



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 7, 2021 – 01:35 PM EDT

PDB ID : 6VIN
Title : Crystallographic structure of the circularly permuted human Taspase1 protein
Authors : Martin-Garcia, J.M.; Fromme, P.
Deposited on : 2020-01-13
Resolution : 3.04 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 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 : 2.23.1
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) : 2.23.1

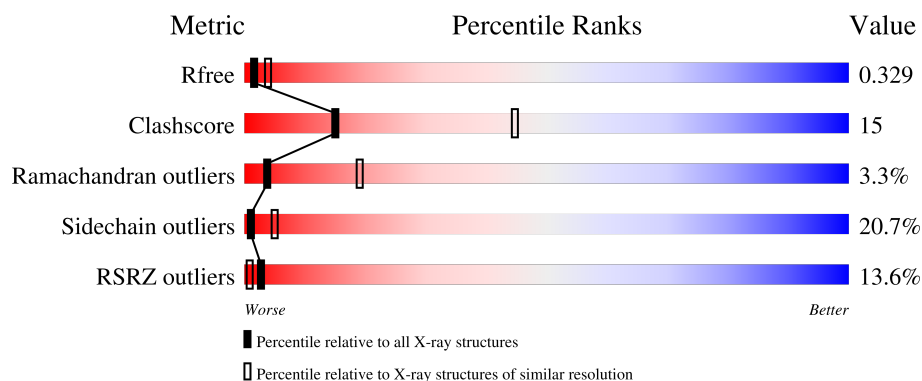
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.04 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	2752 (3.08-3.00)
Clashscore	141614	3096 (3.08-3.00)
Ramachandran outliers	138981	2986 (3.08-3.00)
Sidechain outliers	138945	2988 (3.08-3.00)
RSRZ outliers	127900	2636 (3.08-3.00)

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	389	<div> <div>14%</div> <div>65%</div> <div>22%</div> <div>7%</div> <div>6%</div> </div>
1	B	389	<div> <div>12%</div> <div>60%</div> <div>29%</div> <div>5%</div> <div>6%</div> </div>

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 10500 atoms, of which 5151 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 Threonine aspartase 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	364	Total	C	H	N	O	S	129	0	0
			5257	1659	2582	484	511	21			
1	B	364	Total	C	H	N	O	S	129	0	0
			5243	1659	2569	483	511	21			

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP Q9H6P5
A	185	GLY	-	linker	UNP Q9H6P5
A	186	SER	-	linker	UNP Q9H6P5
A	187	GLY	-	linker	UNP Q9H6P5
A	188	SER	-	linker	UNP Q9H6P5
A	382	LEU	-	expression tag	UNP Q9H6P5
A	383	GLU	-	expression tag	UNP Q9H6P5
A	384	HIS	-	expression tag	UNP Q9H6P5
A	385	HIS	-	expression tag	UNP Q9H6P5
A	386	HIS	-	expression tag	UNP Q9H6P5
A	387	HIS	-	expression tag	UNP Q9H6P5
A	388	HIS	-	expression tag	UNP Q9H6P5
A	389	HIS	-	expression tag	UNP Q9H6P5
B	1	MET	-	initiating methionine	UNP Q9H6P5
B	185	GLY	-	linker	UNP Q9H6P5
B	186	SER	-	linker	UNP Q9H6P5
B	187	GLY	-	linker	UNP Q9H6P5
B	188	SER	-	linker	UNP Q9H6P5
B	382	LEU	-	expression tag	UNP Q9H6P5
B	383	GLU	-	expression tag	UNP Q9H6P5
B	384	HIS	-	expression tag	UNP Q9H6P5
B	385	HIS	-	expression tag	UNP Q9H6P5
B	386	HIS	-	expression tag	UNP Q9H6P5
B	387	HIS	-	expression tag	UNP Q9H6P5
B	388	HIS	-	expression tag	UNP Q9H6P5

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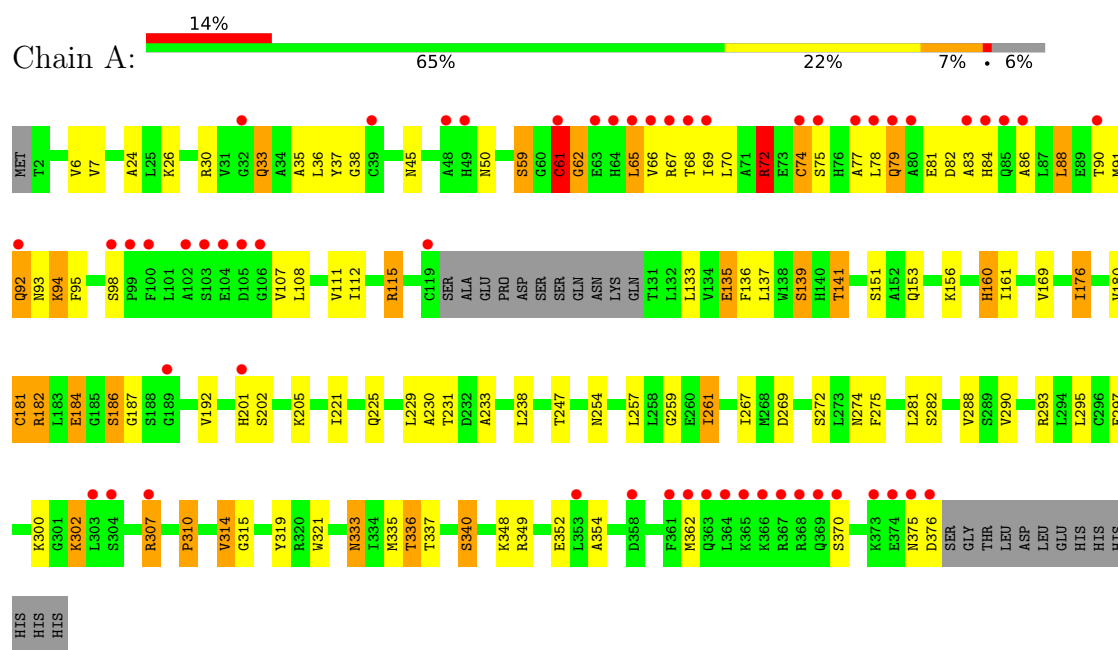
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Chain	Residue	Modelled	Actual	Comment	Reference
B	389	HIS	-	expression tag	UNP Q9H6P5

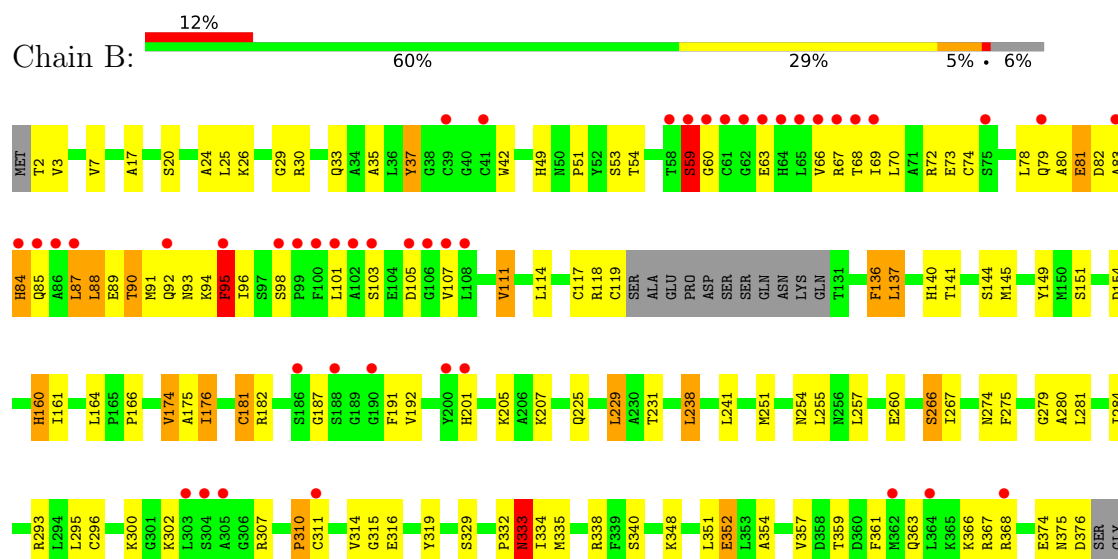
3 Residue-property plots [i](#)

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: Threonine aspartase 1



• Molecule 1: Threonine aspartase 1



THR
LEU
ASP
LEU
GLU
HIS
HIS
HIS
HIS
HIS

4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	196.00Å 196.00Å 196.91Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.05 – 3.04 49.00 – 3.04	Depositor EDS
% Data completeness (in resolution range)	54.4 (49.05-3.04) 54.4 (49.00-3.04)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.37 (at 3.07Å)	Xtriage
Refinement program	REFMAC 5.8.0253	Depositor
R, R_{free}	0.237 , 0.307 0.266 , 0.329	Depositor DCC
R_{free} test set	1568 reflections (10.24%)	wwPDB-VP
Wilson B-factor (Å ²)	140.2	Xtriage
Anisotropy	0.261	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 133.6	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	0.91	EDS
Total number of atoms	10500	wwPDB-VP
Average B, all atoms (Å ²)	165.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.48% 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	0.75	0/2718	1.03	4/3667 (0.1%)
1	B	0.79	1/2718 (0.0%)	1.03	2/3668 (0.1%)
All	All	0.77	1/5436 (0.0%)	1.03	6/7335 (0.1%)

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

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	59	SER	C-N	13.28	1.56	1.33

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	59	SER	C-N-CA	-10.32	100.63	122.30
1	B	49	HIS	C-N-CA	8.17	142.13	121.70
1	A	335	MET	O-C-N	-6.38	112.49	122.70
1	A	68	THR	CB-CA-C	-6.33	94.51	111.60
1	A	74	CYS	N-CA-CB	5.54	120.58	110.60
1	A	335	MET	CA-C-N	5.15	128.54	117.20

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	72	ARG	Sidechain
1	B	375	ASN	Peptide

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	2675	2582	2662	75	15
1	B	2674	2569	2656	99	10
All	All	5349	5151	5318	161	16

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 15.

All (161) 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:B:88:LEU:HG	1:B:91:MET:HB2	1.27	1.14
1:B:7:VAL:HG23	1:B:192:VAL:HG23	1.29	1.13
1:B:88:LEU:HG	1:B:91:MET:CB	1.91	1.00
1:A:70:LEU:HD11	1:A:91:MET:HG3	1.45	0.98
1:B:51:PRO:HD2	1:B:79:GLN:OE1	1.68	0.93
1:B:7:VAL:CG2	1:B:192:VAL:HG23	2.00	0.92
1:A:95:PHE:CE2	1:A:107:VAL:HG12	2.06	0.91
1:A:50:ASN:OD1	1:A:79:GLN:NE2	2.05	0.89
1:B:88:LEU:CG	1:B:91:MET:HB2	2.02	0.88
1:B:84:HIS:CD2	1:B:137:LEU:HD21	2.10	0.85
1:A:72:ARG:NH2	1:B:68:THR:O	2.12	0.83
1:A:78:LEU:HD23	1:A:86:ALA:O	1.80	0.81
1:B:137:LEU:HD12	1:B:137:LEU:N	1.96	0.80
1:A:59:SER:C	1:A:65:LEU:HD21	2.02	0.79
1:A:302:LYS:HA	1:A:302:LYS:HE2	1.66	0.76
1:A:95:PHE:CD2	1:A:107:VAL:CG1	2.68	0.76
1:A:72:ARG:HD2	1:B:69:ILE:HD12	1.69	0.75
1:B:88:LEU:HD21	1:B:92:GLN:HG3	1.67	0.75
1:B:83:ALA:HB3	1:B:84:HIS:HD2	1.53	0.73
1:B:88:LEU:HA	1:B:91:MET:HB2	1.69	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:33:GLN:NE2	1:A:65:LEU:HD13	2.03	0.72
1:B:83:ALA:HB3	1:B:84:HIS:CD2	2.24	0.71
1:A:74:CYS:SG	1:A:90:THR:OG1	2.38	0.71
1:A:33:GLN:CD	1:A:65:LEU:HD13	2.11	0.71
1:A:69:ILE:CD1	1:B:72:ARG:CD	2.69	0.70
1:B:89:GLU:HG2	1:B:93:ASN:HD22	1.56	0.69
1:A:302:LYS:O	1:A:302:LYS:HG3	1.93	0.69
1:A:95:PHE:CD2	1:A:107:VAL:HG12	2.28	0.69
1:B:63:GLU:OE1	1:B:63:GLU:HA	1.92	0.67
1:B:88:LEU:HD11	1:B:174:VAL:HG21	1.78	0.66
1:A:95:PHE:CE2	1:A:107:VAL:CG1	2.79	0.66
1:B:74:CYS:O	1:B:78:LEU:HD12	1.97	0.65
1:A:82:ASP:OD1	1:A:83:ALA:N	2.28	0.64
1:B:88:LEU:HD11	1:B:174:VAL:CG2	2.27	0.64
1:B:79:GLN:OE1	1:B:79:GLN:HA	1.97	0.64
1:B:84:HIS:CD2	1:B:84:HIS:N	2.65	0.64
1:B:284:ILE:O	1:B:335:MET:HE1	1.98	0.64
1:B:164:LEU:HD23	1:B:175:ALA:CB	2.29	0.63
1:A:69:ILE:HD12	1:B:72:ARG:NE	2.14	0.63
1:B:68:THR:HG21	1:B:95:PHE:CD1	2.35	0.62
1:B:144:SER:OG	1:B:161:ILE:HD11	1.99	0.61
1:A:337:THR:HG1	1:A:340:SER:HG	1.44	0.61
1:A:75:SER:HA	1:A:78:LEU:HD12	1.81	0.61
1:A:107:VAL:HG12	1:A:107:VAL:O	2.03	0.58
1:A:314:VAL:HG11	1:B:314:VAL:HG21	1.83	0.58
1:B:136:PHE:C	1:B:137:LEU:HD12	2.23	0.58
1:B:284:ILE:O	1:B:335:MET:CE	2.52	0.58
1:A:38:GLY:C	1:A:69:ILE:HG23	2.24	0.57
1:B:88:LEU:CA	1:B:91:MET:HB2	2.35	0.57
1:B:107:VAL:HG12	1:B:107:VAL:O	2.05	0.56
1:B:84:HIS:HA	1:B:137:LEU:CD2	2.35	0.56
1:B:88:LEU:CD2	1:B:92:GLN:HG3	2.33	0.56
1:B:266:SER:HG	1:B:275:PHE:HE1	1.53	0.56
1:B:92:GLN:HA	1:B:96:ILE:HB	1.87	0.55
1:A:95:PHE:CD2	1:A:107:VAL:HG11	2.42	0.55
1:A:352:GLU:HA	1:A:352:GLU:OE1	2.07	0.55
1:B:88:LEU:HA	1:B:91:MET:CB	2.37	0.55
1:B:164:LEU:HD23	1:B:175:ALA:HB2	1.88	0.55
1:B:88:LEU:CD2	1:B:92:GLN:CG	2.85	0.54
1:B:181:CYS:SG	1:B:182:ARG:N	2.81	0.54
1:A:281:LEU:HD21	1:A:290:VAL:HG21	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:70:LEU:C	1:B:70:LEU:HD23	2.29	0.53
1:B:83:ALA:CB	1:B:84:HIS:HD2	2.21	0.53
1:B:333:ASN:C	1:B:333:ASN:ND2	2.62	0.53
1:A:69:ILE:HD12	1:B:72:ARG:CD	2.38	0.53
1:A:267:ILE:HG23	1:A:295:LEU:HD12	1.91	0.53
1:B:174:VAL:HG22	1:B:174:VAL:O	2.08	0.53
1:A:141:THR:O	1:A:141:THR:OG1	2.23	0.52
1:A:69:ILE:CD1	1:B:72:ARG:HD3	2.39	0.52
1:A:83:ALA:HB2	1:A:135:GLU:OE1	2.09	0.52
1:B:2:THR:HG23	1:B:20:SER:HB2	1.91	0.52
1:B:89:GLU:O	1:B:93:ASN:HB2	2.09	0.52
1:A:83:ALA:HA	1:A:115:ARG:NH2	2.26	0.51
1:A:139:SER:HB2	1:A:176:ILE:HG23	1.92	0.51
1:B:352:GLU:HA	1:B:352:GLU:OE2	2.11	0.51
1:A:297:GLU:OE1	1:A:297:GLU:HA	2.10	0.51
1:A:37:TYR:HA	1:A:275:PHE:CZ	2.46	0.51
1:B:88:LEU:O	1:B:88:LEU:HD23	2.11	0.51
1:A:307:ARG:NH2	1:B:351:LEU:HD11	2.25	0.51
1:B:136:PHE:CD1	1:B:191:PHE:CZ	2.99	0.50
1:A:267:ILE:HG23	1:A:295:LEU:CD1	2.42	0.50
1:B:357:VAL:O	1:B:357:VAL:HG12	2.11	0.50
1:A:261:ILE:O	1:A:282:SER:OG	2.28	0.50
1:B:7:VAL:CG2	1:B:192:VAL:CG2	2.84	0.50
1:A:77:ALA:O	1:A:86:ALA:HB1	2.12	0.49
1:B:84:HIS:CD2	1:B:137:LEU:CD2	2.90	0.49
1:B:164:LEU:HD23	1:B:175:ALA:HB3	1.94	0.49
1:A:36:LEU:HD12	1:A:66:VAL:HG22	1.95	0.49
1:A:61:CYS:SG	1:A:62:GLY:N	2.84	0.49
1:B:95:PHE:N	1:B:98:SER:OG	2.46	0.49
1:A:90:THR:HA	1:A:94:LYS:HB2	1.95	0.49
1:A:247:THR:O	1:A:247:THR:OG1	2.28	0.48
1:A:261:ILE:O	1:A:261:ILE:HG22	2.13	0.48
1:B:17:ALA:HB3	1:B:238:LEU:CD2	2.42	0.48
1:B:136:PHE:CD1	1:B:136:PHE:N	2.81	0.48
1:B:17:ALA:HB3	1:B:238:LEU:HD23	1.94	0.48
1:B:51:PRO:CD	1:B:79:GLN:OE1	2.52	0.48
1:B:82:ASP:OD1	1:B:83:ALA:N	2.42	0.48
1:A:93:ASN:O	1:A:98:SER:CB	2.62	0.47
1:B:137:LEU:N	1:B:137:LEU:CD1	2.70	0.47
1:A:83:ALA:HA	1:A:115:ARG:CZ	2.44	0.47
1:B:88:LEU:C	1:B:91:MET:H	2.18	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6:VAL:HG22	1:A:112:ILE:CD1	2.45	0.47
1:A:69:ILE:CD1	1:B:72:ARG:HD2	2.43	0.47
1:A:69:ILE:HD13	1:B:72:ARG:HD3	1.96	0.47
1:A:307:ARG:HH22	1:B:351:LEU:HD11	1.78	0.47
1:A:333:ASN:O	1:A:336:THR:N	2.48	0.47
1:B:70:LEU:HD23	1:B:70:LEU:O	2.14	0.47
1:B:251:MET:HE2	1:B:334:ILE:HG22	1.98	0.46
1:A:181:CYS:SG	1:A:182:ARG:N	2.89	0.46
1:A:192:VAL:HG13	1:A:192:VAL:O	2.15	0.45
1:A:307:ARG:HD2	1:A:307:ARG:N	2.31	0.45
1:B:95:PHE:HD2	1:B:107:VAL:CG1	2.29	0.45
1:B:25:LEU:HD23	1:B:351:LEU:HD12	1.98	0.45
1:B:238:LEU:HD13	1:B:238:LEU:HA	1.84	0.45
1:B:267:ILE:HG23	1:B:295:LEU:HD12	1.99	0.45
1:A:184:GLU:OE2	1:A:186:SER:OG	2.33	0.45
1:B:68:THR:HG1	1:B:95:PHE:HE1	1.64	0.45
1:A:202:SER:O	1:A:202:SER:OG	2.29	0.45
1:A:321:TRP:O	1:A:321:TRP:CE3	2.70	0.45
1:B:67:ARG:HB2	1:B:101:LEU:HD21	1.97	0.45
1:B:84:HIS:HD2	1:B:84:HIS:N	2.12	0.45
1:B:293:ARG:HD2	1:B:293:ARG:HA	1.84	0.44
1:A:67:ARG:O	1:B:37:TYR:OH	2.24	0.44
1:A:93:ASN:O	1:A:98:SER:HB3	2.18	0.44
1:A:88:LEU:HD22	1:A:88:LEU:HA	1.78	0.44
1:B:302:LYS:HD2	1:B:302:LYS:HA	1.75	0.44
1:A:33:GLN:HB3	1:A:66:VAL:HG23	1.99	0.44
1:A:201:HIS:CG	1:A:201:HIS:O	2.70	0.43
1:A:38:GLY:O	1:A:69:ILE:CG2	2.66	0.43
1:B:88:LEU:HD23	1:B:92:GLN:HB2	1.98	0.43
1:A:69:ILE:HG22	1:A:69:ILE:O	2.19	0.43
1:B:279:GLY:O	1:B:281:LEU:N	2.50	0.43
1:B:201:HIS:CG	1:B:201:HIS:O	2.71	0.43
1:B:316:GLU:OE1	1:B:316:GLU:N	2.52	0.43
1:B:363:GLN:O	1:B:367:ARG:HD2	2.19	0.43
1:B:84:HIS:HD2	1:B:137:LEU:HD21	1.76	0.43
1:A:238:LEU:HB3	1:A:288:VAL:HG11	2.01	0.42
1:A:269:ASP:OD2	1:A:272:SER:OG	2.37	0.42
1:A:59:SER:C	1:A:65:LEU:CD2	2.82	0.42
1:B:59:SER:OG	1:B:60:GLY:N	2.53	0.42
1:B:88:LEU:HD11	1:B:174:VAL:HG22	2.00	0.42
1:B:73:GLU:OE2	1:B:73:GLU:HA	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:229:LEU:N	1:B:229:LEU:CD1	2.83	0.42
1:A:37:TYR:CD1	1:B:66:VAL:HG12	2.55	0.41
1:B:95:PHE:CD2	1:B:107:VAL:HG13	2.55	0.41
1:A:93:ASN:O	1:A:98:SER:HB2	2.21	0.41
1:B:90:THR:HA	1:B:94:LYS:HB2	2.01	0.41
1:B:3:VAL:HG21	1:B:241:LEU:HB3	2.03	0.41
1:A:281:LEU:CD2	1:A:290:VAL:HG21	2.50	0.41
1:A:38:GLY:O	1:A:69:ILE:HG23	2.19	0.41
1:B:140:HIS:CD2	1:B:140:HIS:N	2.85	0.41
1:B:87:LEU:HD11	1:B:111:VAL:HG21	2.02	0.41
1:B:94:LYS:HA	1:B:98:SER:CB	2.51	0.41
1:B:24:ALA:O	1:B:255:LEU:O	2.39	0.40
1:B:80:ALA:O	1:B:82:ASP:N	2.53	0.40
1:B:284:ILE:HD13	1:B:284:ILE:HA	1.92	0.40
1:A:88:LEU:O	1:A:92:GLN:N	2.54	0.40
1:A:230:ALA:O	1:A:233:ALA:HB3	2.22	0.40
1:A:79:GLN:OE1	1:A:79:GLN:CA	2.70	0.40
1:A:91:MET:HA	1:A:95:PHE:HB3	2.03	0.40

All (16) 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:93:ASN:HD21	1:B:81:GLU:O[3_555]	0.24	1.36
1:A:93:ASN:ND2	1:B:81:GLU:O[3_555]	0.99	1.21
1:A:187:GLY:N	1:A:187:GLY:N[5_555]	1.70	0.50
1:A:187:GLY:N	1:A:187:GLY:H[5_555]	1.14	0.46
1:A:187:GLY:H	1:A:187:GLY:H[5_555]	1.15	0.45
1:A:176:ILE:O	1:B:176:ILE:O[3_555]	1.83	0.37
1:A:93:ASN:ND2	1:B:81:GLU:C[3_555]	1.88	0.32
1:B:361:PHE:HE1	1:B:361:PHE:HE1[4_555]	1.29	0.31
1:A:187:GLY:CA	1:A:187:GLY:H[5_555]	1.42	0.18
1:A:187:GLY:N	1:A:187:GLY:CA[5_555]	2.04	0.16
1:A:259:GLY:O	1:A:333:ASN:OD1[18_444]	2.05	0.15
1:A:93:ASN:HD21	1:B:81:GLU:C[3_555]	1.48	0.12
1:A:93:ASN:CG	1:B:81:GLU:O[3_555]	2.10	0.10
1:A:93:ASN:HD22	1:B:81:GLU:O[3_555]	1.53	0.07
1:A:84:HIS:NE2	1:B:176:ILE:HG22[3_555]	1.57	0.03
1:A:81:GLU:O	1:B:93:ASN:OD1[3_555]	2.18	0.02

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	360/389 (92%)	309 (86%)	42 (12%)	9 (2%)	5	25
1	B	360/389 (92%)	300 (83%)	45 (12%)	15 (4%)	3	14
All	All	720/778 (92%)	609 (85%)	87 (12%)	24 (3%)	4	19

All (24) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	24	ALA
1	B	81	GLU
1	B	87	LEU
1	B	280	ALA
1	A	35	ALA
1	A	160	HIS
1	A	169	VAL
1	A	315	GLY
1	A	354	ALA
1	B	35	ALA
1	B	315	GLY
1	B	354	ALA
1	B	29	GLY
1	B	187	GLY
1	B	333	ASN
1	A	61	CYS
1	B	37	TYR
1	B	95	PHE
1	B	160	HIS
1	B	166	PRO
1	B	332	PRO
1	B	310	PRO
1	A	310	PRO
1	A	62	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	280/304 (92%)	222 (79%)	58 (21%)	1	5
1	B	280/304 (92%)	222 (79%)	58 (21%)	1	5
All	All	560/608 (92%)	444 (79%)	116 (21%)	1	5

All (116) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	VAL
1	A	26	LYS
1	A	30	ARG
1	A	33	GLN
1	A	45	ASN
1	A	59	SER
1	A	61	CYS
1	A	65	LEU
1	A	72	ARG
1	A	79	GLN
1	A	88	LEU
1	A	92	GLN
1	A	94	LYS
1	A	108	LEU
1	A	111	VAL
1	A	115	ARG
1	A	133	LEU
1	A	135	GLU
1	A	136	PHE
1	A	137	LEU
1	A	139	SER
1	A	141	THR
1	A	151	SER
1	A	153	GLN
1	A	156	LYS
1	A	160	HIS
1	A	161	ILE

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Mol	Chain	Res	Type
1	A	176	ILE
1	A	180	VAL
1	A	181	CYS
1	A	182	ARG
1	A	184	GLU
1	A	186	SER
1	A	205	LYS
1	A	221	ILE
1	A	225	GLN
1	A	229	LEU
1	A	231	THR
1	A	254	ASN
1	A	257	LEU
1	A	261	ILE
1	A	274	ASN
1	A	293	ARG
1	A	300	LYS
1	A	302	LYS
1	A	307	ARG
1	A	310	PRO
1	A	314	VAL
1	A	319	TYR
1	A	333	ASN
1	A	336	THR
1	A	340	SER
1	A	348	LYS
1	A	349	ARG
1	A	362	MET
1	A	370	SER
1	A	375	ASN
1	A	376	ASP
1	B	26	LYS
1	B	30	ARG
1	B	33	GLN
1	B	42	TRP
1	B	53	SER
1	B	54	THR
1	B	59	SER
1	B	84	HIS
1	B	85	GLN
1	B	88	LEU
1	B	90	THR

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Mol	Chain	Res	Type
1	B	95	PHE
1	B	103	SER
1	B	105	ASP
1	B	111	VAL
1	B	114	LEU
1	B	117	CYS
1	B	118	ARG
1	B	119	CYS
1	B	136	PHE
1	B	137	LEU
1	B	141	THR
1	B	145	MET
1	B	149	TYR
1	B	151	SER
1	B	154	ASP
1	B	160	HIS
1	B	174	VAL
1	B	176	ILE
1	B	181	CYS
1	B	205	LYS
1	B	207	LYS
1	B	225	GLN
1	B	229	LEU
1	B	231	THR
1	B	238	LEU
1	B	254	ASN
1	B	257	LEU
1	B	260	GLU
1	B	266	SER
1	B	274	ASN
1	B	296	CYS
1	B	300	LYS
1	B	307	ARG
1	B	310	PRO
1	B	311	CYS
1	B	319	TYR
1	B	329	SER
1	B	333	ASN
1	B	338	ARG
1	B	340	SER
1	B	348	LYS
1	B	352	GLU

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Mol	Chain	Res	Type
1	B	359	THR
1	B	366	LYS
1	B	368	ARG
1	B	374	GLU
1	B	376	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	33	GLN
1	A	64	HIS
1	A	92	GLN
1	A	160	HIS
1	A	333	ASN
1	A	347	ASN
1	B	33	GLN
1	B	84	HIS
1	B	93	ASN
1	B	347	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 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, 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	364/389 (93%)	0.86	54 (14%) 2 1	115, 157, 235, 305	0
1	B	364/389 (93%)	0.80	45 (12%) 4 1	115, 160, 225, 378	0
All	All	728/778 (93%)	0.83	99 (13%) 3 1	115, 159, 231, 378	0

All (99) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	99	PRO	19.2
1	A	104	GLU	16.1
1	A	99	PRO	12.3
1	A	68	THR	10.7
1	B	86	ALA	10.2
1	B	102	ALA	9.3
1	A	106	GLY	8.4
1	A	100	PHE	8.2
1	B	84	HIS	8.1
1	B	85	GLN	8.1
1	B	303	LEU	7.5
1	B	100	PHE	7.5
1	A	105	ASP	7.5
1	B	98	SER	7.4
1	B	105	ASP	7.4
1	B	62	GLY	6.6
1	A	80	ALA	6.6
1	A	66	VAL	6.2
1	A	84	HIS	5.7
1	B	364	LEU	5.5
1	A	67	ARG	5.4
1	A	79	GLN	5.4
1	A	375	ASN	5.2
1	A	304	SER	4.9

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Mol	Chain	Res	Type	RSRZ
1	A	103	SER	4.9
1	B	107	VAL	4.7
1	B	66	VAL	4.6
1	A	369	GLN	4.3
1	B	63	GLU	4.3
1	A	92	GLN	4.2
1	A	83	ALA	4.2
1	A	48	ALA	4.2
1	B	362	MET	4.1
1	B	69	ILE	4.0
1	A	362	MET	3.9
1	A	98	SER	3.9
1	B	106	GLY	3.8
1	A	64	HIS	3.7
1	B	304	SER	3.6
1	B	67	ARG	3.5
1	A	201	HIS	3.5
1	B	61	CYS	3.5
1	B	201	HIS	3.5
1	B	103	SER	3.4
1	A	363	GLN	3.4
1	A	370	SER	3.3
1	B	101	LEU	3.3
1	A	61	CYS	3.2
1	A	102	ALA	3.2
1	B	95	PHE	3.2
1	B	59	SER	3.2
1	A	364	LEU	3.1
1	A	376	ASP	3.1
1	A	365	LYS	3.0
1	B	92	GLN	2.9
1	B	200	TYR	2.9
1	B	108	LEU	2.9
1	A	374	GLU	2.8
1	B	190	GLY	2.8
1	A	39	CYS	2.7
1	A	49	HIS	2.7
1	B	60	GLY	2.7
1	B	83	ALA	2.7
1	B	87	LEU	2.6
1	B	68	THR	2.6
1	A	307	ARG	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	189	GLY	2.5
1	A	32	GLY	2.5
1	A	78	LEU	2.5
1	A	75	SER	2.5
1	A	77	ALA	2.5
1	A	86	ALA	2.5
1	A	85	GLN	2.4
1	A	358	ASP	2.4
1	A	361	PHE	2.4
1	B	186	SER	2.4
1	B	39	CYS	2.4
1	B	79	GLN	2.3
1	B	368	ARG	2.3
1	A	119	CYS	2.3
1	B	305	ALA	2.3
1	B	75	SER	2.3
1	A	367	ARG	2.3
1	A	353	LEU	2.3
1	A	69	ILE	2.3
1	A	65	LEU	2.3
1	A	74	CYS	2.2
1	B	58	THR	2.2
1	A	373	LYS	2.2
1	A	63	GLU	2.2
1	B	64	HIS	2.2
1	B	188	SER	2.1
1	B	41	CYS	2.1
1	A	90	THR	2.1
1	A	368	ARG	2.0
1	A	303	LEU	2.0
1	A	366	LYS	2.0
1	B	311	CYS	2.0
1	B	65	LEU	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.