



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 09:39 PM UTC

PDB ID : 4HYD / pdb_00004hyd
Title : Structure of a presenilin family intramembrane aspartate protease in C2221 space group
Authors : Li, X.; Dang, S.; Yan, C.; Wang, J.; Shi, Y.
Deposited on : 2012-11-13
Resolution : 3.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

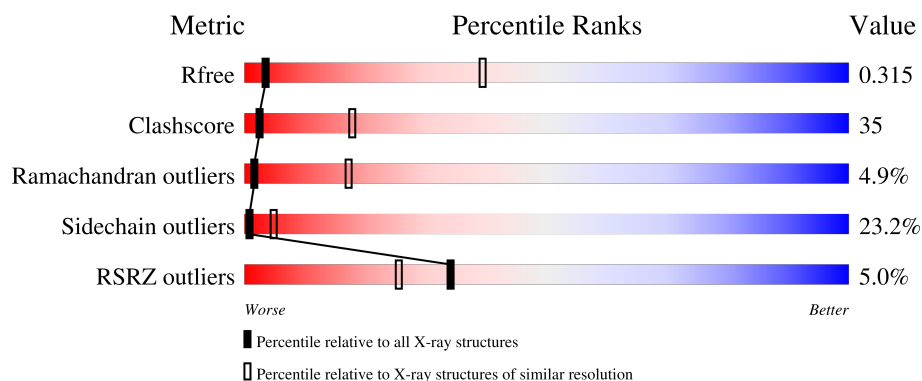
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1065 (3.96-3.64)
Clashscore	190562	1012 (3.94-3.66)
Ramachandran outliers	187476	1048 (3.96-3.64)
Sidechain outliers	187428	1043 (3.96-3.64)
RSRZ outliers	180081	1064 (3.96-3.64)

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	301	5% 30% 37% 14% • 18%
1	B	301	3% 33% 38% 10% • 18%
1	C	301	3% 30% 38% 13% • 18%
1	D	301	5% 30% 38% 12% • 18%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 7199 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 Putative uncharacterized protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	246	Total	C	N	O	S	0	0	0
			1802	1219	276	296	11			
1	B	246	Total	C	N	O	S	0	0	0
			1799	1216	276	296	11			
1	C	246	Total	C	N	O	S	0	0	0
			1799	1216	276	296	11			
1	D	246	Total	C	N	O	S	0	0	0
			1799	1216	276	296	11			

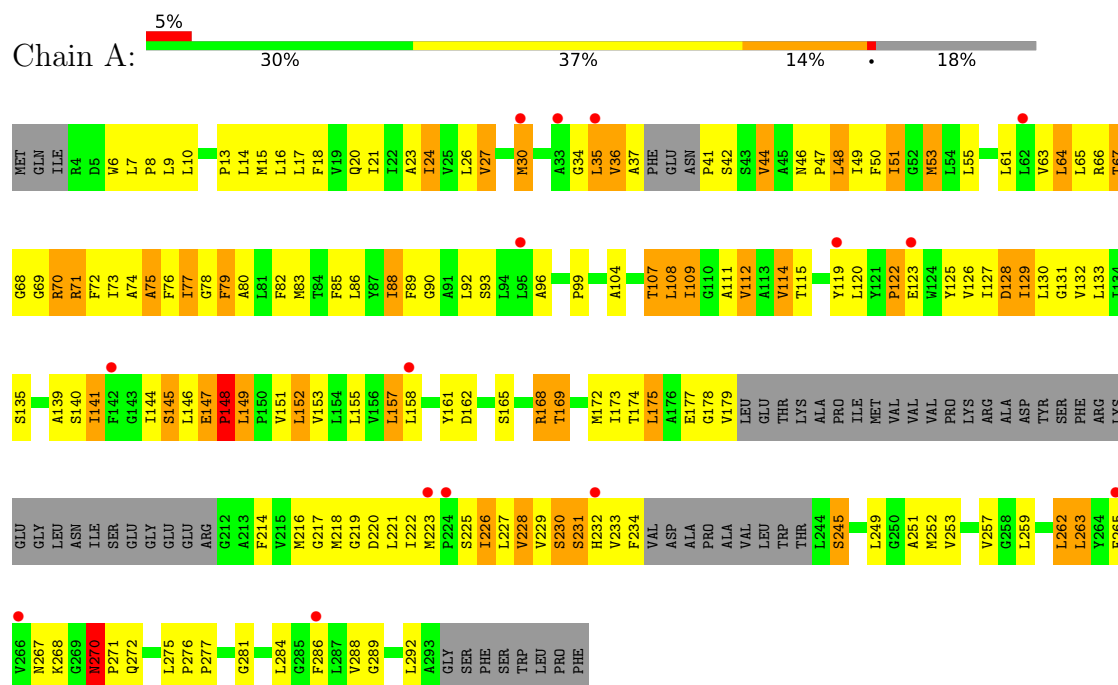
There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	40	ASN	ASP	engineered mutation	UNP A3CWV0
A	42	SER	GLU	engineered mutation	UNP A3CWV0
A	147	GLU	ALA	engineered mutation	UNP A3CWV0
A	148	PRO	VAL	engineered mutation	UNP A3CWV0
A	229	VAL	ALA	engineered mutation	UNP A3CWV0
B	40	ASN	ASP	engineered mutation	UNP A3CWV0
B	42	SER	GLU	engineered mutation	UNP A3CWV0
B	147	GLU	ALA	engineered mutation	UNP A3CWV0
B	148	PRO	VAL	engineered mutation	UNP A3CWV0
B	229	VAL	ALA	engineered mutation	UNP A3CWV0
C	40	ASN	ASP	engineered mutation	UNP A3CWV0
C	42	SER	GLU	engineered mutation	UNP A3CWV0
C	147	GLU	ALA	engineered mutation	UNP A3CWV0
C	148	PRO	VAL	engineered mutation	UNP A3CWV0
C	229	VAL	ALA	engineered mutation	UNP A3CWV0
D	40	ASN	ASP	engineered mutation	UNP A3CWV0
D	42	SER	GLU	engineered mutation	UNP A3CWV0
D	147	GLU	ALA	engineered mutation	UNP A3CWV0
D	148	PRO	VAL	engineered mutation	UNP A3CWV0
D	229	VAL	ALA	engineered mutation	UNP A3CWV0

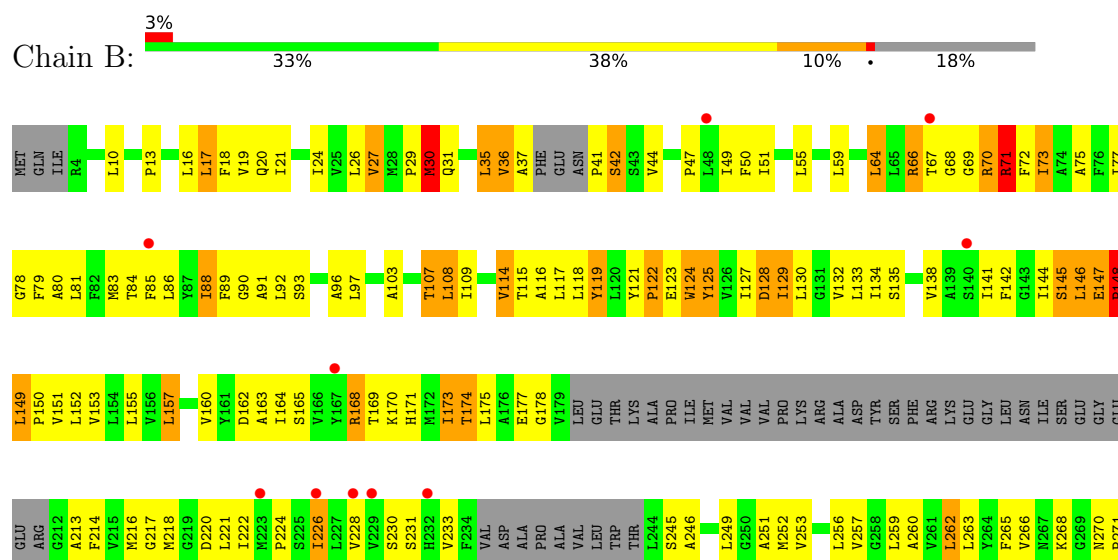
3 Residue-property plots

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: Putative uncharacterized protein

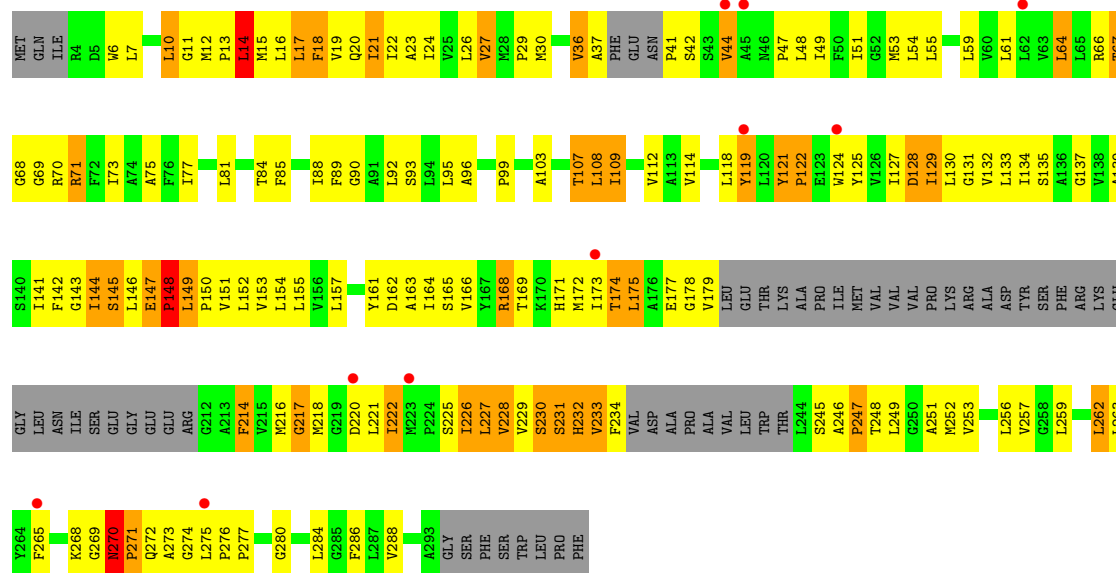


• Molecule 1: Putative uncharacterized protein

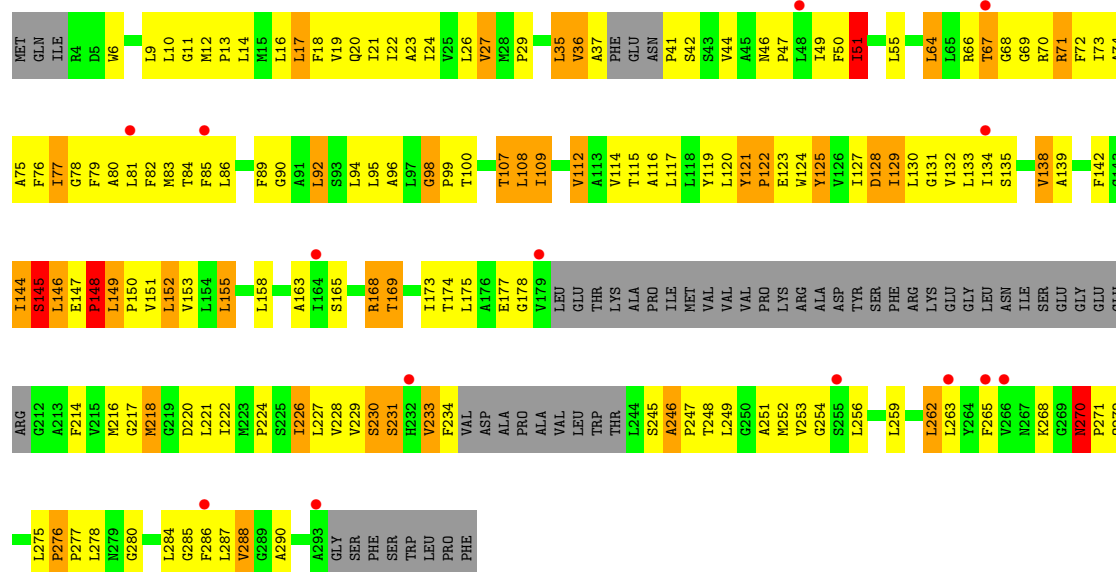




• Molecule 1: Putative uncharacterized protein



• Molecule 1: Putative uncharacterized protein



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	114.08Å 170.03Å 181.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	31.58 – 3.80 31.58 – 3.80	Depositor EDS
% Data completeness (in resolution range)	84.7 (31.58-3.80) 84.1 (31.58-3.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.18 (at 3.75Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8_1069)	Depositor
R, R_{free}	0.307 , 0.308 0.311 , 0.315	Depositor DCC
R_{free} test set	650 reflections (4.24%)	wwPDB-VP
Wilson B-factor (Å ²)	122.8	Xtriage
Anisotropy	0.572	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , 148.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	7199	wwPDB-VP
Average B, all atoms (Å ²)	184.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 14.35% 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	0.66	0/1839	1.20	12/2511 (0.5%)
1	B	0.67	0/1836	1.20	16/2507 (0.6%)
1	C	0.61	0/1836	1.23	19/2507 (0.8%)
1	D	0.67	0/1836	1.24	17/2507 (0.7%)
All	All	0.65	0/7347	1.22	64/10032 (0.6%)

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
1	C	0	1
1	D	0	1
All	All	0	4

There are no bond length outliers.

All (64) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	125	TYR	N-CA-C	-9.36	101.63	113.23
1	D	249	LEU	N-CA-C	-9.12	101.34	111.28
1	A	125	TYR	N-CA-C	-9.11	102.29	113.50
1	B	125	TYR	N-CA-C	-8.67	102.47	113.23
1	C	27	VAL	N-CA-C	-8.29	102.46	110.42
1	B	71	ARG	N-CA-C	-8.18	102.53	114.39
1	A	270	ASN	CA-C-N	8.13	130.01	119.84
1	A	270	ASN	C-N-CA	8.13	130.01	119.84
1	D	125	TYR	N-CA-C	-7.43	104.02	113.23
1	A	30	MET	N-CA-C	7.37	120.36	111.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	160	VAL	N-CA-C	7.30	118.10	110.72
1	C	270	ASN	CA-C-N	7.30	128.97	119.84
1	C	270	ASN	C-N-CA	7.30	128.97	119.84
1	D	71	ARG	N-CA-C	-7.24	103.89	114.39
1	C	249	LEU	N-CA-C	-6.93	103.72	111.28
1	D	51	ILE	N-CA-C	6.51	117.26	110.62
1	A	276	PRO	N-CA-C	6.45	118.57	110.70
1	A	114	VAL	N-CA-C	6.35	116.46	110.30
1	B	27	VAL	N-CA-C	-6.33	104.35	110.42
1	C	127	ILE	N-CA-C	6.32	116.49	110.42
1	B	146	LEU	N-CA-C	6.29	124.19	110.80
1	C	222	ILE	CB-CA-C	-6.23	104.00	111.97
1	C	7	LEU	CA-C-N	-6.17	113.59	119.76
1	C	7	LEU	C-N-CA	-6.17	113.59	119.76
1	C	232	HIS	N-CA-C	-6.15	105.04	112.54
1	D	146	LEU	N-CA-C	6.13	123.86	110.80
1	C	14	LEU	N-CA-C	-6.12	104.52	111.07
1	D	270	ASN	CA-C-N	5.97	127.31	119.84
1	D	270	ASN	C-N-CA	5.97	127.31	119.84
1	D	98	GLY	CA-C-N	5.97	127.30	119.84
1	D	98	GLY	C-N-CA	5.97	127.30	119.84
1	A	27	VAL	CB-CA-C	-5.97	104.33	111.97
1	D	248	THR	N-CA-C	-5.85	106.05	112.72
1	A	24	ILE	N-CA-C	5.85	116.49	110.82
1	D	112	VAL	CB-CA-C	-5.85	104.48	111.97
1	B	30	MET	N-CA-C	5.83	118.58	111.82
1	B	213	ALA	N-CA-C	-5.75	106.11	113.01
1	C	147	GLU	CA-C-N	5.74	127.02	119.84
1	C	147	GLU	C-N-CA	5.74	127.02	119.84
1	A	147	GLU	CA-C-N	5.73	127.01	119.84
1	A	147	GLU	C-N-CA	5.73	127.01	119.84
1	C	121	TYR	CA-C-N	5.69	126.95	119.84
1	C	121	TYR	C-N-CA	5.69	126.95	119.84
1	D	112	VAL	N-CA-C	-5.65	105.00	110.42
1	D	276	PRO	N-CA-C	5.60	117.53	110.70
1	B	114	VAL	N-CA-C	5.51	115.64	110.30
1	B	51	ILE	N-CA-C	5.50	116.23	110.62
1	A	112	VAL	CB-CA-C	-5.43	105.02	111.81
1	C	10	LEU	N-CA-C	-5.40	106.56	114.39
1	D	27	VAL	N-CA-C	-5.34	105.29	110.42
1	B	276	PRO	N-CA-C	5.32	117.19	110.70
1	C	54	LEU	N-CA-C	-5.30	105.58	111.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	147	GLU	CA-C-N	5.28	126.44	119.84
1	B	147	GLU	C-N-CA	5.28	126.44	119.84
1	C	247	PRO	N-CA-C	-5.28	108.64	114.92
1	D	246	ALA	CA-C-N	-5.23	114.93	121.00
1	D	246	ALA	C-N-CA	-5.23	114.93	121.00
1	B	260	ALA	N-CA-C	-5.21	105.49	111.07
1	A	53	MET	N-CA-C	5.13	117.61	111.71
1	B	73	ILE	CB-CA-C	-5.11	105.28	111.87
1	D	145	SER	N-CA-C	-5.04	103.88	110.53
1	C	53	MET	N-CA-C	5.03	117.66	111.82
1	B	276	PRO	CA-C-N	-5.03	113.86	119.19
1	B	276	PRO	C-N-CA	-5.03	113.86	119.19

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	145	SER	Peptide
1	B	145	SER	Peptide
1	C	145	SER	Peptide
1	D	145	SER	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	1802	0	1949	134	3
1	B	1799	0	1940	136	0
1	C	1799	0	1940	145	0
1	D	1799	0	1940	124	0
All	All	7199	0	7769	519	3

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 35.

All (519) 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:36:VAL:CG2	1:B:144:ILE:HG22	1.65	1.25
1:C:36:VAL:CG2	1:C:144:ILE:HG22	1.68	1.21
1:C:36:VAL:HG12	1:C:37:ALA:H	1.01	1.17
1:B:36:VAL:HG22	1:B:144:ILE:HG22	1.18	1.15
1:C:36:VAL:HG13	1:C:145:SER:HA	1.15	1.13
1:B:36:VAL:CG1	1:B:37:ALA:H	1.61	1.12
1:A:36:VAL:CG1	1:A:37:ALA:H	1.63	1.10
1:D:36:VAL:CG1	1:D:37:ALA:H	1.62	1.10
1:A:36:VAL:HG12	1:A:37:ALA:H	0.97	1.09
1:B:36:VAL:HG13	1:B:145:SER:HA	1.31	1.09
1:A:36:VAL:HG12	1:A:37:ALA:N	1.61	1.08
1:D:36:VAL:HG12	1:D:37:ALA:N	1.60	1.08
1:B:36:VAL:CG2	1:B:144:ILE:CG2	2.32	1.07
1:C:36:VAL:HG22	1:C:144:ILE:CG2	1.84	1.06
1:C:36:VAL:HG12	1:C:37:ALA:N	1.66	1.06
1:B:36:VAL:HG12	1:B:37:ALA:N	1.59	1.05
1:D:36:VAL:HG12	1:D:37:ALA:H	0.90	1.04
1:B:36:VAL:HG12	1:B:37:ALA:H	0.88	1.04
1:C:36:VAL:CG1	1:C:37:ALA:H	1.71	1.04
1:C:36:VAL:HG22	1:C:144:ILE:HG22	1.02	1.02
1:A:64:LEU:HD21	1:A:73:ILE:HA	1.42	1.01
1:C:36:VAL:CG2	1:C:144:ILE:CG2	2.37	1.00
1:B:36:VAL:HG22	1:B:144:ILE:CG2	1.93	0.97
1:A:36:VAL:CG1	1:A:146:LEU:HD23	1.99	0.91
1:C:30:MET:HE1	1:C:141:ILE:HA	1.51	0.91
1:A:161:TYR:OH	1:D:71:ARG:NH1	2.05	0.89
1:B:36:VAL:HG23	1:B:144:ILE:HG22	1.53	0.89
1:C:36:VAL:CG1	1:C:145:SER:HA	2.04	0.86
1:A:36:VAL:HG13	1:A:146:LEU:HD23	1.55	0.85
1:A:70:ARG:HH12	1:B:168:ARG:HH12	1.24	0.84
1:D:36:VAL:HG13	1:D:145:SER:HA	1.59	0.84
1:A:86:LEU:HD21	1:A:108:LEU:HA	1.57	0.83
1:C:36:VAL:HG13	1:C:145:SER:CA	2.05	0.81
1:B:30:MET:HE1	1:B:141:ILE:HA	1.61	0.81
1:B:36:VAL:HG23	1:B:144:ILE:CG2	2.09	0.80
1:A:262:LEU:HD23	1:A:277:PRO:HG2	1.62	0.80
1:B:41:PRO:C	1:B:146:LEU:HG	2.06	0.80
1:B:42:SER:HA	1:B:146:LEU:HB3	1.64	0.80
1:C:64:LEU:HD21	1:C:73:ILE:HA	1.63	0.79
1:B:128:ASP:N	1:B:128:ASP:OD2	2.15	0.79
1:B:220:ASP:HB2	1:B:275:LEU:HD21	1.66	0.78
1:D:262:LEU:HD23	1:D:277:PRO:HG2	1.66	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:253:VAL:HA	1:B:256:LEU:HD12	1.66	0.77
1:D:121:TYR:HD1	1:D:124:TRP:HZ2	1.31	0.76
1:C:139:ALA:HA	1:C:226:ILE:HD12	1.68	0.76
1:D:124:TRP:HA	1:D:127:ILE:HG12	1.67	0.76
1:A:36:VAL:HA	1:A:145:SER:HA	1.68	0.75
1:D:148:PRO:O	1:D:151:VAL:N	2.20	0.75
1:B:71:ARG:O	1:B:75:ALA:N	2.20	0.74
1:D:86:LEU:HD21	1:D:108:LEU:HA	1.70	0.74
1:B:70:ARG:HH12	1:C:169:THR:HG21	1.50	0.74
1:D:253:VAL:HA	1:D:256:LEU:HD12	1.69	0.73
1:D:151:VAL:HG11	1:D:227:LEU:HB2	1.71	0.72
1:B:16:LEU:HD23	1:B:256:LEU:HD23	1.71	0.72
1:C:118:LEU:HD12	1:C:130:LEU:HD12	1.70	0.71
1:B:80:ALA:HA	1:B:83:MET:HE2	1.71	0.71
1:D:135:SER:HA	1:D:222:ILE:HD11	1.72	0.71
1:C:36:VAL:CG1	1:C:146:LEU:HD22	2.20	0.71
1:D:64:LEU:HD11	1:D:73:ILE:HG23	1.73	0.70
1:D:41:PRO:C	1:D:146:LEU:HG	2.17	0.69
1:A:36:VAL:HG11	1:A:146:LEU:HD23	1.73	0.69
1:A:148:PRO:O	1:A:151:VAL:N	2.24	0.69
1:C:144:ILE:HA	1:C:233:VAL:HG21	1.75	0.69
1:A:36:VAL:HG11	1:A:41:PRO:HD2	1.74	0.69
1:D:165:SER:OG	1:D:168:ARG:NH2	2.26	0.68
1:C:30:MET:HE2	1:C:144:ILE:HD12	1.76	0.67
1:D:128:ASP:OD2	1:D:128:ASP:N	2.27	0.67
1:A:139:ALA:HA	1:A:226:ILE:HD12	1.77	0.66
1:B:36:VAL:HG22	1:B:144:ILE:C	2.19	0.66
1:A:36:VAL:CG1	1:A:41:PRO:HD2	2.25	0.66
1:D:155:LEU:HD11	1:D:224:PRO:HA	1.76	0.66
1:C:165:SER:O	1:C:169:THR:HG23	1.96	0.66
1:A:129:ILE:HG22	1:A:130:LEU:HD23	1.78	0.65
1:D:138:VAL:HB	1:D:222:ILE:HD13	1.77	0.65
1:D:90:GLY:O	1:D:94:LEU:HB2	1.97	0.65
1:C:142:PHE:HB2	1:C:226:ILE:HD11	1.79	0.65
1:B:118:LEU:HD12	1:B:130:LEU:HD12	1.80	0.64
1:C:71:ARG:HG2	1:D:168:ARG:HH22	1.62	0.64
1:D:268:LYS:HE2	1:D:272:GLN:HE22	1.62	0.64
1:D:24:ILE:HD13	1:D:229:VAL:HG22	1.80	0.64
1:A:220:ASP:HB2	1:A:275:LEU:HD21	1.79	0.64
1:C:89:PHE:HD1	1:C:92:LEU:HD12	1.63	0.64
1:B:19:VAL:HG22	1:B:132:VAL:HA	1.81	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:121:TYR:HA	1:C:124:TRP:CZ2	2.33	0.63
1:C:129:ILE:HG22	1:C:130:LEU:HD23	1.80	0.63
1:D:246:ALA:HB1	1:D:247:PRO:HD2	1.81	0.63
1:B:17:LEU:HD23	1:B:256:LEU:HD11	1.80	0.63
1:D:218:MET:HA	1:D:221:LEU:HB3	1.81	0.63
1:B:20:GLN:NE2	1:B:252:MET:HG2	2.13	0.63
1:A:63:VAL:O	1:A:67:THR:OG1	2.12	0.63
1:D:142:PHE:HB2	1:D:226:ILE:HD11	1.81	0.63
1:B:85:PHE:HE1	1:B:133:LEU:HD23	1.62	0.62
1:A:13:PRO:HB3	1:A:259:LEU:HD22	1.81	0.62
1:B:142:PHE:HB2	1:B:226:ILE:HD11	1.79	0.62
1:B:13:PRO:HB3	1:B:259:LEU:HD22	1.81	0.62
1:B:265:PHE:O	1:B:268:LYS:HG2	1.99	0.62
1:C:41:PRO:C	1:C:146:LEU:HG	2.24	0.62
1:D:254:GLY:HA3	1:D:285:GLY:HA3	1.82	0.62
1:A:265:PHE:HB3	1:A:272:GLN:OE1	1.99	0.61
1:D:42:SER:HA	1:D:146:LEU:HB3	1.82	0.61
1:C:21:ILE:HG22	1:C:22:ILE:HD13	1.82	0.61
1:C:108:LEU:O	1:C:112:VAL:HG23	2.00	0.61
1:D:29:PRO:HA	1:D:95:LEU:HD22	1.82	0.61
1:B:135:SER:HA	1:B:222:ILE:HD11	1.82	0.61
1:A:64:LEU:HD11	1:A:73:ILE:HG23	1.81	0.61
1:A:89:PHE:HD1	1:A:92:LEU:HD12	1.65	0.61
1:C:20:GLN:CD	1:C:252:MET:HG2	2.26	0.61
1:C:231:SER:HB2	1:C:248:THR:HG21	1.82	0.61
1:B:122:PRO:HD2	1:B:124:TRP:CE2	2.35	0.60
1:C:147:GLU:HB3	1:C:148:PRO:HD2	1.83	0.60
1:C:36:VAL:HG23	1:C:144:ILE:HG22	1.76	0.60
1:A:68:GLY:HA3	1:A:73:ILE:HD11	1.83	0.60
1:A:108:LEU:O	1:A:112:VAL:HG23	2.02	0.60
1:B:116:ALA:HA	1:C:164:ILE:HD11	1.82	0.60
1:C:16:LEU:HA	1:C:221:LEU:HD11	1.82	0.60
1:D:68:GLY:HA3	1:D:73:ILE:HD11	1.84	0.60
1:B:36:VAL:CG2	1:B:144:ILE:HG23	2.30	0.60
1:A:251:ALA:HB2	1:A:286:PHE:HB2	1.83	0.60
1:C:253:VAL:HA	1:C:256:LEU:HD12	1.83	0.59
1:D:64:LEU:HD21	1:D:73:ILE:HA	1.84	0.59
1:D:69:GLY:HA2	1:D:72:PHE:HB3	1.84	0.59
1:A:229:VAL:O	1:A:231:SER:N	2.35	0.59
1:B:171:HIS:HA	1:B:174:THR:HG23	1.85	0.59
1:D:149:LEU:HB3	1:D:150:PRO:HD3	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:42:SER:HA	1:A:146:LEU:HB3	1.84	0.59
1:A:120:LEU:HD21	1:B:164:ILE:HG12	1.84	0.59
1:A:23:ALA:O	1:A:27:VAL:HG23	2.02	0.58
1:A:72:PHE:CE2	1:B:59:LEU:HD23	2.38	0.58
1:D:16:LEU:HD23	1:D:256:LEU:HD23	1.85	0.58
1:D:151:VAL:CG1	1:D:227:LEU:HB2	2.33	0.58
1:D:24:ILE:O	1:D:27:VAL:N	2.36	0.58
1:B:41:PRO:O	1:B:146:LEU:HG	2.04	0.58
1:B:170:LYS:HZ1	1:B:271:PRO:C	2.10	0.58
1:B:122:PRO:HD2	1:B:124:TRP:CZ2	2.39	0.58
1:C:15:MET:SD	1:C:128:ASP:HB3	2.45	0.57
1:A:24:ILE:HD13	1:A:229:VAL:HG22	1.86	0.57
1:B:262:LEU:HD21	1:B:274:GLY:HA2	1.85	0.57
1:D:75:ALA:O	1:D:78:GLY:N	2.38	0.57
1:B:138:VAL:HB	1:B:222:ILE:HD13	1.87	0.56
1:C:36:VAL:HG23	1:C:144:ILE:CG2	2.32	0.56
1:C:121:TYR:HA	1:C:124:TRP:HZ2	1.68	0.56
1:D:108:LEU:O	1:D:112:VAL:HG23	2.06	0.56
1:D:220:ASP:HB2	1:D:275:LEU:HD21	1.87	0.56
1:D:139:ALA:HA	1:D:226:ILE:HD12	1.88	0.56
1:A:135:SER:HA	1:A:222:ILE:HD11	1.87	0.56
1:C:90:GLY:HA2	1:C:107:THR:HG21	1.88	0.56
1:B:165:SER:O	1:B:169:THR:HG23	2.06	0.56
1:C:265:PHE:HB3	1:C:272:GLN:CD	2.31	0.56
1:D:47:PRO:O	1:D:50:PHE:HB3	2.05	0.56
1:A:89:PHE:HZ	1:A:114:VAL:HG21	1.70	0.55
1:A:165:SER:O	1:A:169:THR:HG23	2.07	0.55
1:A:85:PHE:HE1	1:A:133:LEU:HD23	1.70	0.55
1:A:128:ASP:OD2	1:A:128:ASP:N	2.40	0.55
1:B:89:PHE:HA	1:B:92:LEU:HB2	1.88	0.55
1:B:121:TYR:HA	1:B:124:TRP:CZ2	2.41	0.55
1:C:85:PHE:HE1	1:C:133:LEU:HD23	1.71	0.55
1:C:225:SER:HA	1:C:228:VAL:HB	1.89	0.55
1:D:26:LEU:C	1:D:29:PRO:HD2	2.31	0.55
1:B:69:GLY:HA2	1:B:72:PHE:HB3	1.89	0.55
1:A:70:ARG:HH22	1:B:168:ARG:CZ	2.20	0.55
1:C:24:ILE:O	1:C:27:VAL:HB	2.07	0.55
1:B:81:LEU:O	1:B:84:THR:OG1	2.18	0.55
1:B:149:LEU:HB3	1:B:150:PRO:HD3	1.87	0.55
1:D:163:ALA:HB2	1:D:276:PRO:HG2	1.88	0.55
1:B:220:ASP:O	1:B:224:PRO:HD3	2.08	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:259:LEU:HB2	1:B:278:LEU:HD11	1.89	0.55
1:C:131:GLY:O	1:C:135:SER:OG	2.11	0.54
1:D:42:SER:HB2	1:D:146:LEU:HD12	1.87	0.54
1:A:158:LEU:HG	1:A:223:MET:HE1	1.89	0.54
1:B:35:LEU:HD23	1:B:141:ILE:HG12	1.89	0.54
1:B:42:SER:OG	1:B:147:GLU:HB2	2.07	0.54
1:C:6:TRP:O	1:C:10:LEU:HB2	2.08	0.54
1:C:36:VAL:HG22	1:C:144:ILE:C	2.31	0.54
1:D:129:ILE:HG22	1:D:130:LEU:HD23	1.88	0.54
1:B:26:LEU:C	1:B:29:PRO:HD2	2.32	0.54
1:B:147:GLU:HB3	1:B:148:PRO:HD2	1.90	0.54
1:D:13:PRO:HB3	1:D:259:LEU:HD22	1.90	0.54
1:B:89:PHE:HD1	1:B:92:LEU:HD12	1.73	0.54
1:C:147:GLU:O	1:C:150:PRO:HD2	2.08	0.54
1:B:97:LEU:HB2	1:B:103:ALA:HB2	1.89	0.54
1:C:226:ILE:O	1:C:230:SER:N	2.33	0.54
1:C:231:SER:HA	1:C:234:PHE:HB3	1.88	0.54
1:C:84:THR:O	1:C:88:ILE:HG13	2.08	0.54
1:A:48:LEU:HD23	1:D:83:MET:HG2	1.88	0.54
1:D:85:PHE:HE1	1:D:133:LEU:HD23	1.73	0.54
1:A:168:ARG:HH12	1:A:169:THR:HG22	1.73	0.54
1:B:148:PRO:O	1:B:151:VAL:N	2.41	0.53
1:D:24:ILE:HA	1:D:27:VAL:HG23	1.90	0.53
1:D:280:GLY:O	1:D:284:LEU:HG	2.08	0.53
1:A:70:ARG:O	1:A:70:ARG:NE	2.39	0.53
1:A:146:LEU:HD12	1:A:147:GLU:HG2	1.90	0.53
1:C:44:VAL:HG12	1:C:150:PRO:HA	1.89	0.53
1:D:146:LEU:HD12	1:D:147:GLU:HG2	1.90	0.53
1:C:13:PRO:HB3	1:C:259:LEU:HD22	1.89	0.53
1:B:71:ARG:NH1	1:C:161:TYR:OH	2.35	0.53
1:C:27:VAL:HG22	1:C:229:VAL:HG13	1.91	0.53
1:A:128:ASP:O	1:A:132:VAL:HG23	2.09	0.53
1:D:122:PRO:HG3	1:D:127:ILE:HD11	1.90	0.53
1:C:81:LEU:O	1:C:84:THR:OG1	2.24	0.53
1:A:169:THR:HA	1:A:172:MET:HB2	1.91	0.53
1:C:42:SER:HA	1:C:146:LEU:HB3	1.91	0.53
1:A:157:LEU:HD21	1:D:82:PHE:CE2	2.44	0.52
1:B:86:LEU:HD21	1:B:108:LEU:HA	1.91	0.52
1:C:20:GLN:O	1:C:24:ILE:HG12	2.09	0.52
1:A:15:MET:SD	1:A:128:ASP:HB3	2.50	0.52
1:B:36:VAL:HG22	1:B:144:ILE:O	2.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:68:GLY:HA3	1:C:73:ILE:HD11	1.90	0.52
1:D:69:GLY:HA3	1:D:73:ILE:HG13	1.91	0.52
1:D:122:PRO:HD2	1:D:124:TRP:CE2	2.45	0.52
1:A:30:MET:HE1	1:A:141:ILE:HA	1.92	0.52
1:A:265:PHE:O	1:A:268:LYS:HG2	2.09	0.52
1:B:75:ALA:O	1:B:78:GLY:N	2.43	0.52
1:C:128:ASP:OD2	1:C:128:ASP:N	2.39	0.52
1:C:142:PHE:HA	1:C:145:SER:HB3	1.92	0.52
1:B:90:GLY:HA2	1:B:107:THR:HG21	1.92	0.52
1:C:11:GLY:HA2	1:C:14:LEU:HB3	1.91	0.52
1:C:18:PHE:O	1:C:22:ILE:HG12	2.10	0.52
1:C:24:ILE:HA	1:C:27:VAL:HG23	1.92	0.51
1:D:51:ILE:HD11	1:D:158:LEU:HD21	1.92	0.51
1:A:148:PRO:HD3	1:A:230:SER:OG	2.10	0.51
1:C:151:VAL:HG11	1:C:227:LEU:HB2	1.91	0.51
1:A:82:PHE:HE1	1:A:112:VAL:HA	1.75	0.51
1:A:89:PHE:CZ	1:A:114:VAL:HG21	2.45	0.51
1:C:93:SER:HB2	1:C:103:ALA:HB1	1.92	0.51
1:C:128:ASP:O	1:C:132:VAL:HG23	2.09	0.51
1:B:20:GLN:CD	1:B:252:MET:HG2	2.36	0.51
1:B:218:MET:HA	1:B:221:LEU:HB3	1.91	0.51
1:D:36:VAL:CG1	1:D:37:ALA:N	2.32	0.51
1:A:20:GLN:CD	1:A:252:MET:HG2	2.35	0.51
1:A:36:VAL:HA	1:A:145:SER:CA	2.38	0.51
1:C:143:GLY:HA2	1:C:230:SER:HB2	1.92	0.51
1:D:71:ARG:O	1:D:75:ALA:N	2.38	0.51
1:A:36:VAL:CG1	1:A:37:ALA:N	2.31	0.51
1:D:148:PRO:O	1:D:149:LEU:C	2.54	0.51
1:A:24:ILE:HA	1:A:27:VAL:HG23	1.92	0.50
1:D:19:VAL:HG21	1:D:221:LEU:HD21	1.93	0.50
1:D:42:SER:N	1:D:146:LEU:HG	2.26	0.50
1:A:225:SER:HA	1:A:228:VAL:HB	1.93	0.50
1:D:44:VAL:O	1:D:47:PRO:HD2	2.11	0.50
1:C:44:VAL:O	1:C:47:PRO:HD2	2.10	0.50
1:B:70:ARG:NH1	1:C:169:THR:HG21	2.23	0.50
1:D:44:VAL:C	1:D:47:PRO:HD2	2.36	0.50
1:D:275:LEU:O	1:D:278:LEU:N	2.44	0.50
1:D:168:ARG:HH12	1:D:169:THR:HG22	1.77	0.50
1:A:16:LEU:HA	1:A:221:LEU:HD11	1.92	0.50
1:B:35:LEU:HD23	1:B:141:ILE:HG23	1.93	0.50
1:A:245:SER:H	1:A:249:LEU:HD11	1.77	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:20:GLN:O	1:B:24:ILE:HG12	2.12	0.50
1:A:107:THR:OG1	1:A:108:LEU:N	2.45	0.50
1:C:14:LEU:O	1:C:18:PHE:HB2	2.12	0.50
1:A:44:VAL:O	1:A:47:PRO:HD2	2.12	0.49
1:A:61:LEU:O	1:A:64:LEU:HB3	2.11	0.49
1:B:36:VAL:CG1	1:B:37:ALA:N	2.32	0.49
1:A:66:ARG:CZ	1:B:66:ARG:HE	2.26	0.49
1:C:139:ALA:HB2	1:C:222:ILE:HG23	1.95	0.49
1:D:168:ARG:HB3	1:D:168:ARG:NH1	2.28	0.49
1:B:274:GLY:O	1:B:277:PRO:HD2	2.13	0.49
1:A:47:PRO:O	1:A:50:PHE:HB3	2.13	0.49
1:A:89:PHE:HD2	1:A:111:ALA:N	2.10	0.49
1:C:89:PHE:HZ	1:C:114:VAL:HG21	1.76	0.49
1:C:93:SER:CB	1:C:103:ALA:HB1	2.43	0.49
1:C:218:MET:O	1:C:222:ILE:HG13	2.12	0.49
1:A:74:ALA:HA	1:A:77:ILE:HD11	1.94	0.49
1:B:16:LEU:HD23	1:B:256:LEU:CD2	2.42	0.49
1:B:31:GLN:OE1	1:B:31:GLN:N	2.40	0.49
1:D:74:ALA:HA	1:D:77:ILE:HD11	1.95	0.49
1:A:36:VAL:HG11	1:A:41:PRO:CD	2.43	0.48
1:A:72:PHE:CZ	1:B:59:LEU:HD23	2.48	0.48
1:B:42:SER:HA	1:B:146:LEU:CB	2.39	0.48
1:B:128:ASP:O	1:B:132:VAL:HG23	2.13	0.48
1:A:36:VAL:HG13	1:A:146:LEU:CD2	2.35	0.48
1:B:129:ILE:HG22	1:B:130:LEU:HD23	1.95	0.48
1:C:36:VAL:CG1	1:C:146:LEU:CD2	2.89	0.48
1:C:47:PRO:HG3	1:C:154:LEU:HB2	1.96	0.48
1:A:64:LEU:HD13	1:A:65:LEU:HG	1.96	0.48
1:A:162:ASP:HB2	1:A:275:LEU:HD12	1.96	0.48
1:C:162:ASP:HB2	1:C:275:LEU:HD12	1.94	0.48
1:D:66:ARG:NH1	1:D:67:THR:OG1	2.46	0.48
1:D:230:SER:HA	1:D:233:VAL:HG12	1.95	0.48
1:C:66:ARG:HH12	1:C:67:THR:HG23	1.79	0.48
1:C:262:LEU:HD21	1:C:274:GLY:HA2	1.95	0.48
1:A:44:VAL:C	1:A:47:PRO:HD2	2.38	0.48
1:A:70:ARG:HH12	1:B:169:THR:HG22	1.78	0.48
1:A:245:SER:HB2	1:A:249:LEU:HD21	1.96	0.48
1:A:268:LYS:HE2	1:A:272:GLN:HE22	1.79	0.48
1:C:36:VAL:CG1	1:C:37:ALA:N	2.37	0.48
1:A:24:ILE:O	1:A:27:VAL:N	2.46	0.47
1:B:24:ILE:O	1:B:27:VAL:HB	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:47:PRO:O	1:B:50:PHE:HB3	2.13	0.47
1:B:274:GLY:C	1:B:277:PRO:HD2	2.39	0.47
1:D:148:PRO:HD3	1:D:230:SER:O	2.14	0.47
1:B:220:ASP:C	1:B:220:ASP:OD2	2.58	0.47
1:D:165:SER:O	1:D:169:THR:HG23	2.14	0.47
1:D:231:SER:HA	1:D:234:PHE:HB3	1.96	0.47
1:B:114:VAL:O	1:B:117:LEU:HB3	2.14	0.47
1:C:12:MET:HA	1:C:15:MET:HE3	1.96	0.47
1:C:71:ARG:O	1:C:75:ALA:N	2.44	0.47
1:C:168:ARG:HH12	1:C:169:THR:HG22	1.79	0.47
1:D:20:GLN:O	1:D:24:ILE:HG12	2.14	0.47
1:C:88:ILE:O	1:C:92:LEU:HG	2.13	0.47
1:B:64:LEU:HD21	1:B:73:ILE:HA	1.96	0.47
1:C:231:SER:OG	1:C:232:HIS:N	2.47	0.47
1:D:80:ALA:O	1:D:84:THR:HG23	2.15	0.47
1:B:170:LYS:NZ	1:B:271:PRO:HG2	2.30	0.47
1:C:19:VAL:HG22	1:C:132:VAL:HG13	1.97	0.47
1:C:61:LEU:O	1:C:64:LEU:HB3	2.14	0.47
1:C:163:ALA:HB2	1:C:276:PRO:HG2	1.97	0.47
1:C:227:LEU:O	1:C:231:SER:HB3	2.15	0.47
1:D:24:ILE:O	1:D:27:VAL:HB	2.15	0.47
1:A:64:LEU:HD23	1:A:72:PHE:CE2	2.50	0.47
1:B:36:VAL:CG1	1:B:146:LEU:HD22	2.44	0.47
1:B:124:TRP:HA	1:B:127:ILE:HG12	1.97	0.47
1:C:217:GLY:O	1:C:220:ASP:OD2	2.33	0.47
1:D:76:PHE:C	1:D:76:PHE:CD2	2.93	0.47
1:B:24:ILE:HA	1:B:27:VAL:HG23	1.97	0.47
1:C:129:ILE:HD13	1:C:129:ILE:HA	1.76	0.46
1:A:68:GLY:HA2	1:A:69:GLY:HA3	1.61	0.46
1:A:131:GLY:O	1:A:135:SER:OG	2.17	0.46
1:C:36:VAL:HG11	1:C:146:LEU:HD22	1.94	0.46
1:A:36:VAL:CG1	1:A:146:LEU:CD2	2.83	0.46
1:C:109:ILE:HA	1:C:109:ILE:HD13	1.67	0.46
1:B:115:THR:O	1:B:119:TYR:HB2	2.16	0.46
1:B:251:ALA:HA	1:B:285:GLY:HA3	1.97	0.46
1:C:166:VAL:HG11	1:C:273:ALA:HA	1.98	0.46
1:D:42:SER:HA	1:D:146:LEU:CB	2.43	0.46
1:A:35:LEU:HB3	1:A:36:VAL:H	1.48	0.46
1:A:70:ARG:HE	1:A:70:ARG:C	2.23	0.46
1:A:175:LEU:HD22	1:A:179:VAL:HB	1.98	0.46
1:B:162:ASP:HB2	1:B:275:LEU:HD12	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:85:PHE:CE1	1:C:133:LEU:HD23	2.49	0.46
1:D:151:VAL:HG11	1:D:227:LEU:CB	2.42	0.46
1:C:24:ILE:HD13	1:C:229:VAL:HG22	1.97	0.46
1:D:148:PRO:HA	1:D:151:VAL:HB	1.98	0.46
1:A:93:SER:HB3	1:A:107:THR:HG22	1.97	0.45
1:A:289:GLY:HA2	1:A:292:LEU:HD12	1.97	0.45
1:C:20:GLN:HE22	1:C:252:MET:HA	1.82	0.45
1:C:23:ALA:HB1	1:C:229:VAL:HG21	1.97	0.45
1:D:134:ILE:O	1:D:138:VAL:HG23	2.15	0.45
1:A:13:PRO:HD3	1:A:259:LEU:HD21	1.97	0.45
1:D:116:ALA:O	1:D:120:LEU:HG	2.16	0.45
1:B:13:PRO:HD3	1:B:259:LEU:HD21	1.98	0.45
1:B:29:PRO:HB2	1:B:91:ALA:O	2.16	0.45
1:B:173:ILE:HD12	1:B:173:ILE:HA	1.72	0.45
1:D:23:ALA:O	1:D:27:VAL:HG23	2.17	0.45
1:A:64:LEU:C	1:A:64:LEU:HD22	2.41	0.45
1:A:75:ALA:O	1:A:78:GLY:N	2.49	0.45
1:D:20:GLN:CD	1:D:252:MET:HG2	2.41	0.45
1:D:89:PHE:HZ	1:D:114:VAL:HG21	1.81	0.45
1:D:109:ILE:HD13	1:D:109:ILE:HA	1.70	0.45
1:A:26:LEU:O	1:A:30:MET:HG3	2.17	0.45
1:A:85:PHE:HD2	1:A:115:THR:HG1	1.62	0.45
1:A:90:GLY:HA2	1:A:107:THR:HG21	1.99	0.45
1:B:275:LEU:HB2	1:B:276:PRO:HD3	1.98	0.45
1:C:148:PRO:O	1:C:151:VAL:N	2.50	0.45
1:A:79:PHE:CD1	1:A:80:ALA:N	2.85	0.45
1:C:36:VAL:CG2	1:C:144:ILE:HG23	2.40	0.45
1:C:64:LEU:HD11	1:C:73:ILE:HG23	1.99	0.45
1:C:122:PRO:HD2	1:C:124:TRP:CE2	2.51	0.45
1:C:268:LYS:HG3	1:C:269:GLY:O	2.17	0.45
1:A:72:PHE:CZ	1:B:59:LEU:HB2	2.52	0.44
1:A:218:MET:HA	1:A:221:LEU:HB3	1.98	0.44
1:C:27:VAL:CG2	1:C:229:VAL:HG13	2.47	0.44
1:C:246:ALA:HB1	1:C:247:PRO:HD2	1.99	0.44
1:D:89:PHE:HB2	1:D:107:THR:HB	1.98	0.44
1:A:76:PHE:O	1:A:79:PHE:HB3	2.16	0.44
1:C:232:HIS:HA	1:C:246:ALA:HB2	1.99	0.44
1:B:280:GLY:O	1:B:284:LEU:HG	2.17	0.44
1:B:70:ARG:HH22	1:C:169:THR:HG22	1.83	0.44
1:C:92:LEU:HD23	1:C:92:LEU:HA	1.76	0.44
1:A:72:PHE:O	1:A:75:ALA:HB3	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:88:ILE:HD12	1:B:133:LEU:HG	2.00	0.44
1:B:246:ALA:O	1:B:249:LEU:HG	2.18	0.44
1:C:118:LEU:HD23	1:C:119:TYR:HD1	1.83	0.44
1:C:275:LEU:HD23	1:C:275:LEU:HA	1.74	0.44
1:B:123:GLU:O	1:B:125:TYR:N	2.51	0.44
1:D:125:TYR:HA	1:D:128:ASP:OD2	2.18	0.44
1:D:251:ALA:HB2	1:D:286:PHE:HB2	1.99	0.44
1:A:80:ALA:HA	1:A:83:MET:HE2	1.99	0.44
1:B:30:MET:HE2	1:B:144:ILE:HD12	2.00	0.44
1:B:138:VAL:HA	1:B:141:ILE:HD12	2.00	0.44
1:C:284:LEU:HD23	1:C:284:LEU:HA	1.83	0.44
1:D:82:PHE:HE1	1:D:112:VAL:HA	1.82	0.44
1:A:263:LEU:O	1:A:267:ASN:HB2	2.18	0.44
1:A:270:ASN:HA	1:A:271:PRO:HD2	1.65	0.44
1:B:89:PHE:O	1:B:107:THR:HG22	2.18	0.44
1:B:118:LEU:HA	1:B:122:PRO:HG3	2.00	0.44
1:B:118:LEU:CD1	1:B:130:LEU:HD12	2.45	0.44
1:C:88:ILE:HD11	1:C:134:ILE:HD13	1.99	0.43
1:A:71:ARG:O	1:A:75:ALA:N	2.51	0.43
1:C:103:ALA:O	1:C:107:THR:HG23	2.18	0.43
1:C:149:LEU:HB3	1:C:150:PRO:HD3	2.00	0.43
1:A:42:SER:CA	1:A:146:LEU:HB3	2.48	0.43
1:B:30:MET:HB3	1:B:30:MET:HE3	1.58	0.43
1:B:134:ILE:O	1:B:138:VAL:HG23	2.18	0.43
1:C:175:LEU:HD22	1:C:179:VAL:HB	1.99	0.43
1:A:64:LEU:HD22	1:A:64:LEU:O	2.18	0.43
1:A:148:PRO:O	1:A:149:LEU:C	2.61	0.43
1:B:171:HIS:HA	1:B:174:THR:CG2	2.47	0.43
1:C:64:LEU:HD22	1:C:64:LEU:O	2.19	0.43
1:A:30:MET:HE1	1:A:140:SER:C	2.43	0.43
1:A:70:ARG:HB3	1:A:71:ARG:HE	1.84	0.43
1:B:121:TYR:HA	1:B:124:TRP:HZ2	1.84	0.43
1:B:148:PRO:O	1:B:149:LEU:C	2.61	0.43
1:B:270:ASN:HA	1:B:271:PRO:HD2	1.81	0.43
1:A:14:LEU:HA	1:A:14:LEU:HD22	1.77	0.43
1:A:53:MET:HE3	1:A:53:MET:HB3	1.94	0.43
1:A:127:ILE:HD13	1:A:127:ILE:HA	1.85	0.43
1:B:83:MET:HG2	1:C:48:LEU:HD21	2.00	0.43
1:C:231:SER:H	1:C:233:VAL:H	1.66	0.43
1:D:35:LEU:HB3	1:D:36:VAL:H	1.67	0.43
1:D:122:PRO:HG2	1:D:124:TRP:CG	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:131:GLY:O	1:D:135:SER:OG	2.26	0.43
1:B:36:VAL:HG13	1:B:146:LEU:HD22	2.00	0.43
1:B:275:LEU:O	1:B:278:LEU:N	2.49	0.43
1:C:275:LEU:HB2	1:C:276:PRO:HD3	2.01	0.43
1:C:280:GLY:O	1:C:284:LEU:HG	2.19	0.43
1:D:165:SER:HG	1:D:168:ARG:NH2	2.15	0.43
1:A:122:PRO:HB2	1:A:123:GLU:H	1.69	0.43
1:B:93:SER:O	1:B:97:LEU:N	2.52	0.43
1:C:30:MET:HE3	1:C:30:MET:HB3	1.69	0.43
1:D:92:LEU:HA	1:D:92:LEU:HD23	1.67	0.43
1:D:128:ASP:O	1:D:132:VAL:HG23	2.19	0.43
1:D:288:VAL:C	1:D:290:ALA:H	2.26	0.43
1:A:219:GLY:O	1:A:222:ILE:HB	2.18	0.43
1:C:217:GLY:O	1:C:220:ASP:N	2.52	0.43
1:D:6:TRP:O	1:D:10:LEU:HB2	2.18	0.43
1:B:42:SER:HB2	1:B:147:GLU:HG2	2.00	0.42
1:C:89:PHE:HA	1:C:92:LEU:HB2	2.01	0.42
1:C:270:ASN:HA	1:C:271:PRO:HD2	1.67	0.42
1:A:104:ALA:HA	1:A:107:THR:HG23	2.00	0.42
1:A:152:LEU:HD22	1:A:152:LEU:HA	1.76	0.42
1:C:169:THR:HA	1:C:172:MET:HB2	2.00	0.42
1:D:68:GLY:HA2	1:D:69:GLY:HA3	1.81	0.42
1:A:6:TRP:O	1:A:10:LEU:HB2	2.19	0.42
1:A:9:LEU:O	1:A:263:LEU:HD12	2.19	0.42
1:B:275:LEU:HD23	1:B:275:LEU:HA	1.75	0.42
1:C:26:LEU:HD23	1:C:26:LEU:HA	1.79	0.42
1:D:90:GLY:HA2	1:D:107:THR:HG21	1.99	0.42
1:A:24:ILE:O	1:A:27:VAL:HB	2.19	0.42
1:A:34:GLY:O	1:A:35:LEU:HB2	2.19	0.42
1:A:168:ARG:NH1	1:A:169:THR:HG22	2.35	0.42
1:C:68:GLY:HA2	1:C:69:GLY:HA3	1.50	0.42
1:D:51:ILE:HD11	1:D:158:LEU:CD2	2.50	0.42
1:D:270:ASN:HA	1:D:271:PRO:HD2	1.81	0.42
1:D:278:LEU:HA	1:D:278:LEU:HD13	1.82	0.42
1:A:263:LEU:HD22	1:A:263:LEU:HA	1.86	0.42
1:B:29:PRO:HB2	1:B:91:ALA:C	2.45	0.42
1:B:70:ARG:CZ	1:B:71:ARG:HG3	2.49	0.42
1:C:29:PRO:HB3	1:C:95:LEU:HD22	2.02	0.42
1:C:89:PHE:CZ	1:C:114:VAL:HG21	2.55	0.42
1:D:17:LEU:HD22	1:D:17:LEU:HA	1.81	0.42
1:D:11:GLY:O	1:D:14:LEU:HB3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:129:ILE:HD13	1:D:129:ILE:HA	1.83	0.42
1:A:88:ILE:O	1:A:92:LEU:HG	2.18	0.42
1:B:92:LEU:HD23	1:B:92:LEU:HA	1.76	0.42
1:C:143:GLY:HA3	1:C:229:VAL:CG1	2.50	0.42
1:D:98:GLY:O	1:D:100:THR:N	2.53	0.42
1:D:123:GLU:HB3	1:D:125:TYR:CE2	2.55	0.42
1:B:165:SER:OG	1:B:168:ARG:NH2	2.52	0.41
1:B:281:GLY:O	1:B:284:LEU:HB2	2.20	0.41
1:C:214:PHE:CD2	1:C:218:MET:HG2	2.55	0.41
1:D:121:TYR:HA	1:D:122:PRO:HD2	1.89	0.41
1:D:122:PRO:HB2	1:D:123:GLU:H	1.67	0.41
1:C:229:VAL:O	1:C:231:SER:N	2.50	0.41
1:D:152:LEU:HD22	1:D:152:LEU:HA	1.82	0.41
1:A:281:GLY:O	1:A:284:LEU:HB2	2.20	0.41
1:B:79:PHE:CD1	1:B:80:ALA:N	2.88	0.41
1:B:117:LEU:HD13	1:B:118:LEU:N	2.36	0.41
1:D:24:ILE:C	1:D:27:VAL:H	2.28	0.41
1:D:246:ALA:HB1	1:D:247:PRO:CD	2.49	0.41
1:B:10:LEU:HD23	1:B:10:LEU:HA	1.71	0.41
1:B:287:LEU:HD23	1:B:287:LEU:HA	1.86	0.41
1:B:59:LEU:HD13	1:B:59:LEU:O	2.21	0.41
1:C:134:ILE:O	1:C:137:GLY:N	2.53	0.41
1:D:13:PRO:HB3	1:D:259:LEU:CD2	2.51	0.41
1:D:36:VAL:HG22	1:D:144:ILE:HG22	2.02	0.41
1:D:287:LEU:HD23	1:D:287:LEU:HA	1.88	0.41
1:A:30:MET:HE2	1:A:144:ILE:HD12	2.03	0.41
1:B:36:VAL:CG1	1:B:146:LEU:CD2	2.98	0.41
1:B:157:LEU:HD12	1:B:157:LEU:HA	1.78	0.41
1:B:163:ALA:HB2	1:B:276:PRO:HG2	2.03	0.41
1:C:20:GLN:NE2	1:C:252:MET:HA	2.35	0.41
1:C:122:PRO:HD2	1:C:124:TRP:CZ2	2.55	0.41
1:A:114:VAL:HG13	1:A:130:LEU:HD13	2.02	0.41
1:A:147:GLU:HB3	1:A:148:PRO:HD2	2.03	0.41
1:C:251:ALA:HB2	1:C:286:PHE:HB2	2.03	0.41
1:A:51:ILE:HG22	1:D:79:PHE:HD2	1.86	0.41
1:A:70:ARG:HH22	1:B:168:ARG:NH1	2.19	0.41
1:A:85:PHE:HD2	1:A:115:THR:OG1	2.04	0.41
1:B:68:GLY:HA2	1:B:69:GLY:HA3	1.71	0.41
1:B:122:PRO:HB2	1:B:123:GLU:H	1.69	0.41
1:C:11:GLY:C	1:C:14:LEU:HB3	2.46	0.41
1:C:17:LEU:O	1:C:21:ILE:HD13	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:171:HIS:HA	1:C:174:THR:HG23	2.03	0.41
1:C:246:ALA:C	1:C:248:THR:H	2.29	0.41
1:D:114:VAL:O	1:D:117:LEU:HB3	2.20	0.41
1:A:30:MET:HE1	1:A:141:ILE:N	2.36	0.41
1:A:86:LEU:HA	1:A:86:LEU:HD23	1.81	0.41
1:C:121:TYR:HD1	1:C:124:TRP:CZ2	2.39	0.41
1:C:274:GLY:O	1:C:277:PRO:HD2	2.21	0.41
1:C:143:GLY:N	1:C:226:ILE:HG13	2.36	0.40
1:D:22:ILE:O	1:D:26:LEU:HB2	2.21	0.40
1:A:7:LEU:HB2	1:A:8:PRO:HD3	2.03	0.40
1:A:30:MET:HB3	1:A:30:MET:HE3	1.70	0.40
1:A:146:LEU:HD13	1:A:147:GLU:OE2	2.20	0.40
1:B:35:LEU:O	1:B:36:VAL:HB	2.22	0.40
1:C:88:ILE:HD13	1:C:133:LEU:HG	2.03	0.40
1:C:121:TYR:C	1:C:124:TRP:HE1	2.29	0.40
1:D:168:ARG:HB3	1:D:168:ARG:CZ	2.52	0.40
1:A:36:VAL:HA	1:A:145:SER:CB	2.50	0.40
1:A:109:ILE:HD13	1:A:109:ILE:HA	1.74	0.40
1:C:64:LEU:CD2	1:C:73:ILE:HA	2.44	0.40
1:D:10:LEU:HA	1:D:10:LEU:HD23	1.79	0.40
1:D:81:LEU:HD13	1:D:115:THR:HG23	2.03	0.40
1:D:214:PHE:O	1:D:218:MET:HG3	2.21	0.40
1:A:41:PRO:N	1:A:146:LEU:HG	2.37	0.40
1:B:13:PRO:O	1:B:16:LEU:HB3	2.21	0.40
1:D:9:LEU:HD22	1:D:12:MET:HE3	2.04	0.40
1:D:265:PHE:HB3	1:D:272:GLN:OE1	2.22	0.40

All (3) 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:292:LEU:CD2	1:A:292:LEU:CD2[3_654]	1.64	0.56
1:A:292:LEU:CD1	1:A:292:LEU:CD2[3_654]	1.90	0.30
1:A:292:LEU:CG	1:A:292:LEU:CD2[3_654]	2.00	0.20

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	238/301 (79%)	196 (82%)	30 (13%)	12 (5%)	1	17
1	B	238/301 (79%)	198 (83%)	29 (12%)	11 (5%)	2	18
1	C	238/301 (79%)	196 (82%)	31 (13%)	11 (5%)	2	18
1	D	238/301 (79%)	198 (83%)	27 (11%)	13 (6%)	1	16
All	All	952/1204 (79%)	788 (83%)	117 (12%)	47 (5%)	1	17

All (47) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	36	VAL
1	A	96	ALA
1	B	36	VAL
1	C	36	VAL
1	D	36	VAL
1	A	230	SER
1	B	96	ALA
1	B	231	SER
1	C	96	ALA
1	C	122	PRO
1	C	230	SER
1	C	231	SER
1	D	96	ALA
1	D	217	GLY
1	D	231	SER
1	A	35	LEU
1	A	122	PRO
1	A	148	PRO
1	A	217	GLY
1	A	231	SER
1	B	35	LEU
1	B	122	PRO

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Mol	Chain	Res	Type
1	B	148	PRO
1	B	217	GLY
1	C	99	PRO
1	C	148	PRO
1	D	122	PRO
1	D	148	PRO
1	D	230	SER
1	A	99	PRO
1	A	178	GLY
1	B	178	GLY
1	B	230	SER
1	C	178	GLY
1	C	217	GLY
1	D	35	LEU
1	D	92	LEU
1	D	99	PRO
1	D	178	GLY
1	A	75	ALA
1	B	42	SER
1	B	124	TRP
1	A	270	ASN
1	C	270	ASN
1	D	270	ASN
1	C	271	PRO
1	D	121	TYR

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	188/236 (80%)	138 (73%)	50 (27%)	0	3
1	B	187/236 (79%)	145 (78%)	42 (22%)	1	5
1	C	187/236 (79%)	144 (77%)	43 (23%)	1	5
1	D	187/236 (79%)	148 (79%)	39 (21%)	1	7
All	All	749/944 (79%)	575 (77%)	174 (23%)	1	5

All (174) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	LEU
1	A	18	PHE
1	A	21	ILE
1	A	44	VAL
1	A	46	ASN
1	A	48	LEU
1	A	49	ILE
1	A	51	ILE
1	A	55	LEU
1	A	64	LEU
1	A	67	THR
1	A	70	ARG
1	A	71	ARG
1	A	77	ILE
1	A	79	PHE
1	A	88	ILE
1	A	107	THR
1	A	108	LEU
1	A	109	ILE
1	A	119	TYR
1	A	126	VAL
1	A	128	ASP
1	A	129	ILE
1	A	141	ILE
1	A	148	PRO
1	A	149	LEU
1	A	152	LEU
1	A	153	VAL
1	A	155	LEU
1	A	157	LEU
1	A	168	ARG
1	A	169	THR
1	A	173	ILE
1	A	174	THR
1	A	175	LEU
1	A	177	GLU
1	A	214	PHE
1	A	216	MET
1	A	226	ILE
1	A	227	LEU
1	A	228	VAL
1	A	232	HIS

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Mol	Chain	Res	Type
1	A	233	VAL
1	A	234	PHE
1	A	245	SER
1	A	253	VAL
1	A	257	VAL
1	A	262	LEU
1	A	263	LEU
1	A	288	VAL
1	B	17	LEU
1	B	18	PHE
1	B	21	ILE
1	B	30	MET
1	B	44	VAL
1	B	49	ILE
1	B	55	LEU
1	B	64	LEU
1	B	66	ARG
1	B	67	THR
1	B	70	ARG
1	B	71	ARG
1	B	77	ILE
1	B	88	ILE
1	B	107	THR
1	B	108	LEU
1	B	109	ILE
1	B	119	TYR
1	B	128	ASP
1	B	129	ILE
1	B	148	PRO
1	B	149	LEU
1	B	152	LEU
1	B	153	VAL
1	B	155	LEU
1	B	157	LEU
1	B	168	ARG
1	B	173	ILE
1	B	174	THR
1	B	175	LEU
1	B	177	GLU
1	B	214	PHE
1	B	216	MET
1	B	226	ILE

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Mol	Chain	Res	Type
1	B	228	VAL
1	B	233	VAL
1	B	245	SER
1	B	257	VAL
1	B	262	LEU
1	B	263	LEU
1	B	266	VAL
1	B	288	VAL
1	C	14	LEU
1	C	17	LEU
1	C	18	PHE
1	C	21	ILE
1	C	44	VAL
1	C	49	ILE
1	C	51	ILE
1	C	55	LEU
1	C	59	LEU
1	C	64	LEU
1	C	67	THR
1	C	70	ARG
1	C	71	ARG
1	C	77	ILE
1	C	107	THR
1	C	108	LEU
1	C	109	ILE
1	C	119	TYR
1	C	128	ASP
1	C	129	ILE
1	C	144	ILE
1	C	148	PRO
1	C	149	LEU
1	C	152	LEU
1	C	153	VAL
1	C	155	LEU
1	C	157	LEU
1	C	168	ARG
1	C	173	ILE
1	C	174	THR
1	C	175	LEU
1	C	177	GLU
1	C	214	PHE
1	C	216	MET

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Mol	Chain	Res	Type
1	C	226	ILE
1	C	227	LEU
1	C	228	VAL
1	C	233	VAL
1	C	245	SER
1	C	257	VAL
1	C	262	LEU
1	C	263	LEU
1	C	288	VAL
1	D	17	LEU
1	D	18	PHE
1	D	21	ILE
1	D	46	ASN
1	D	49	ILE
1	D	51	ILE
1	D	55	LEU
1	D	64	LEU
1	D	67	THR
1	D	70	ARG
1	D	77	ILE
1	D	107	THR
1	D	108	LEU
1	D	109	ILE
1	D	119	TYR
1	D	128	ASP
1	D	129	ILE
1	D	138	VAL
1	D	144	ILE
1	D	148	PRO
1	D	149	LEU
1	D	152	LEU
1	D	153	VAL
1	D	155	LEU
1	D	168	ARG
1	D	169	THR
1	D	173	ILE
1	D	174	THR
1	D	175	LEU
1	D	177	GLU
1	D	216	MET
1	D	218	MET
1	D	226	ILE

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Mol	Chain	Res	Type
1	D	228	VAL
1	D	233	VAL
1	D	245	SER
1	D	262	LEU
1	D	263	LEU
1	D	288	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	46	ASN
1	B	232	HIS
1	C	20	GLN
1	C	46	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	246/301 (81%)	0.30	15 (6%)	27 22	102, 158, 255, 434	0
1	B	246/301 (81%)	0.22	10 (4%)	41 30	144, 184, 265, 350	0
1	C	246/301 (81%)	0.22	10 (4%)	41 30	147, 196, 279, 349	0
1	D	246/301 (81%)	0.34	14 (5%)	29 23	97, 154, 257, 361	0
All	All	984/1204 (81%)	0.27	49 (4%)	34 25	97, 176, 268, 434	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	228	VAL	7.6
1	C	44	VAL	6.3
1	B	48	LEU	4.9
1	C	223	MET	4.6
1	A	232	HIS	4.6
1	D	266	VAL	3.6
1	D	293	ALA	3.4
1	D	286	PHE	3.3
1	D	134	ILE	3.2
1	D	179	VAL	3.2
1	B	140	SER	3.2
1	D	265	PHE	3.1
1	A	223	MET	3.1
1	A	123	GLU	3.0
1	D	263	LEU	3.0
1	D	85	PHE	2.9
1	A	142	PHE	2.8
1	A	286	PHE	2.8
1	B	232	HIS	2.8
1	D	232	HIS	2.8
1	A	265	PHE	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	119	TYR	2.8
1	C	220	ASP	2.7
1	B	85	PHE	2.7
1	A	224	PRO	2.6
1	B	67	THR	2.6
1	C	62	LEU	2.5
1	D	81	LEU	2.5
1	B	229	VAL	2.5
1	D	67	THR	2.4
1	C	119	TYR	2.4
1	A	35	LEU	2.3
1	C	265	PHE	2.3
1	A	62	LEU	2.3
1	B	223	MET	2.2
1	A	30	MET	2.2
1	A	95	LEU	2.2
1	D	164	ILE	2.1
1	A	158	LEU	2.1
1	A	33	ALA	2.1
1	C	45	ALA	2.1
1	C	173	ILE	2.1
1	D	48	LEU	2.1
1	C	124	TRP	2.0
1	C	275	LEU	2.0
1	A	266	VAL	2.0
1	B	167	TYR	2.0
1	B	226	ILE	2.0
1	D	255	SER	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.