



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:47 AM UTC

PDB ID : 9J0A / pdb_00009j0a
Title : Complex structure of ANKRD11/STAG2/RAD21
Authors : Liu, H.; Cai, Q.; Zhang, M.
Deposited on : 2024-08-02
Resolution : 3.30 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

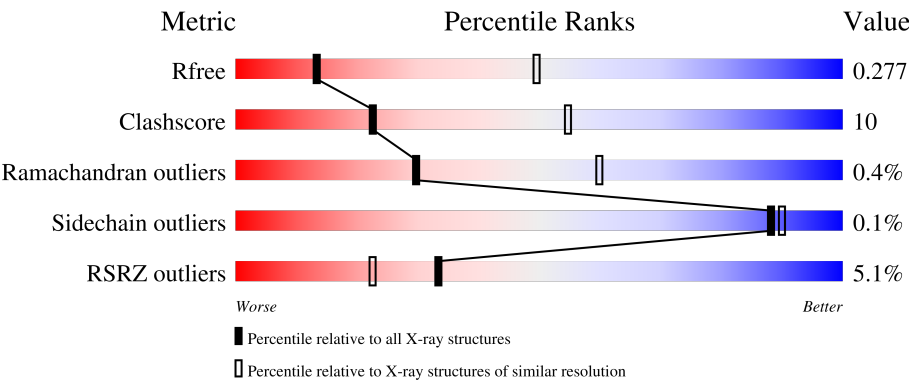
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)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

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1169 (3.32-3.28)
Clashscore	190562	1209 (3.32-3.28)
Ramachandran outliers	187476	1188 (3.32-3.28)
Sidechain outliers	187428	1187 (3.32-3.28)
RSRZ outliers	180081	1169 (3.32-3.28)

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	985	5% 69% 25% 6%
1	D	985	5% 70% 25% ..
2	B	143	% 41% 11% 48%
2	E	143	% 45% 6% 48%
3	C	41	7% 54% 5% 41%

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Mol	Chain	Length	Quality of chain
3	F	41	7% 39% 20% 41%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 16806 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 Cohesin subunit SA-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	926	Total	C	N	O	S	0	0	0
			7541	4800	1264	1423	54			
1	D	941	Total	C	N	O	S	0	0	0
			7635	4861	1279	1439	56			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	76	GLY	-	expression tag	UNP O35638
A	77	PRO	-	expression tag	UNP O35638
A	78	GLY	-	expression tag	UNP O35638
A	79	SER	-	expression tag	UNP O35638
D	76	GLY	-	expression tag	UNP O35638
D	77	PRO	-	expression tag	UNP O35638
D	78	GLY	-	expression tag	UNP O35638
D	79	SER	-	expression tag	UNP O35638

- Molecule 2 is a protein called 64-kDa C-terminal product.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	74	Total	C	N	O	S	0	0	0
			601	391	101	106	3			
2	E	74	Total	C	N	O	S	0	0	0
			601	391	101	106	3			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	278	GLY	-	expression tag	UNP Q61550
B	279	PRO	-	expression tag	UNP Q61550
B	280	GLY	-	expression tag	UNP Q61550
B	281	SER	-	expression tag	UNP Q61550

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Chain	Residue	Modelled	Actual	Comment	Reference
E	278	GLY	-	expression tag	UNP Q61550
E	279	PRO	-	expression tag	UNP Q61550
E	280	GLY	-	expression tag	UNP Q61550
E	281	SER	-	expression tag	UNP Q61550

- Molecule 3 is a protein called Ankyrin repeat domain-containing protein 11.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	24	Total	C	N	O	0	0	0
			207	127	32	48			
3	F	24	Total	C	N	O	0	0	0
			207	127	32	48			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	217	GLY	-	expression tag	UNP E9Q4F7
C	218	PRO	-	expression tag	UNP E9Q4F7
C	219	GLY	-	expression tag	UNP E9Q4F7
C	220	SER	-	expression tag	UNP E9Q4F7
F	217	GLY	-	expression tag	UNP E9Q4F7
F	218	PRO	-	expression tag	UNP E9Q4F7
F	219	GLY	-	expression tag	UNP E9Q4F7
F	220	SER	-	expression tag	UNP E9Q4F7

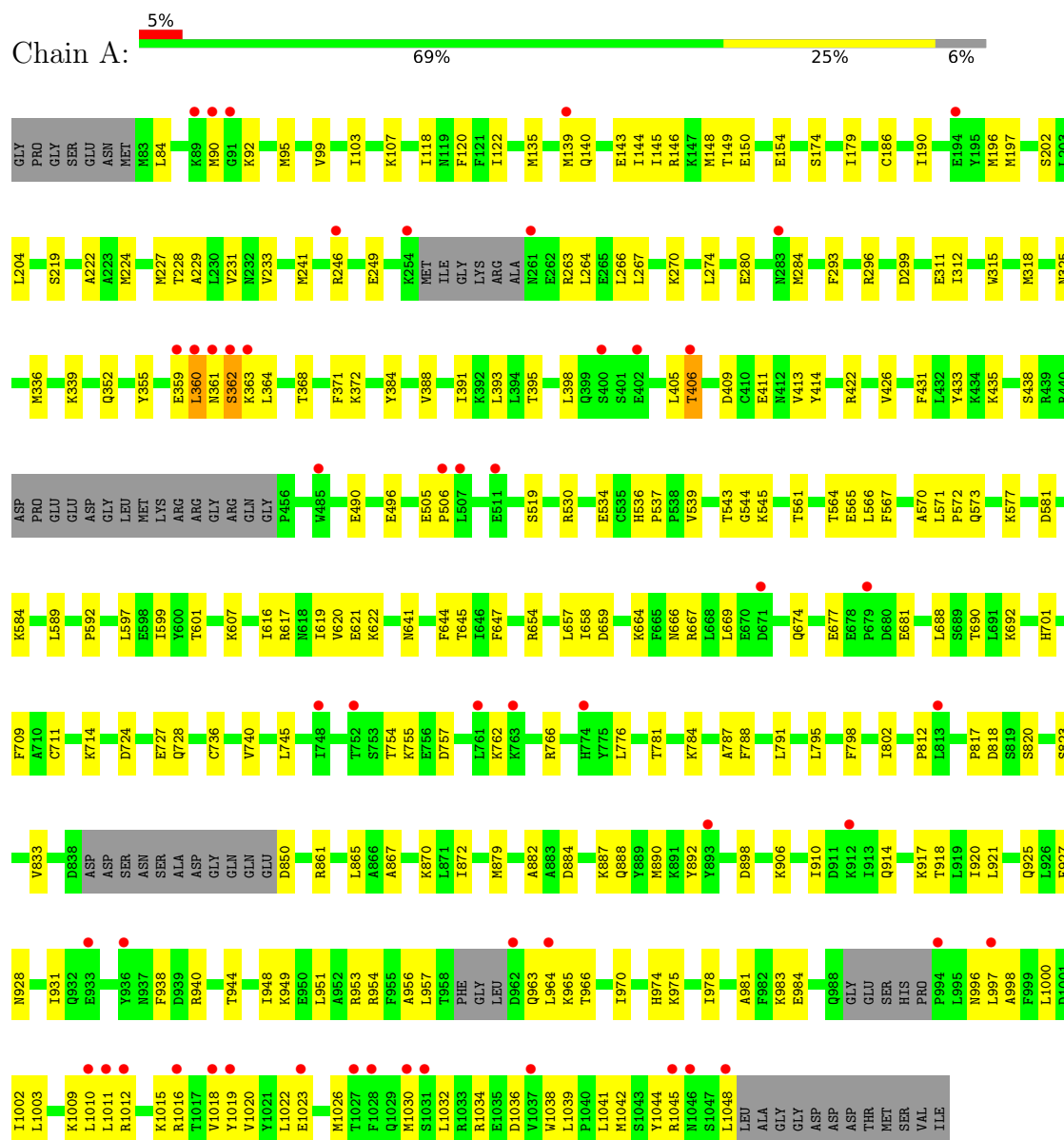
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	6	Total	O	0	0
			6	6		
4	C	1	Total	O	0	0
			1	1		
4	D	6	Total	O	0	0
			6	6		
4	E	1	Total	O	0	0
			1	1		

3 Residue-property plots

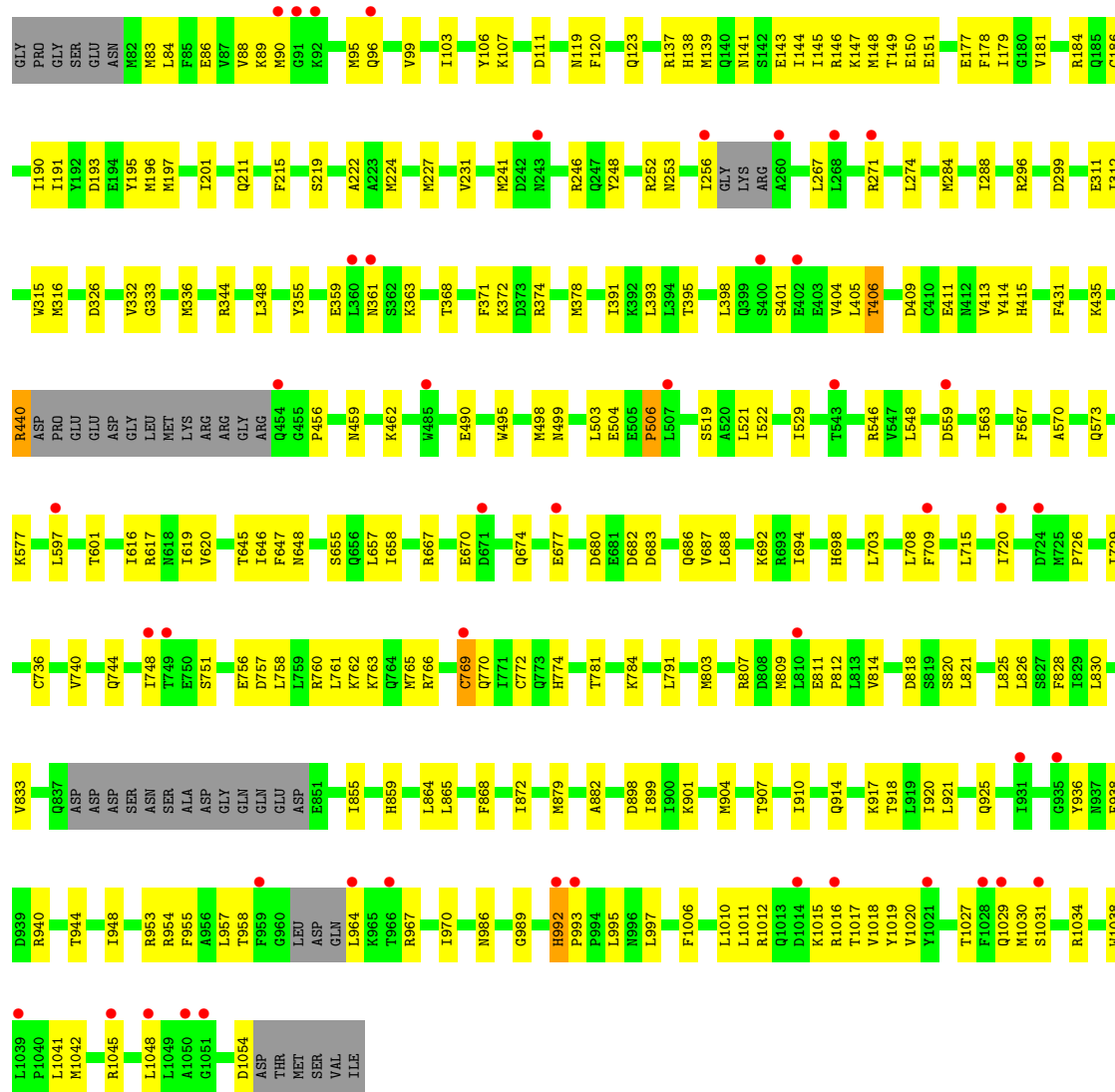
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Cohesin subunit SA-2

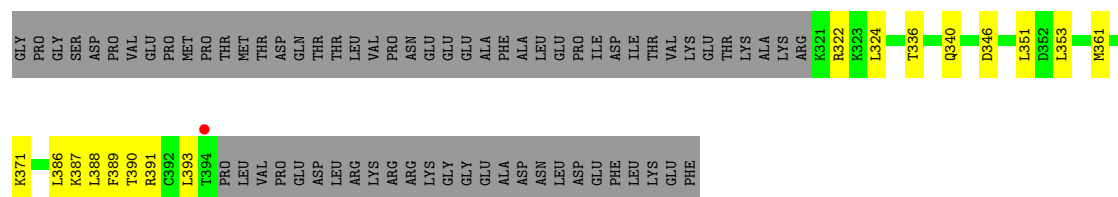
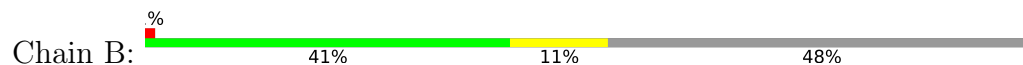


• Molecule 1: Cohesin subunit SA-2

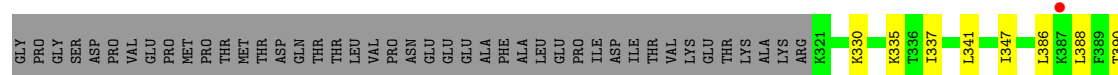


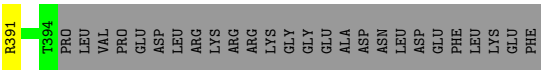


• Molecule 2: 64-kDa C-terminal product



• Molecule 2: 64-kDa C-terminal product





● Molecule 3: Ankyrin repeat domain-containing protein 11



● Molecule 3: Ankyrin repeat domain-containing protein 11



4 Data and refinement statistics

Property	Value	Source
Space group	I 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	280.54Å 48.77Å 290.91Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	31.88 – 3.30 31.88 – 3.30	Depositor EDS
% Data completeness (in resolution range)	99.8 (31.88-3.30) 99.8 (31.88-3.30)	Depositor EDS
R_{merge}	0.20	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.56 (at 3.33Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.233 , 0.279 0.233 , 0.277	Depositor DCC
R_{free} test set	3198 reflections (5.25%)	wwPDB-VP
Wilson B-factor (Å ²)	84.8	Xtriage
Anisotropy	0.431	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 78.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	0.040 for -l,k,h 0.034 for -l,-k,-h 0.046 for h,-k,-l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	16806	wwPDB-VP
Average B, all atoms (Å ²)	96.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.34% 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	0.16	0/7671	0.37	0/10338
1	D	0.16	0/7768	0.36	0/10471
2	B	0.13	0/612	0.37	0/826
2	E	0.13	0/612	0.36	0/826
3	C	0.10	0/211	0.36	0/285
3	F	0.12	0/211	0.37	0/285
All	All	0.16	0/17085	0.36	0/23031

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	D	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	440	ARG	Sidechain
1	D	992	HIS	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	7541	0	7539	178	0
1	D	7635	0	7633	161	1
2	B	601	0	643	14	0
2	E	601	0	643	6	0
3	C	207	0	177	2	0
3	F	207	0	177	5	1
4	A	6	0	0	0	0
4	C	1	0	0	0	0
4	D	6	0	0	0	0
4	E	1	0	0	0	0
All	All	16806	0	16812	352	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 10.

All (352) 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:938:PHE:O	1:A:940:ARG:NH2	2.13	0.81
1:A:1022:LEU:HG	1:A:1026:MET:HE1	1.64	0.79
2:B:351:LEU:HD22	2:B:353:LEU:HD12	1.65	0.79
1:D:141:ASN:ND2	3:F:242:ASP:OD1	2.17	0.77
1:A:325:ASN:HA	1:A:363:LYS:HG3	1.66	0.77
1:A:246:ARG:HH21	1:A:249:GLU:HB3	1.50	0.76
1:A:186:CYS:HB3	1:A:190:ILE:HB	1.68	0.76
1:D:879:MET:HE3	1:D:907:THR:HB	1.69	0.75
1:D:88:VAL:HG22	1:D:95:MET:HE1	1.68	0.73
1:D:440:ARG:NH2	1:D:490:GLU:OE2	2.21	0.73
1:D:1012:ARG:HH21	1:D:1054:ASP:HA	1.52	0.72
1:A:1034:ARG:HG3	1:A:1042:MET:HG3	1.71	0.72
1:A:669:LEU:HD12	1:A:711:CYS:HB3	1.70	0.72
1:D:405:LEU:O	1:D:406:THR:HG22	1.90	0.72
1:D:191:ILE:HG13	1:D:197:MET:HE3	1.72	0.71
1:D:709:PHE:HB2	1:D:740:VAL:HG11	1.73	0.70
1:D:252:ARG:HE	1:D:252:ARG:C	2.00	0.69
1:A:970:ILE:CD1	1:A:1009:LYS:HB2	2.21	0.69
1:D:1012:ARG:H	1:D:1012:ARG:HD2	1.58	0.68
1:A:398:LEU:HD11	1:A:435:LYS:HD3	1.76	0.68
1:A:519:SER:OG	1:A:584:LYS:NZ	2.22	0.68
1:D:197:MET:O	1:D:201:ILE:HD12	1.96	0.66
1:A:336:MET:HE1	1:A:371:PHE:HB3	1.78	0.66
1:A:228:THR:OG1	1:A:311:GLU:OE1	2.12	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:405:LEU:O	1:A:406:THR:HG22	1.97	0.65
1:D:917:LYS:O	1:D:920:ILE:HG22	1.97	0.65
1:D:964:LEU:HB2	1:D:967:ARG:HH12	1.61	0.65
1:A:95:MET:HE3	1:A:174:SER:HB3	1.78	0.64
1:A:917:LYS:O	1:A:920:ILE:HG22	1.98	0.64
1:D:252:ARG:NH2	1:D:256:ILE:HA	2.13	0.64
1:A:983:LYS:HD2	1:A:984:GLU:HG3	1.80	0.64
1:D:667:ARG:HH12	1:D:670:GLU:HG3	1.63	0.64
1:D:760:ARG:O	1:D:763:LYS:HG3	1.97	0.63
1:A:149:THR:O	2:B:322:ARG:NH1	2.31	0.63
1:A:1010:LEU:HD21	1:A:1018:VAL:HG21	1.81	0.63
1:D:95:MET:HE2	1:D:178:PHE:HB2	1.80	0.63
1:A:228:THR:HG23	1:A:318:MET:HE1	1.80	0.63
1:D:967:ARG:H	1:D:967:ARG:HD2	1.64	0.63
1:A:135:MET:HA	1:A:139:MET:HB2	1.81	0.63
1:A:1012:ARG:HG3	1:A:1016:ARG:NH1	2.13	0.62
1:D:119:ASN:O	1:D:123:GLN:HG3	1.99	0.62
1:A:823:SER:HB2	1:D:1045:ARG:NH2	2.14	0.62
1:D:181:VAL:HG22	1:D:184:ARG:HH12	1.64	0.61
1:D:215:PHE:O	1:D:219:SER:OG	2.15	0.61
1:D:601:THR:HG21	1:D:645:THR:OG1	2.01	0.61
1:A:890:MET:HB2	1:A:951:LEU:HD13	1.82	0.61
1:A:1036:ASP:O	1:A:1039:LEU:HD23	2.01	0.61
1:A:296:ARG:NH1	1:A:299:ASP:OD2	2.32	0.61
1:A:135:MET:HG2	1:A:139:MET:SD	2.40	0.61
1:A:867:ALA:HA	1:A:870:LYS:HE2	1.82	0.61
1:A:997:LEU:HD11	1:A:1041:LEU:HB2	1.83	0.61
1:A:1034:ARG:HG3	1:A:1042:MET:CG	2.30	0.60
1:D:186:CYS:HB3	1:D:190:ILE:HB	1.82	0.60
1:A:1023:GLU:HA	1:A:1026:MET:SD	2.41	0.60
1:A:617:ARG:NH1	1:A:621:GLU:OE2	2.34	0.60
1:A:490:GLU:N	1:A:490:GLU:OE1	2.34	0.60
1:D:224:MET:HB3	1:D:311:GLU:HG3	1.83	0.60
1:D:336:MET:HE1	1:D:371:PHE:HB3	1.83	0.60
1:D:1034:ARG:HG2	1:D:1042:MET:HG3	1.84	0.60
1:A:536:HIS:HB2	1:A:544:GLY:HA2	1.83	0.60
1:A:543:THR:HG23	1:A:545:LYS:H	1.65	0.60
1:A:709:PHE:HB2	1:A:740:VAL:HG11	1.83	0.60
1:A:601:THR:HG21	1:A:645:THR:HB	1.83	0.59
1:A:953:ARG:HH21	1:A:954:ARG:HG2	1.67	0.59
1:A:617:ARG:O	1:A:620:VAL:HG22	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1034:ARG:O	1:A:1034:ARG:HG2	2.03	0.59
1:A:145:ILE:O	1:A:149:THR:HG23	2.03	0.59
1:A:981:ALA:HA	1:A:996:ASN:HB2	1.85	0.59
1:D:252:ARG:O	1:D:252:ARG:NE	2.31	0.59
1:D:398:LEU:HB2	1:D:405:LEU:HD11	1.83	0.59
1:A:293:PHE:CZ	1:A:312:ILE:HD12	2.38	0.58
1:A:1026:MET:HB2	1:A:1030:MET:HB2	1.85	0.58
1:A:872:ILE:HG21	1:A:882:ALA:HB2	1.86	0.58
1:A:280:GLU:HG3	1:A:284:MET:HE3	1.85	0.58
1:A:879:MET:HG2	1:A:918:THR:HG21	1.86	0.58
1:D:145:ILE:O	1:D:149:THR:HG23	2.03	0.58
1:D:414:TYR:OH	1:D:435:LYS:HG2	2.04	0.58
1:D:694:ILE:HG21	1:D:708:LEU:HD22	1.86	0.58
1:D:898:ASP:OD1	1:D:899:ILE:N	2.37	0.57
1:A:1034:ARG:HA	1:A:1038:TRP:HB2	1.86	0.57
1:A:1000:LEU:HD12	1:A:1041:LEU:HA	1.87	0.57
1:D:147:LYS:O	1:D:151:GLU:HG2	2.05	0.57
1:A:669:LEU:CD1	1:A:711:CYS:HB3	2.34	0.57
1:D:440:ARG:HH22	1:D:490:GLU:CD	2.12	0.57
1:A:664:LYS:HG2	1:A:690:THR:HG21	1.87	0.56
1:D:241:MET:HE2	1:D:274:LEU:HB3	1.86	0.56
1:A:391:ILE:O	1:A:395:THR:HG23	2.05	0.56
1:D:617:ARG:O	1:D:620:VAL:HG22	2.05	0.56
1:D:970:ILE:HD11	1:D:1006:PHE:HB3	1.88	0.56
1:D:111:ASP:OD2	1:D:195:TYR:OH	2.18	0.56
1:D:391:ILE:O	1:D:395:THR:HG23	2.06	0.56
1:D:296:ARG:NH1	1:D:299:ASP:OD2	2.38	0.56
1:D:698:HIS:CD2	1:D:703:LEU:HB2	2.41	0.56
1:A:224:MET:HB3	1:A:311:GLU:HG3	1.87	0.56
1:D:667:ARG:NH1	1:D:670:GLU:HG3	2.21	0.56
1:D:137:ARG:HG2	1:D:138:HIS:CD2	2.41	0.55
1:D:248:TYR:HA	1:D:267:LEU:HD13	1.89	0.55
1:D:683:ASP:O	1:D:687:VAL:HG12	2.05	0.55
1:A:190:ILE:HG22	1:A:196:MET:HG2	1.89	0.55
1:A:795:LEU:HD22	1:A:817:PRO:HG3	1.88	0.55
1:A:898:ASP:OD2	1:A:898:ASP:N	2.40	0.55
1:D:1011:LEU:HD12	1:D:1012:ARG:NH1	2.22	0.55
1:D:901:LYS:HA	1:D:904:MET:HE2	1.89	0.54
1:A:745:LEU:HD23	2:B:388:LEU:HD13	1.88	0.54
1:D:715:LEU:HD21	1:D:729:ILE:HG22	1.88	0.54
1:D:997:LEU:HD11	1:D:1041:LEU:HB2	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:833:VAL:HG11	1:D:865:LEU:HB2	1.90	0.54
1:D:954:ARG:O	1:D:958:THR:HG23	2.08	0.54
1:A:964:LEU:C	1:A:966:THR:H	2.16	0.54
1:A:674:GLN:HB3	1:A:677:GLU:HB2	1.89	0.53
2:B:346:ASP:OD1	2:B:346:ASP:N	2.37	0.53
1:D:106:TYR:OH	1:D:193:ASP:OD2	2.21	0.53
1:A:263:ARG:HA	1:A:266:LEU:HD12	1.88	0.53
1:D:519:SER:O	1:D:522:ILE:HG13	2.07	0.53
1:D:986:ASN:HB2	1:D:993:PRO:O	2.07	0.53
1:D:504:GLU:HG3	1:D:506:PRO:HD3	1.91	0.53
1:D:821:LEU:O	1:D:825:LEU:HG	2.08	0.53
1:D:103:ILE:HG22	1:D:107:LYS:HE2	1.89	0.53
1:D:898:ASP:OD1	1:D:899:ILE:HG13	2.08	0.53
1:A:754:THR:HG21	1:A:757:ASP:HB2	1.91	0.53
1:D:333:GLY:HA2	1:D:336:MET:HE2	1.91	0.52
1:A:940:ARG:NH1	1:A:998:ALA:HB1	2.24	0.52
1:A:1012:ARG:HG3	1:A:1016:ARG:HH11	1.73	0.52
1:A:496:GLU:HG2	1:A:566:LEU:HD11	1.92	0.52
1:A:970:ILE:HG23	1:A:1010:LEU:HD13	1.90	0.52
1:A:1032:LEU:HA	1:A:1034:ARG:HH11	1.75	0.52
1:A:762:LYS:O	1:A:766:ARG:HG3	2.10	0.52
1:A:144:ILE:O	1:A:148:MET:HG2	2.10	0.52
1:D:1017:THR:O	1:D:1020:VAL:HG22	2.09	0.52
1:D:495:TRP:HH2	1:D:559:ASP:HB2	1.76	0.51
1:A:409:ASP:O	1:A:413:VAL:HG23	2.10	0.51
1:A:888:GLN:HB3	1:A:892:TYR:HD1	1.75	0.51
1:D:1034:ARG:HA	1:D:1038:TRP:HB2	1.93	0.51
1:A:762:LYS:NZ	1:A:812:PRO:O	2.42	0.51
1:A:411:GLU:HA	1:A:414:TYR:CD2	2.45	0.51
1:A:654:ARG:HH11	1:A:701:HIS:CD2	2.29	0.51
1:A:561:THR:O	1:A:565:GLU:HG3	2.10	0.50
1:D:84:LEU:HD23	1:D:120:PHE:HB2	1.93	0.50
1:A:589:LEU:O	1:A:592:PRO:HD2	2.12	0.50
1:D:818:ASP:OD2	1:D:820:SER:OG	2.28	0.50
1:D:872:ILE:HG21	1:D:882:ALA:HB2	1.94	0.50
1:A:620:VAL:HG21	1:A:657:LEU:HD21	1.94	0.50
1:A:150:GLU:HA	2:B:322:ARG:HH12	1.76	0.50
1:A:724:ASP:OD1	1:A:724:ASP:N	2.37	0.50
1:A:154:GLU:OE1	1:A:154:GLU:N	2.42	0.49
1:A:938:PHE:HE1	1:A:940:ARG:HH12	1.59	0.49
1:D:409:ASP:O	1:D:413:VAL:HG23	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:83:MET:HA	1:D:86:GLU:HG2	1.93	0.49
1:D:940:ARG:HD3	1:D:940:ARG:N	2.27	0.49
1:D:744:GLN:O	1:D:748:ILE:HG12	2.12	0.49
1:A:368:THR:O	1:A:372:LYS:HB2	2.13	0.49
1:A:1032:LEU:HA	1:A:1034:ARG:NH1	2.28	0.49
1:D:751:SER:HA	1:D:809:MET:HE1	1.94	0.49
1:A:607:LYS:HD3	1:A:607:LYS:C	2.38	0.49
1:D:196:MET:SD	1:D:197:MET:HE2	2.53	0.49
1:A:818:ASP:OD2	1:A:820:SER:OG	2.29	0.49
1:D:762:LYS:NZ	1:D:812:PRO:O	2.46	0.49
1:A:736:CYS:O	1:A:740:VAL:HG23	2.13	0.48
1:A:918:THR:HA	1:A:921:LEU:HD12	1.96	0.48
1:D:1018:VAL:HG23	1:D:1048:LEU:HD13	1.93	0.48
1:A:1026:MET:HE2	1:A:1041:LEU:HD11	1.94	0.48
1:D:440:ARG:HD3	1:D:456:PRO:HB2	1.95	0.48
1:A:970:ILE:HD13	1:A:1009:LYS:HB2	1.94	0.48
1:A:352:GLN:HG3	1:A:393:LEU:HD13	1.95	0.48
1:D:989:GLY:HA3	1:D:992:HIS:CB	2.44	0.48
1:D:1029:GLN:HG2	1:D:1030:MET:N	2.28	0.48
1:A:820:SER:HB3	1:D:1019:TYR:CZ	2.49	0.48
1:A:581:ASP:CG	1:A:584:LYS:HG3	2.37	0.48
1:A:927:PHE:O	1:A:931:ILE:HG23	2.14	0.48
1:D:938:PHE:O	1:D:940:ARG:NH1	2.45	0.48
1:A:140:GLN:OE1	3:C:241:ASP:N	2.47	0.48
1:A:589:LEU:HD11	1:A:619:ILE:HD11	1.95	0.48
1:A:776:LEU:HD11	1:A:788:PHE:HB2	1.95	0.48
1:D:828:PHE:HE1	1:D:864:LEU:HD13	1.77	0.48
1:A:266:LEU:O	1:A:270:LYS:HG3	2.14	0.48
1:A:956:ALA:HB1	1:A:1009:LYS:HE3	1.95	0.48
1:A:103:ILE:HG22	1:A:107:LYS:HE3	1.96	0.48
1:A:573:GLN:O	1:A:577:LYS:HG2	2.14	0.48
1:D:769:CYS:O	1:D:772:CYS:HB2	2.14	0.48
1:A:202:SER:HA	3:C:244:HIS:ND1	2.29	0.47
1:A:567:PHE:HA	1:A:570:ALA:HB3	1.96	0.47
1:A:1030:MET:HG2	1:A:1038:TRP:HZ3	1.79	0.47
1:A:914:GLN:O	1:A:918:THR:HG23	2.14	0.47
1:A:949:LYS:HG3	1:A:1002:ILE:HD13	1.96	0.47
1:A:1015:LYS:HD3	1:A:1048:LEU:HA	1.97	0.47
1:D:655:SER:HA	1:D:658:ILE:HD12	1.95	0.47
1:A:622:LYS:HD2	1:A:622:LYS:O	2.14	0.47
1:A:1012:ARG:O	1:A:1016:ARG:HG2	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:374:ARG:O	1:D:378:MET:HG3	2.14	0.47
1:A:154:GLU:OE2	2:B:324:LEU:HB2	2.15	0.47
1:A:339:LYS:HE2	2:B:340:GLN:HG2	1.97	0.47
1:A:395:THR:HG22	1:A:431:PHE:HB2	1.96	0.47
1:D:765:MET:SD	1:D:766:ARG:NE	2.88	0.47
1:A:361:ASN:O	1:A:363:LYS:N	2.45	0.47
1:A:641:ASN:O	1:A:647:PHE:HB2	2.15	0.47
1:A:928:ASN:O	1:A:931:ILE:HG12	2.14	0.47
1:A:84:LEU:HD23	1:A:120:PHE:HB2	1.96	0.47
1:A:798:PHE:HA	1:A:802:ILE:HD12	1.97	0.46
1:A:833:VAL:HA	1:A:861:ARG:HG2	1.97	0.46
1:A:1020:VAL:O	1:A:1023:GLU:HG2	2.15	0.46
1:D:573:GLN:O	1:D:577:LYS:HG2	2.14	0.46
1:D:921:LEU:O	1:D:925:GLN:HG3	2.14	0.46
1:A:1019:TYR:OH	1:A:1045:ARG:HG2	2.15	0.46
1:D:688:LEU:HD13	1:D:729:ILE:HG12	1.98	0.46
1:D:756:GLU:OE2	1:D:756:GLU:N	2.47	0.46
1:D:757:ASP:HA	1:D:760:ARG:HG2	1.97	0.46
3:F:233:GLU:H	3:F:236:ARG:HH21	1.63	0.46
1:A:781:THR:HA	1:A:784:LYS:HD3	1.97	0.46
1:D:144:ILE:O	1:D:148:MET:HG2	2.15	0.46
1:D:803:MET:HA	1:D:807:ARG:HB2	1.98	0.46
1:A:204:LEU:O	1:A:219:SER:OG	2.33	0.46
1:A:754:THR:O	1:A:755:LYS:HD2	2.15	0.46
1:D:344:ARG:HD3	2:E:347:ILE:HD12	1.98	0.46
1:D:766:ARG:NH2	1:D:814:VAL:O	2.49	0.45
1:A:1019:TYR:OH	1:A:1045:ARG:NH1	2.50	0.45
1:D:227:MET:HB2	1:D:288:ILE:HG21	1.98	0.45
1:D:914:GLN:O	1:D:918:THR:HG23	2.16	0.45
1:D:139:MET:SD	1:D:143:GLU:HG2	2.56	0.45
1:A:229:ALA:O	1:A:233:VAL:HG23	2.16	0.45
1:A:666:ASN:HB3	1:A:667:ARG:NH2	2.32	0.45
1:A:884:ASP:O	1:A:887:LYS:HG2	2.17	0.45
2:B:387:LYS:O	2:B:391:ARG:HG2	2.16	0.45
1:D:855:ILE:HG22	1:D:859:HIS:CE1	2.51	0.45
1:D:1027:THR:O	1:D:1031:SER:N	2.45	0.45
1:A:241:MET:HE2	1:A:274:LEU:HB3	1.97	0.45
1:A:530:ARG:O	1:A:534:GLU:HG2	2.16	0.45
1:A:143:GLU:HG2	1:A:146:ARG:NH2	2.31	0.45
1:D:498:MET:HG2	1:D:521:LEU:HD11	1.98	0.45
1:D:736:CYS:O	1:D:740:VAL:HG23	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:90:MET:O	1:A:92:LYS:HG2	2.16	0.44
1:A:974:HIS:HA	1:A:1003:LEU:HD11	1.98	0.44
1:D:89:LYS:HG2	1:D:90:MET:HG2	1.98	0.44
1:D:146:ARG:O	1:D:150:GLU:HG3	2.16	0.44
1:D:411:GLU:HG3	1:D:415:HIS:CE1	2.52	0.44
1:A:246:ARG:NH2	1:A:249:GLU:HB3	2.27	0.44
1:A:362:SER:C	1:A:364:LEU:H	2.24	0.44
1:A:953:ARG:NH2	1:A:954:ARG:HG2	2.30	0.44
1:A:906:LYS:O	1:A:910:ILE:HG12	2.18	0.44
1:D:312:ILE:O	1:D:316:MET:HG3	2.17	0.44
1:D:682:ASP:O	1:D:686:GLN:HG2	2.18	0.44
2:E:337:ILE:O	2:E:341:LEU:HG	2.16	0.44
1:D:688:LEU:O	1:D:692:LYS:HG3	2.17	0.44
1:D:826:LEU:O	1:D:830:LEU:HG	2.17	0.44
1:D:758:LEU:HD21	1:D:812:PRO:HB2	2.00	0.44
2:B:371:LYS:HB3	2:B:371:LYS:HE3	1.86	0.44
1:D:348:LEU:HD23	1:D:348:LEU:HA	1.82	0.44
3:F:242:ASP:OD2	3:F:243:LYS:NZ	2.49	0.44
1:A:597:LEU:HB2	1:A:644:PHE:CD2	2.53	0.44
1:A:197:MET:HG2	1:A:284:MET:SD	2.58	0.43
1:A:787:ALA:O	1:A:791:LEU:HB2	2.18	0.43
1:A:888:GLN:HB3	1:A:892:TYR:CD1	2.53	0.43
1:D:791:LEU:HD21	1:D:821:LEU:HD11	2.00	0.43
1:A:539:VAL:HG12	2:B:361:MET:HG3	2.01	0.43
1:D:597:LEU:HD13	1:D:646:ILE:HD11	1.99	0.43
1:D:211:GLN:HG2	2:E:330:LYS:HD3	2.00	0.43
1:D:227:MET:O	1:D:231:VAL:HG23	2.18	0.43
1:A:983:LYS:NZ	1:A:984:GLU:OE2	2.46	0.43
1:A:1015:LYS:O	1:A:1019:TYR:HB2	2.18	0.43
1:A:1026:MET:HB2	1:A:1030:MET:CB	2.48	0.43
1:D:970:ILE:HG23	1:D:1010:LEU:HD12	2.00	0.43
1:A:571:LEU:HB3	1:A:572:PRO:HD3	2.01	0.43
1:D:944:THR:O	1:D:948:ILE:HG13	2.19	0.43
2:E:388:LEU:HD22	2:E:391:ARG:NH2	2.34	0.43
1:D:99:VAL:O	1:D:103:ILE:HG13	2.19	0.43
1:D:359:GLU:HB2	1:D:361:ASN:OD1	2.18	0.43
1:A:727:GLU:OE2	1:A:728:GLN:HG2	2.18	0.43
1:A:143:GLU:HG2	1:A:146:ARG:HH22	1.83	0.42
1:D:312:ILE:HA	1:D:315:TRP:CE3	2.54	0.42
1:D:907:THR:HA	1:D:910:ILE:HD12	2.01	0.42
1:A:887:LYS:HD3	1:A:925:GLN:OE1	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:499:ASN:O	1:D:503:LEU:HB2	2.19	0.42
1:D:616:ILE:O	1:D:620:VAL:HG13	2.19	0.42
1:D:766:ARG:HA	1:D:766:ARG:HD3	1.87	0.42
1:D:953:ARG:O	1:D:957:LEU:HG	2.19	0.42
3:F:246:LEU:HD23	3:F:246:LEU:HA	1.93	0.42
1:A:422:ARG:O	1:A:426:VAL:HG22	2.19	0.42
2:B:387:LYS:HB2	2:B:387:LYS:HE2	1.71	0.42
1:D:1010:LEU:O	1:D:1015:LYS:NZ	2.51	0.42
1:D:529:ILE:HG23	1:D:563:ILE:HD13	2.01	0.42
2:E:386:LEU:O	2:E:390:THR:HG23	2.19	0.42
1:A:1044:TYR:O	1:A:1048:LEU:HD23	2.19	0.42
1:D:953:ARG:HH21	1:D:954:ARG:HG3	1.84	0.42
1:D:955:PHE:O	1:D:958:THR:OG1	2.33	0.42
1:A:688:LEU:O	1:A:692:LYS:HG3	2.20	0.42
1:A:714:LYS:HB2	1:A:714:LYS:HE2	1.76	0.42
1:A:833:VAL:HG11	1:A:865:LEU:HB2	2.01	0.42
1:D:177:GLU:O	1:D:181:VAL:HG23	2.20	0.42
1:D:253:ASN:HA	1:D:256:ILE:HD12	2.02	0.42
1:D:1012:ARG:NH2	1:D:1054:ASP:O	2.53	0.42
1:A:99:VAL:O	1:A:103:ILE:HG13	2.19	0.42
1:A:433:TYR:O	1:A:438:SER:N	2.53	0.42
1:D:355:TYR:CD2	1:D:393:LEU:HD11	2.55	0.42
1:D:917:LYS:HE2	1:D:917:LYS:HB3	1.93	0.42
1:A:920:ILE:HD12	1:A:920:ILE:HA	1.88	0.41
1:A:964:LEU:C	1:A:966:THR:N	2.78	0.41
2:B:389:PHE:O	2:B:393:LEU:HG	2.20	0.41
1:A:545:LYS:HE2	1:A:545:LYS:HB2	1.92	0.41
1:A:564:THR:HB	1:A:599:ILE:HD12	2.01	0.41
1:D:241:MET:HG2	1:D:271:ARG:HD3	2.02	0.41
1:D:326:ASP:H	1:D:363:LYS:HE2	1.84	0.41
1:D:726:PRO:HB2	1:D:729:ILE:HG13	2.02	0.41
1:D:781:THR:HA	1:D:784:LYS:HD3	2.02	0.41
1:D:936:TYR:CE1	1:D:995:LEU:HB3	2.55	0.41
1:A:118:ILE:O	1:A:122:ILE:HG12	2.21	0.41
1:A:944:THR:O	1:A:948:ILE:HG13	2.20	0.41
2:B:336:THR:O	2:B:340:GLN:HG3	2.20	0.41
1:D:96:GLN:HA	1:D:181:VAL:HG11	2.02	0.41
1:A:384:TYR:O	1:A:388:VAL:HG23	2.20	0.41
1:D:1016:ARG:NH2	1:D:1019:TYR:HD2	2.19	0.41
1:A:90:MET:O	1:A:90:MET:HE2	2.19	0.41
1:A:1015:LYS:HA	1:A:1018:VAL:HG22	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:401:SER:OG	1:D:404:VAL:HG12	2.20	0.41
1:D:459:ASN:HA	1:D:462:LYS:HD3	2.03	0.41
1:D:803:MET:HE1	1:D:811:GLU:HG2	2.02	0.41
1:A:264:LEU:HA	1:A:267:LEU:HD12	2.02	0.41
1:A:616:ILE:O	1:A:620:VAL:HG13	2.20	0.41
1:A:818:ASP:HB2	1:D:1016:ARG:NH1	2.35	0.41
2:B:386:LEU:O	2:B:390:THR:HG23	2.20	0.41
1:D:680:ASP:OD2	1:D:683:ASP:N	2.47	0.41
2:E:335:LYS:HA	2:E:335:LYS:HD3	1.81	0.41
1:A:536:HIS:HA	1:A:537:PRO:HD3	1.92	0.41
1:A:681:GLU:H	1:A:681:GLU:HG3	1.67	0.41
1:D:332:VAL:O	1:D:336:MET:HG3	2.21	0.41
1:D:395:THR:HG22	1:D:431:PHE:HB2	2.01	0.41
1:D:567:PHE:HA	1:D:570:ALA:HB3	2.01	0.41
1:D:179:ILE:HD13	1:D:222:ALA:HA	2.03	0.41
1:D:197:MET:HG2	1:D:284:MET:SD	2.60	0.41
1:D:616:ILE:O	1:D:619:ILE:HG12	2.21	0.41
1:D:748:ILE:HG13	1:D:761:LEU:HD22	2.02	0.41
1:A:179:ILE:HD13	1:A:222:ALA:HA	2.02	0.41
1:A:963:GLN:HE22	1:A:1011:LEU:HD21	1.86	0.41
1:D:368:THR:O	1:D:372:LYS:HB2	2.20	0.41
1:D:674:GLN:HG3	1:D:677:GLU:H	1.86	0.41
1:D:720:ILE:HD13	1:D:720:ILE:HA	1.93	0.41
1:D:1042:MET:SD	1:D:1045:ARG:NH1	2.91	0.41
1:A:227:MET:O	1:A:231:VAL:HG23	2.21	0.40
1:A:355:TYR:CD2	1:A:393:LEU:HD11	2.56	0.40
1:A:658:ILE:O	1:A:659:ASP:C	2.63	0.40
1:D:147:LYS:HA	1:D:147:LYS:HD2	1.93	0.40
1:D:546:ARG:NH1	1:D:548:LEU:HA	2.36	0.40
3:F:232:ASP:O	3:F:234:GLN:N	2.48	0.40
1:A:360:LEU:O	1:A:364:LEU:HB2	2.21	0.40
1:A:975:LYS:O	1:A:978:ILE:HG12	2.22	0.40
1:A:1030:MET:HE3	1:A:1038:TRP:CZ3	2.56	0.40
1:A:312:ILE:HA	1:A:315:TRP:CE3	2.57	0.40
1:D:868:PHE:O	1:D:872:ILE:HG13	2.21	0.40
1:A:359:GLU:O	1:A:361:ASN:N	2.55	0.40
1:A:850:ASP:OD1	1:A:850:ASP:N	2.55	0.40
1:A:957:LEU:HD23	1:A:1009:LYS:HZ2	1.87	0.40
1:D:620:VAL:HG21	1:D:657:LEU:HD21	2.03	0.40
1:D:647:PHE:O	1:D:648:ASN:C	2.64	0.40
1:D:770:GLN:HE21	1:D:774:HIS:CE1	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:D:246:ARG:NH1	3:F:227:GLU:OE2[4_445]	2.17	0.03

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	914/985 (93%)	876 (96%)	32 (4%)	6 (1%)	18	49
1	D	931/985 (94%)	900 (97%)	29 (3%)	2 (0%)	43	71
2	B	72/143 (50%)	70 (97%)	2 (3%)	0	100	100
2	E	72/143 (50%)	71 (99%)	1 (1%)	0	100	100
3	C	22/41 (54%)	19 (86%)	3 (14%)	0	100	100
3	F	22/41 (54%)	16 (73%)	6 (27%)	0	100	100
All	All	2033/2338 (87%)	1952 (96%)	73 (4%)	8 (0%)	30	60

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	406	THR
1	A	965	LYS
1	D	406	THR
1	A	360	LEU
1	A	362	SER
1	A	505	GLU
1	D	506	PRO
1	A	506	PRO

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	837/883 (95%)	837 (100%)	0	100	100
1	D	845/883 (96%)	844 (100%)	1 (0%)	88	90
2	B	69/130 (53%)	69 (100%)	0	100	100
2	E	69/130 (53%)	69 (100%)	0	100	100
3	C	24/37 (65%)	24 (100%)	0	100	100
3	F	24/37 (65%)	24 (100%)	0	100	100
All	All	1868/2100 (89%)	1867 (100%)	1 (0%)	88	90

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	769	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	232	ASN
1	A	247	GLN
1	A	269	GLN
1	A	361	ASN
1	A	389	GLN
1	A	587	ASN
1	A	590	GLN
1	A	593	GLN
1	A	773	GLN
1	A	837	GLN
1	A	880	ASN
1	A	937	ASN
3	C	234	GLN
1	D	141	ASN
1	D	232	ASN
1	D	253	ASN

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Mol	Chain	Res	Type
1	D	415	HIS
1	D	459	ASN
1	D	474	HIS
1	D	536	HIS
1	D	587	ASN
1	D	608	HIS
1	D	615	GLN
1	D	666	ASN
1	D	712	ASN
1	D	735	GLN
1	D	770	GLN
1	D	773	GLN
1	D	880	ASN
3	F	234	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 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	926/985 (94%)	0.39	52 (5%)	30	20	51, 88, 160, 197	0
1	D	941/985 (95%)	0.40	46 (4%)	35	23	47, 96, 151, 220	0
2	B	74/143 (51%)	-0.02	1 (1%)	73	55	53, 77, 101, 131	0
2	E	74/143 (51%)	-0.05	1 (1%)	73	55	54, 79, 106, 124	0
3	C	24/41 (58%)	1.24	3 (12%)	8	6	70, 109, 135, 171	0
3	F	24/41 (58%)	0.91	3 (12%)	8	6	73, 105, 124, 135	0
All	All	2063/2338 (88%)	0.38	106 (5%)	33	22	47, 91, 153, 220	0

All (106) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	362	SER	5.4
1	A	997	LEU	5.0
3	C	241	ASP	4.8
1	D	964	LEU	4.0
1	A	1019	TYR	3.8
1	A	962	ASP	3.6
1	D	507	LEU	3.5
1	A	1045	ARG	3.5
1	D	769	CYS	3.4
1	A	813	LEU	3.4
1	D	749	THR	3.4
3	F	244	HIS	3.4
1	D	992	HIS	3.4
3	C	244	HIS	3.3
1	D	454	GLN	3.3
3	C	231	ASP	3.2
1	A	748	ILE	3.2
1	D	402	GLU	3.1
1	D	677	GLU	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	506	PRO	3.1
1	A	361	ASN	3.0
1	D	361	ASN	3.0
1	A	1011	LEU	2.9
1	D	1031	SER	2.9
1	A	1028	PHE	2.9
3	F	242	ASP	2.9
1	A	359	GLU	2.9
1	A	1016	ARG	2.8
1	A	763	LYS	2.8
1	D	96	GLN	2.8
1	D	400	SER	2.8
1	A	194	GLU	2.8
1	A	402	GLU	2.8
1	A	936	TYR	2.8
1	A	774	HIS	2.7
1	A	964	LEU	2.7
1	A	254	LYS	2.7
1	D	1039	LEU	2.7
1	D	709	PHE	2.7
1	A	246	ARG	2.7
1	A	363	LYS	2.7
1	A	1037	VAL	2.7
1	D	748	ILE	2.6
1	D	935	GLY	2.6
1	D	543	THR	2.6
1	A	761	LEU	2.6
1	A	485	TRP	2.6
1	D	271	ARG	2.6
1	A	994	PRO	2.5
1	D	485	TRP	2.5
1	A	1010	LEU	2.5
1	A	1048	LEU	2.5
1	A	933	GLU	2.5
1	A	1012	ARG	2.5
1	D	91	GLY	2.5
1	A	1030	MET	2.5
1	A	752	THR	2.5
1	D	959	PHE	2.4
1	D	268	LEU	2.4
1	D	1021	TYR	2.4
1	D	1048	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
3	F	241	ASP	2.4
1	D	360	LEU	2.4
1	A	507	LEU	2.4
1	D	1014	ASP	2.4
2	B	394	THR	2.3
1	A	1046	ASN	2.3
1	D	92	LYS	2.3
1	A	89	LYS	2.3
1	D	1028	PHE	2.3
1	A	679	PRO	2.3
1	A	90	MET	2.3
1	A	400	SER	2.3
1	A	406	THR	2.3
1	A	283	ASN	2.3
1	D	243	ASN	2.3
1	A	1027	THR	2.3
1	D	966	THR	2.3
1	D	671	ASP	2.3
1	A	91	GLY	2.2
1	D	260	ALA	2.2
1	A	671	ASP	2.2
1	D	1050	ALA	2.2
1	A	893	TYR	2.2
1	D	90	MET	2.2
1	D	1016	ARG	2.2
1	D	724	ASP	2.2
1	A	1018	VAL	2.1
1	D	256	ILE	2.1
1	A	360	LEU	2.1
1	A	1023	GLU	2.1
1	D	993	PRO	2.1
1	A	261	ASN	2.1
2	E	387	LYS	2.1
1	D	1029	GLN	2.1
1	D	597	LEU	2.1
1	D	720	ILE	2.1
1	A	1031	SER	2.1
1	A	139	MET	2.1
1	D	559	ASP	2.0
1	A	912	LYS	2.0
1	D	1045	ARG	2.0
1	D	1051	GLY	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	511	GLU	2.0
1	D	810	LEU	2.0
1	D	931	ILE	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.