1:A:419:LEU:HD11	2.24	0.71
1:A:273:SER:HB3	1:A:275:ILE:HG22	1.71	0.71
1:A:314:MET:CE	1:A:337:PRO:HA	2.17	0.70
1:A:344:PHE:CD1	1:A:366:LEU:HD11	2.26	0.70
1:A:401:ILE:HD13	1:A:401:ILE:H	1.56	0.70
1:A:395:HIS:CD2	1:A:424:ILE:HG21	2.24	0.70
1:A:430:HIS:O	1:A:432:ASP:N	2.24	0.70
1:A:250:ILE:HG23	1:A:271:GLY:O	1.91	0.70
1:A:514:LYS:HG2	1:A:515:ILE:N	2.06	0.70
1:A:433:LEU:HA	1:A:436:MET:HB3	1.73	0.69
1:A:262:LEU:HD13	1:A:262:LEU:O	1.93	0.69
1:A:527:LYS:C	1:A:529:ASP:H	1.96	0.69
1:A:506:GLU:CG	1:A:507:ALA:H	1.99	0.68
1:A:506:GLU:HG3	1:A:507:ALA:N	2.00	0.68
1:A:335:TYR:HD1	1:A:335:TYR:H	1.40	0.68
1:A:525:ARG:O	1:A:529:ASP:HB2	1.94	0.68
1:A:392:ALA:N	1:A:405:ASP:OD2	2.27	0.68
1:A:422:ARG:CG	1:A:423:SER:H	2.07	0.67
1:A:250:ILE:HA	1:A:273:SER:N	2.11	0.66
1:A:430:HIS:ND1	1:A:431:GLU:N	2.44	0.66
1:A:343:THR:HG23	1:A:367:VAL:HB	1.78	0.66
1:A:432:ASP:OD1	1:A:460:ARG:HG2	1.96	0.65
1:A:413:GLY:HA3	1:A:504:VAL:CG1	2.26	0.65
1:A:413:GLY:HA3	1:A:504:VAL:HG12	1.79	0.65
1:A:412:LEU:HA	1:A:463:ILE:HG23	1.79	0.65
1:A:429:HIS:HB3	1:A:460:ARG:HG3	1.77	0.65
1:A:386:GLN:NE2	1:A:494:ARG:HH22	1.95	0.64
1:A:381:ASP:OD2	1:A:500:LYS:HA	1.97	0.64
1:A:386:GLN:HA	1:A:494:ARG:HH12	1.61	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:277:PHE:O	1:A:316:ARG:HG2	1.98	0.63
1:A:494:ARG:HG3	1:A:494:ARG:NH2	2.12	0.63
1:A:479:VAL:HG11	1:A:538:VAL:HG21	1.80	0.62
1:A:463:ILE:HD11	1:A:469:VAL:HG22	1.80	0.62
1:A:463:ILE:HG22	1:A:464:GLN:N	2.10	0.62
1:A:424:ILE:HG13	1:A:428:TYR:HE1	1.64	0.62
1:A:502:CYS:O	1:A:503:ASN:HB3	1.99	0.62
1:A:433:LEU:H	1:A:433:LEU:HD22	1.65	0.62
1:A:535:ALA:O	1:A:537:THR:HG22	1.99	0.62
1:A:429:HIS:CD2	1:A:460:ARG:CZ	2.83	0.62
1:A:463:ILE:HD11	1:A:469:VAL:CG2	2.29	0.61
1:A:379:VAL:O	1:A:502:CYS:HB2	2.00	0.61
1:A:461:PHE:HD2	1:A:462:LEU:N	1.98	0.61
1:A:243:VAL:C	1:A:244:ILE:HG12	2.21	0.61
1:A:395:HIS:HB2	1:A:400:ILE:O	2.01	0.61
1:A:429:HIS:HD2	1:A:460:ARG:NH1	1.98	0.61
1:A:428:TYR:CD2	1:A:436:MET:HG2	2.36	0.61
1:A:409:VAL:CG2	1:A:416:PRO:HD3	2.31	0.60
1:A:399:GLY:O	1:A:424:ILE:HG22	2.01	0.60
1:A:379:VAL:HG23	1:A:382:GLU:HB2	1.84	0.60
1:A:393:ILE:HD12	1:A:393:ILE:C	2.21	0.60
1:A:337:PRO:HB2	1:A:373:ILE:HG22	1.83	0.60
1:A:392:ALA:HB2	1:A:492:HIS:HA	1.84	0.59
1:A:433:LEU:HD11	1:A:527:LYS:HE3	1.84	0.59
1:A:462:LEU:HB3	1:A:468:TYR:CE1	2.36	0.59
1:A:243:VAL:O	1:A:244:ILE:CG2	2.45	0.59
1:A:243:VAL:HA	1:A:366:LEU:O	2.02	0.59
1:A:488:PHE:HD2	1:A:488:PHE:O	1.85	0.59
1:A:250:ILE:HA	1:A:272:ARG:C	2.24	0.58
1:A:344:PHE:CE1	1:A:366:LEU:HD11	2.38	0.58
1:A:243:VAL:CG1	1:A:253:TYR:H	2.15	0.58
1:A:417:GLN:HA	1:A:420:ILE:CG2	2.33	0.58
1:A:383:ILE:O	1:A:383:ILE:HG13	2.03	0.58
1:A:429:HIS:O	1:A:432:ASP:HB2	2.02	0.58
1:A:422:ARG:CG	1:A:423:SER:N	2.66	0.58
1:A:513:LEU:O	1:A:514:LYS:CB	2.51	0.58
1:A:514:LYS:O	1:A:516:SER:N	2.37	0.58
1:A:380:PRO:O	1:A:382:GLU:N	2.36	0.57
1:A:413:GLY:CA	1:A:504:VAL:HG11	2.33	0.57
1:A:430:HIS:C	1:A:432:ASP:N	2.56	0.57
1:A:395:HIS:HD2	1:A:399:GLY:HA2	1.68	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:264:TYR:H	1:A:264:TYR:HD2	1.50	0.57
1:A:404:VAL:O	1:A:404:VAL:HG13	2.04	0.57
1:A:479:VAL:HA	1:A:485:LYS:O	2.04	0.57
1:A:513:LEU:C	1:A:514:LYS:HE2	2.25	0.57
1:A:250:ILE:CA	1:A:273:SER:HA	2.35	0.57
1:A:534:LEU:C	1:A:534:LEU:HD12	2.26	0.56
1:A:255:THR:O	1:A:258:ILE:HG22	2.06	0.56
1:A:415:LEU:O	1:A:416:PRO:C	2.42	0.56
1:A:391:PHE:CD2	1:A:408:ALA:HB2	2.40	0.56
1:A:423:SER:O	1:A:426:ASP:HB2	2.05	0.56
1:A:264:TYR:HD2	1:A:264:TYR:N	2.03	0.56
1:A:476:THR:HG23	1:A:490:VAL:HB	1.87	0.56
1:A:527:LYS:O	1:A:529:ASP:N	2.38	0.56
1:A:444:MET:O	1:A:447:GLY:N	2.39	0.56
1:A:276:ASP:HA	1:A:283:ARG:NH2	2.21	0.55
1:A:244:ILE:O	1:A:245:SER:HB2	2.06	0.55
1:A:255:THR:HG22	1:A:257:SER:H	1.71	0.55
1:A:314:MET:HE1	1:A:415:LEU:HD11	1.89	0.55
1:A:464:GLN:HG2	1:A:505:PHE:HA	1.88	0.55
1:A:264:TYR:N	1:A:264:TYR:CD2	2.72	0.55
1:A:244:ILE:HD13	1:A:250:ILE:O	2.06	0.55
1:A:375:SER:OG	1:A:504:VAL:HG23	2.05	0.55
1:A:274:PHE:O	1:A:277:PHE:HB2	2.07	0.55
1:A:254:THR:O	1:A:255:THR:C	2.45	0.54
1:A:317:ARG:HD3	1:A:319:ARG:NH2	2.17	0.54
1:A:345:ARG:NH2	1:A:367:VAL:HG21	2.22	0.54
1:A:387:LYS:O	1:A:388:SER:C	2.46	0.54
1:A:463:ILE:HD11	1:A:469:VAL:CG1	2.38	0.54
1:A:279:HIS:HA	1:A:335:TYR:CE2	2.36	0.54
1:A:281:LYS:HB3	1:A:281:LYS:HZ2	1.73	0.54
1:A:446:LYS:HE3	1:A:454:PHE:CD1	2.42	0.54
1:A:536:GLU:O	1:A:538:VAL:N	2.41	0.54
1:A:269:TRP:O	1:A:271:GLY:N	2.40	0.54
1:A:377:TYR:OH	1:A:411:ALA:HB1	2.07	0.54
1:A:482:TRP:C	1:A:484:ARG:H	2.11	0.54
1:A:251:VAL:HG13	1:A:272:ARG:O	2.08	0.53
1:A:416:PRO:O	1:A:420:ILE:HG22	2.08	0.53
1:A:470:LEU:HD23	1:A:497:GLN:HB2	1.91	0.53
1:A:374:LYS:NZ	1:A:374:LYS:HB2	2.23	0.53
1:A:377:TYR:HE1	1:A:411:ALA:HA	1.74	0.53
1:A:430:HIS:O	1:A:433:LEU:CD2	2.57	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:504:VAL:O	1:A:505:PHE:C	2.46	0.53
1:A:463:ILE:HD11	1:A:469:VAL:HG13	1.91	0.53
1:A:515:ILE:C	1:A:517:GLU:H	2.12	0.52
1:A:514:LYS:NZ	1:A:520:GLN:HG3	2.23	0.52
1:A:343:THR:CG2	1:A:367:VAL:HB	2.39	0.52
1:A:433:LEU:CD1	1:A:527:LYS:HE3	2.40	0.52
1:A:465:ASN:HD22	1:A:465:ASN:N	2.06	0.52
1:A:264:TYR:CE1	1:A:269:TRP:NE1	2.77	0.52
1:A:365:LEU:HD13	1:A:365:LEU:N	2.24	0.52
1:A:406:SER:C	1:A:408:ALA:H	2.13	0.52
1:A:250:ILE:N	1:A:273:SER:HA	2.25	0.52
1:A:430:HIS:HA	1:A:433:LEU:HD22	1.90	0.52
1:A:470:LEU:N	1:A:470:LEU:HD22	2.25	0.52
1:A:267:ASP:HA	1:A:270:LEU:CD2	2.36	0.52
1:A:523:ASN:C	1:A:525:ARG:N	2.62	0.52
1:A:459:TYR:HD1	1:A:460:ARG:O	1.92	0.51
1:A:243:VAL:HG13	1:A:253:TYR:H	1.74	0.51
1:A:422:ARG:HG3	1:A:423:SER:H	1.73	0.51
1:A:395:HIS:CE1	1:A:489:VAL:HG21	2.45	0.51
1:A:430:HIS:ND1	1:A:431:GLU:HG3	2.25	0.51
1:A:442:THR:O	1:A:446:LYS:HB2	2.11	0.51
1:A:449:THR:HG22	1:A:450:ALA:N	2.26	0.51
1:A:527:LYS:C	1:A:529:ASP:N	2.63	0.51
1:A:269:TRP:C	1:A:271:GLY:H	2.15	0.51
1:A:378:LYS:HE2	1:A:379:VAL:HG13	1.93	0.51
1:A:391:PHE:HB2	1:A:408:ALA:HB2	1.93	0.51
1:A:254:THR:CG2	1:A:258:ILE:HD13	2.41	0.50
1:A:314:MET:HE3	1:A:415:LEU:HD21	1.93	0.50
1:A:309:SER:O	1:A:310:THR:CG2	2.52	0.50
1:A:428:TYR:CD2	1:A:436:MET:SD	3.05	0.50
1:A:394:ARG:HG3	1:A:489:VAL:O	2.11	0.50
1:A:456:SER:HB3	1:A:473:THR:O	2.11	0.50
1:A:279:HIS:ND1	1:A:280:LEU:N	2.59	0.50
1:A:384:LEU:HD22	1:A:384:LEU:H	1.76	0.50
1:A:419:LEU:O	1:A:420:ILE:C	2.49	0.50
1:A:433:LEU:H	1:A:433:LEU:CD2	2.24	0.50
1:A:465:ASN:HB3	1:A:504:VAL:O	2.12	0.50
1:A:269:TRP:C	1:A:271:GLY:N	2.63	0.50
1:A:255:THR:C	1:A:257:SER:H	2.15	0.49
1:A:397:ALA:HB1	1:A:534:LEU:HB2	1.94	0.49
1:A:449:THR:HG22	1:A:450:ALA:H	1.76	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:342:LEU:N	1:A:342:LEU:CD1	2.75	0.49
1:A:379:VAL:CB	1:A:380:PRO:CD	2.84	0.49
1:A:401:ILE:HG12	1:A:401:ILE:O	2.13	0.49
1:A:409:VAL:HG21	1:A:416:PRO:HD3	1.94	0.49
1:A:429:HIS:HD2	1:A:460:ARG:CZ	2.24	0.49
1:A:464:GLN:OE1	1:A:464:GLN:O	2.29	0.49
1:A:258:ILE:C	1:A:260:ASP:H	2.15	0.49
1:A:457:LYS:HA	1:A:457:LYS:HE2	1.94	0.49
1:A:316:ARG:HH12	1:A:318:TYR:HE1	1.60	0.49
1:A:317:ARG:HB3	1:A:319:ARG:HH12	1.77	0.49
1:A:463:ILE:HG21	1:A:504:VAL:HG13	1.95	0.49
1:A:503:ASN:OD1	1:A:506:GLU:HB2	2.13	0.49
1:A:250:ILE:HD12	1:A:272:ARG:HA	1.95	0.49
1:A:420:ILE:HG12	1:A:420:ILE:O	2.12	0.49
1:A:244:ILE:HG21	1:A:249:GLY:O	2.12	0.49
1:A:390:LYS:O	1:A:390:LYS:HG3	2.13	0.48
1:A:422:ARG:HG2	1:A:423:SER:H	1.77	0.48
1:A:519:ALA:C	1:A:521:SER:H	2.15	0.48
1:A:430:HIS:HE1	1:A:431:GLU:HG3	1.73	0.48
1:A:344:PHE:HA	1:A:366:LEU:HD13	1.96	0.48
1:A:409:VAL:HG22	1:A:416:PRO:HD3	1.94	0.48
1:A:482:TRP:C	1:A:484:ARG:N	2.66	0.48
1:A:484:ARG:HD3	1:A:538:VAL:OXT	2.13	0.48
1:A:524:THR:O	1:A:528:GLU:HG2	2.13	0.48
1:A:459:TYR:CD1	1:A:459:TYR:O	2.67	0.48
1:A:503:ASN:O	1:A:504:VAL:C	2.52	0.48
1:A:441:GLU:O	1:A:443:VAL:N	2.47	0.48
1:A:412:LEU:HD23	1:A:463:ILE:HD12	1.96	0.47
1:A:465:ASN:HD22	1:A:466:GLY:N	2.13	0.47
1:A:488:PHE:C	1:A:488:PHE:CD2	2.88	0.47
1:A:262:LEU:CD1	1:A:264:TYR:HE2	2.28	0.47
1:A:266:ARG:HG3	1:A:267:ASP:N	2.25	0.47
1:A:286:PHE:C	1:A:288:SER:H	2.17	0.47
1:A:456:SER:HB2	1:A:475:TRP:HE1	1.80	0.47
1:A:471:LEU:N	1:A:471:LEU:HD22	2.29	0.47
1:A:274:PHE:O	1:A:277:PHE:N	2.48	0.47
1:A:345:ARG:HH21	1:A:367:VAL:HG21	1.80	0.47
1:A:316:ARG:HA	1:A:335:TYR:HA	1.96	0.47
1:A:448:GLN:HG3	1:A:538:VAL:O	2.15	0.47
1:A:433:LEU:CD2	1:A:433:LEU:N	2.78	0.47
1:A:241:CYS:H	1:A:255:THR:CB	2.28	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:244:ILE:O	1:A:366:LEU:N	2.46	0.46
1:A:456:SER:OG	1:A:457:LYS:N	2.49	0.46
1:A:286:PHE:O	1:A:288:SER:N	2.48	0.46
1:A:430:HIS:O	1:A:433:LEU:HD22	2.16	0.46
1:A:391:PHE:O	1:A:493:HIS:HB2	2.15	0.46
1:A:274:PHE:O	1:A:276:ASP:N	2.48	0.46
1:A:243:VAL:HG23	1:A:366:LEU:O	2.16	0.46
1:A:380:PRO:HG2	1:A:381:ASP:N	2.31	0.46
1:A:536:GLU:OE1	1:A:538:VAL:HG12	2.16	0.46
1:A:406:SER:C	1:A:408:ALA:N	2.69	0.45
1:A:465:ASN:N	1:A:465:ASN:ND2	2.64	0.45
1:A:455:CYS:HA	1:A:474:GLU:HA	1.96	0.45
1:A:456:SER:CB	1:A:475:TRP:HE1	2.28	0.45
1:A:250:ILE:CG2	1:A:251:VAL:H	2.07	0.45
1:A:527:LYS:O	1:A:530:ILE:N	2.47	0.45
1:A:258:ILE:C	1:A:260:ASP:N	2.70	0.45
1:A:244:ILE:CG2	1:A:249:GLY:O	2.64	0.45
1:A:275:ILE:O	1:A:278:VAL:HG12	2.17	0.45
1:A:265:PRO:C	1:A:268:MET:HB3	2.37	0.45
1:A:406:SER:O	1:A:408:ALA:N	2.50	0.45
1:A:422:ARG:HG2	1:A:423:SER:N	2.30	0.44
1:A:384:LEU:H	1:A:384:LEU:CD2	2.30	0.44
1:A:484:ARG:HG2	1:A:538:VAL:CG1	2.39	0.44
1:A:484:ARG:HH21	1:A:484:ARG:HB2	1.82	0.44
1:A:512:LYS:N	1:A:512:LYS:HD2	2.32	0.44
1:A:412:LEU:CD2	1:A:463:ILE:HD12	2.47	0.44
1:A:531:VAL:O	1:A:534:LEU:HD23	2.17	0.44
1:A:419:LEU:CD1	1:A:419:LEU:N	2.80	0.44
1:A:519:ALA:C	1:A:521:SER:N	2.70	0.44
1:A:267:ASP:O	1:A:270:LEU:HB2	2.17	0.44
1:A:462:LEU:HD12	1:A:463:ILE:O	2.17	0.44
1:A:513:LEU:HB2	1:A:514:LYS:H	1.40	0.44
1:A:510:THR:HG23	1:A:510:THR:O	2.18	0.44
1:A:377:TYR:O	1:A:502:CYS:SG	2.76	0.43
1:A:378:LYS:HZ3	1:A:379:VAL:CG2	2.31	0.43
1:A:413:GLY:HA2	1:A:505:PHE:CE1	2.52	0.43
1:A:414:TYR:HB3	1:A:419:LEU:HD13	2.00	0.43
1:A:523:ASN:C	1:A:525:ARG:H	2.20	0.43
1:A:413:GLY:CA	1:A:504:VAL:CG1	2.92	0.43
1:A:465:ASN:ND2	1:A:465:ASN:H	2.15	0.43
1:A:316:ARG:O	1:A:317:ARG:O	2.37	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:446:LYS:HD2	1:A:452:ALA:HB1	2.01	0.43
1:A:463:ILE:HB	1:A:465:ASN:ND2	2.33	0.43
1:A:254:THR:CG2	1:A:258:ILE:HG21	2.37	0.43
1:A:444:MET:HE2	1:A:533:ARG:HD2	2.01	0.43
1:A:494:ARG:HD3	1:A:495:VAL:O	2.18	0.43
1:A:429:HIS:CB	1:A:460:ARG:HG3	2.45	0.43
1:A:461:PHE:C	1:A:461:PHE:CD2	2.91	0.43
1:A:512:LYS:N	1:A:512:LYS:CD	2.82	0.42
1:A:430:HIS:HA	1:A:433:LEU:CD2	2.50	0.42
1:A:504:VAL:HG12	1:A:505:PHE:N	2.33	0.42
1:A:314:MET:HE3	1:A:415:LEU:CD2	2.49	0.42
1:A:344:PHE:CG	1:A:366:LEU:HD11	2.54	0.42
1:A:427:PHE:O	1:A:428:TYR:C	2.57	0.42
1:A:476:THR:CG2	1:A:490:VAL:HB	2.47	0.42
1:A:513:LEU:O	1:A:514:LYS:HD3	2.19	0.42
1:A:414:TYR:HB3	1:A:419:LEU:CD1	2.49	0.42
1:A:428:TYR:HD2	1:A:436:MET:SD	2.42	0.42
1:A:250:ILE:HA	1:A:273:SER:HA	2.00	0.42
1:A:428:TYR:CD2	1:A:436:MET:CG	3.02	0.42
1:A:395:HIS:O	1:A:489:VAL:HG23	2.20	0.42
1:A:457:LYS:O	1:A:458:PRO:C	2.57	0.42
1:A:264:TYR:CE2	1:A:318:TYR:HB2	2.54	0.42
1:A:488:PHE:HD2	1:A:488:PHE:C	2.22	0.42
1:A:463:ILE:CG2	1:A:504:VAL:HG13	2.49	0.42
1:A:529:ASP:HA	1:A:532:LYS:HG2	2.01	0.42
1:A:240:PHE:CE1	1:A:258:ILE:HD12	2.55	0.42
1:A:319:ARG:HA	1:A:319:ARG:HD3	1.91	0.42
1:A:378:LYS:HZ3	1:A:379:VAL:HG21	1.85	0.42
1:A:428:TYR:CE2	1:A:436:MET:SD	3.13	0.42
1:A:523:ASN:O	1:A:525:ARG:N	2.52	0.42
1:A:272:ARG:HD3	1:A:272:ARG:N	2.27	0.41
1:A:273:SER:O	1:A:274:PHE:C	2.58	0.41
1:A:420:ILE:O	1:A:420:ILE:HG23	2.20	0.41
1:A:384:LEU:HB3	1:A:388:SER:HB3	2.01	0.41
1:A:279:HIS:HB2	1:A:335:TYR:HD2	1.86	0.41
1:A:344:PHE:CE1	1:A:366:LEU:HD21	2.56	0.41
1:A:514:LYS:HZ2	1:A:520:GLN:HG3	1.85	0.41
1:A:382:GLU:O	1:A:497:GLN:HG2	2.21	0.41
1:A:254:THR:O	1:A:255:THR:O	2.39	0.41
1:A:485:LYS:HG2	1:A:486:LEU:N	2.35	0.41
1:A:536:GLU:CD	1:A:537:THR:N	2.74	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:315:LEU:HD12	1:A:338:PHE:CD2	2.56	0.41
1:A:387:LYS:O	1:A:387:LYS:HD3	2.21	0.41
1:A:241:CYS:H	1:A:255:THR:HB	1.86	0.41
1:A:255:THR:O	1:A:257:SER:N	2.54	0.41
1:A:378:LYS:CD	1:A:379:VAL:HG22	2.46	0.41
1:A:446:LYS:C	1:A:448:GLN:H	2.24	0.41
1:A:238:ASP:HB3	1:A:261:VAL:CG2	2.51	0.41
1:A:404:VAL:HG11	1:A:416:PRO:HA	2.02	0.41
1:A:239:SER:O	1:A:240:PHE:HB3	2.21	0.40
1:A:244:ILE:HG23	1:A:251:VAL:HA	2.04	0.40
1:A:401:ILE:HG12	1:A:420:ILE:HA	2.04	0.40
1:A:318:TYR:C	1:A:319:ARG:HH11	2.24	0.40
1:A:427:PHE:O	1:A:428:TYR:O	2.40	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:450:ALA:O	1:A:450:ALA:O[6_765]	1.63	0.57

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	249/312 (80%)	152 (61%)	55 (22%)	42 (17%)	0 3

All (42) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	244	ILE
1	A	245	SER
1	A	250	ILE

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Mol	Chain	Res	Type
1	A	275	ILE
1	A	317	ARG
1	A	379	VAL
1	A	380	PRO
1	A	385	SER
1	A	420	ILE
1	A	428	TYR
1	A	431	GLU
1	A	504	VAL
1	A	505	PHE
1	A	514	LYS
1	A	515	ILE
1	A	536	GLU
1	A	537	THR
1	A	248	ASP
1	A	255	THR
1	A	270	LEU
1	A	271	GLY
1	A	287	ALA
1	A	293	GLY
1	A	335	TYR
1	A	376	SER
1	A	405	ASP
1	A	442	THR
1	A	450	ALA
1	A	503	ASN
1	A	506	GLU
1	A	516	SER
1	A	528	GLU
1	A	266	ARG
1	A	281	LYS
1	A	384	LEU
1	A	407	ALA
1	A	416	PRO
1	A	381	ASP
1	A	388	SER
1	A	283	ARG
1	A	476	THR
1	A	463	ILE

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	229/271 (84%)	186 (81%)	43 (19%)	1 10

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	238	ASP
1	A	242	CYS
1	A	244	ILE
1	A	262	LEU
1	A	264	TYR
1	A	272	ARG
1	A	281	LYS
1	A	282	ASP
1	A	290	ILE
1	A	314	MET
1	A	335	TYR
1	A	342	LEU
1	A	365	LEU
1	A	374	LYS
1	A	378	LYS
1	A	400	ILE
1	A	401	ILE
1	A	417	GLN
1	A	418	ASP
1	A	419	LEU
1	A	426	ASP
1	A	432	ASP
1	A	433	LEU
1	A	440	TYR
1	A	454	PHE
1	A	456	SER
1	A	457	LYS
1	A	460	ARG
1	A	461	PHE
1	A	464	GLN

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Mol	Chain	Res	Type
1	A	465	ASN
1	A	471	LEU
1	A	484	ARG
1	A	488	PHE
1	A	489	VAL
1	A	494	ARG
1	A	497	GLN
1	A	512	LYS
1	A	513	LEU
1	A	514	LYS
1	A	520	GLN
1	A	533	ARG
1	A	536	GLU

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	386	GLN
1	A	395	HIS
1	A	403	HIS
1	A	417	GLN
1	A	429	HIS
1	A	465	ASN
1	A	493	HIS
1	A	497	GLN

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 [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	257/312 (82%)	-0.68	0 100 100	79, 145, 210, 210	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 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.