



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 09:38 am BST

PDB ID : 1KN0
Title : Crystal Structure of the human Rad52 protein
Authors : Kagawa, W.; Kurumizaka, H.; Ishitani, R.; Fukai, S.; Nureki, O.; Shibata, T.;
Yokoyama, S.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2001-12-18
Resolution : 2.85 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 : 2.11
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) : 2.11

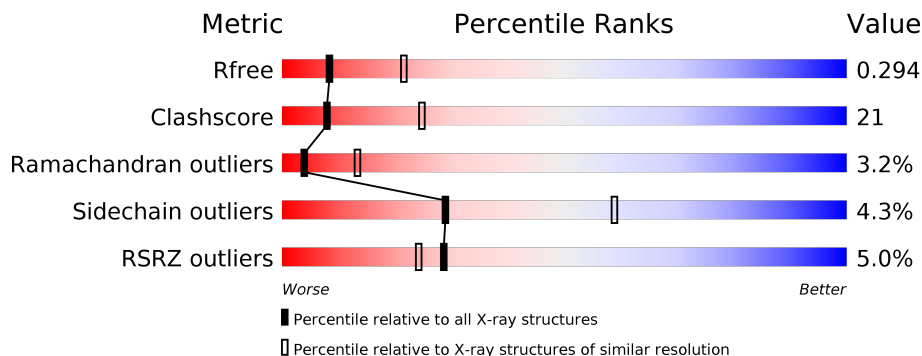
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 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.







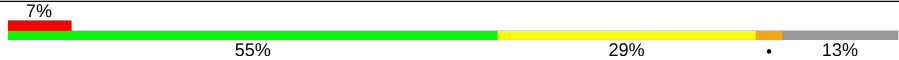
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	3168 (2.90-2.82)
Clashscore	141614	3438 (2.90-2.82)
Ramachandran outliers	138981	3348 (2.90-2.82)
Sidechain outliers	138945	3351 (2.90-2.82)
RSRZ outliers	127900	3103 (2.90-2.82)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 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	212	<div style="display: flex; align-items: center;"> <div style="width: 10px; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 100%; height: 15px; position: relative;"> <div style="width: 49%; height: 100%; background-color: green;"></div> <div style="width: 34%; height: 100%; background-color: yellow;"></div> <div style="width: 13%; height: 100%; background-color: grey;"></div> </div> </div> <p style="margin-left: 20px;">49% 34% 13%</p>
1	B	212	<div style="display: flex; align-items: center;"> <div style="width: 10px; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 100%; height: 15px; position: relative;"> <div style="width: 52%; height: 100%; background-color: green;"></div> <div style="width: 31%; height: 100%; background-color: yellow;"></div> <div style="width: 13%; height: 100%; background-color: grey;"></div> </div> </div> <p style="margin-left: 20px;">52% 31% 13%</p>
1	C	212	<div style="display: flex; align-items: center;"> <div style="width: 10px; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 100%; height: 15px; position: relative;"> <div style="width: 52%; height: 100%; background-color: green;"></div> <div style="width: 31%; height: 100%; background-color: yellow;"></div> <div style="width: 13%; height: 100%; background-color: grey;"></div> </div> </div> <p style="margin-left: 20px;">52% 31% 13%</p>
1	D	212	<div style="display: flex; align-items: center;"> <div style="width: 10px; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 100%; height: 15px; position: relative;"> <div style="width: 4%; height: 100%; background-color: red;"></div> <div style="width: 46%; height: 100%; background-color: green;"></div> <div style="width: 33%; height: 100%; background-color: yellow;"></div> <div style="width: 13%; height: 100%; background-color: grey;"></div> </div> </div> <p style="margin-left: 20px;">4% 50% 33% 13%</p>
1	E	212	<div style="display: flex; align-items: center;"> <div style="width: 10px; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 100%; height: 15px; position: relative;"> <div style="width: 11%; height: 100%; background-color: red;"></div> <div style="width: 37%; height: 100%; background-color: green;"></div> <div style="width: 33%; height: 100%; background-color: yellow;"></div> <div style="width: 6%; height: 100%; background-color: orange;"></div> <div style="width: 13%; height: 100%; background-color: grey;"></div> </div> </div> <p style="margin-left: 20px;">11% 48% 33% 6% 13%</p>
1	F	212	<div style="display: flex; align-items: center;"> <div style="width: 10px; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 100%; height: 15px; position: relative;"> <div style="width: 8%; height: 100%; background-color: red;"></div> <div style="width: 42%; height: 100%; background-color: green;"></div> <div style="width: 32%; height: 100%; background-color: yellow;"></div> <div style="width: 5%; height: 100%; background-color: orange;"></div> <div style="width: 13%; height: 100%; background-color: grey;"></div> </div> </div> <p style="margin-left: 20px;">8% 50% 32% 5% 13%</p>

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Mol	Chain	Length	Quality of chain
1	G	212	
1	H	212	
1	I	212	
1	J	212	
1	K	212	

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 16032 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 Rad52.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	184	1450	907	261	274	8	0	0	0
1	B	184	1450	907	261	274	8	0	0	0
1	C	184	1450	907	261	274	8	0	0	0
1	D	184	1450	907	261	274	8	0	0	0
1	E	184	1450	907	261	274	8	0	0	0
1	F	184	1450	907	261	274	8	0	0	0
1	G	184	1450	907	261	274	8	0	0	0
1	H	184	1450	907	261	274	8	0	0	0
1	I	184	1450	907	261	274	8	0	0	0
1	J	184	1450	907	261	274	8	0	0	0
1	K	184	1450	907	261	274	8	0	0	0

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	9	Total	O	0	0
			9	9		
2	B	12	Total	O	0	0
			12	12		
2	C	9	Total	O	0	0
			9	9		

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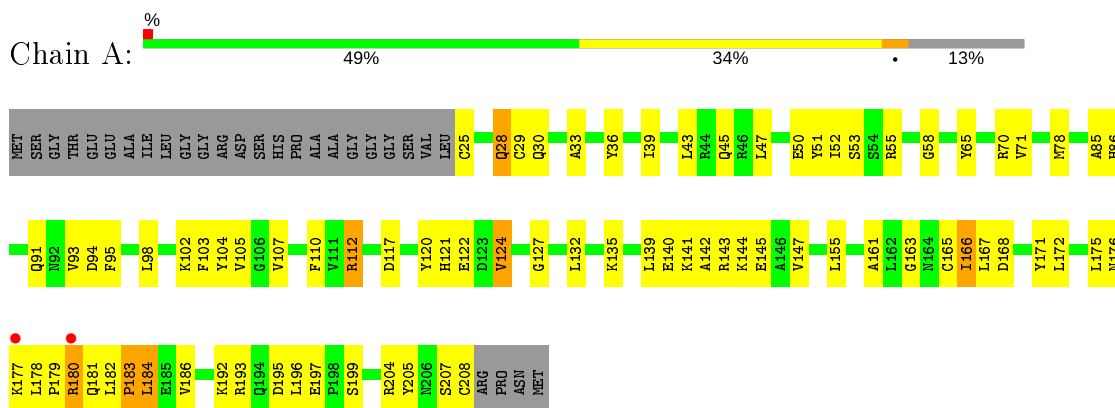
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	D	6	Total O 6 6	0	0
2	E	3	Total O 3 3	0	0
2	F	2	Total O 2 2	0	0
2	G	8	Total O 8 8	0	0
2	H	10	Total O 10 10	0	0
2	I	7	Total O 7 7	0	0
2	J	9	Total O 9 9	0	0
2	K	7	Total O 7 7	0	0

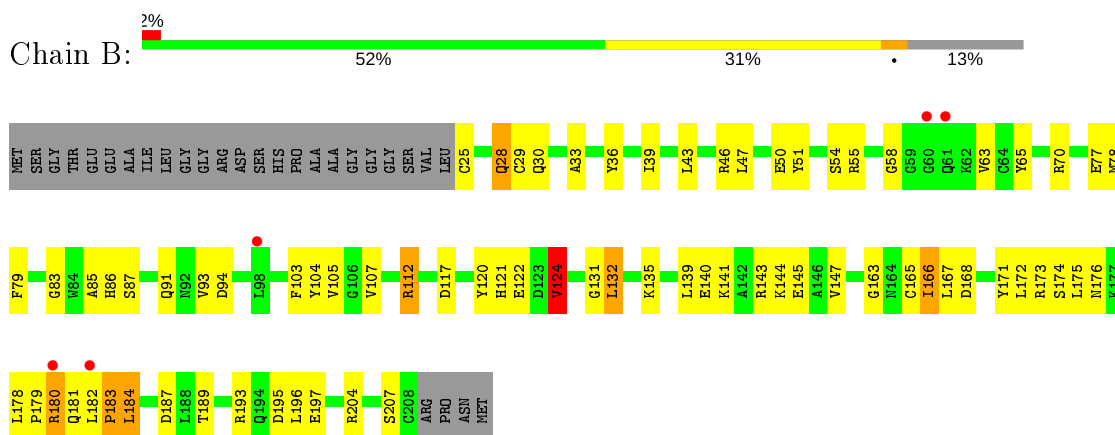
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

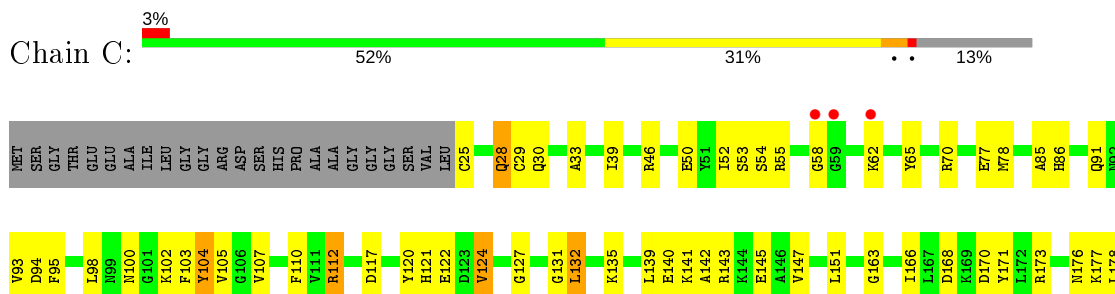
- Molecule 1: Rad52



- Molecule 1: Rad52

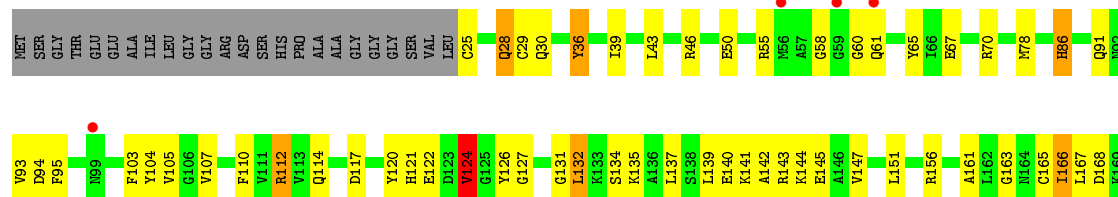


- Molecule 1: Rad52

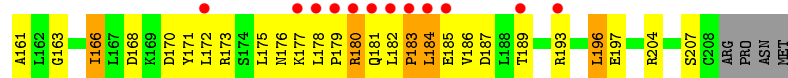
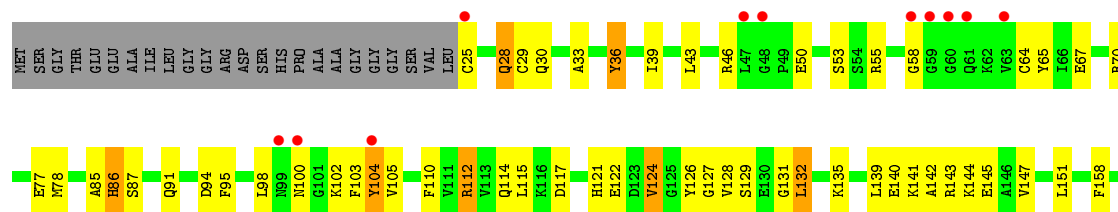




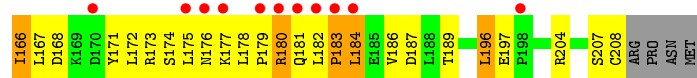
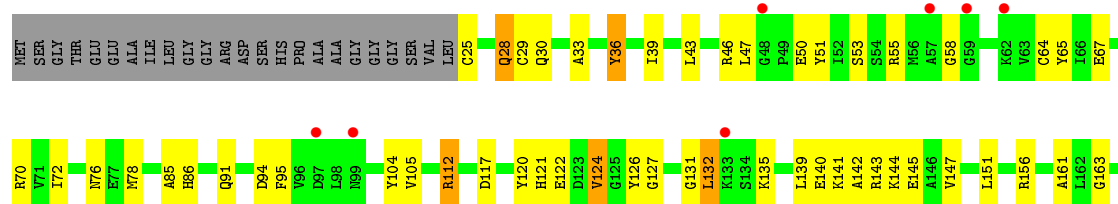
- Molecule 1: Rad52



- Molecule 1: Rad52

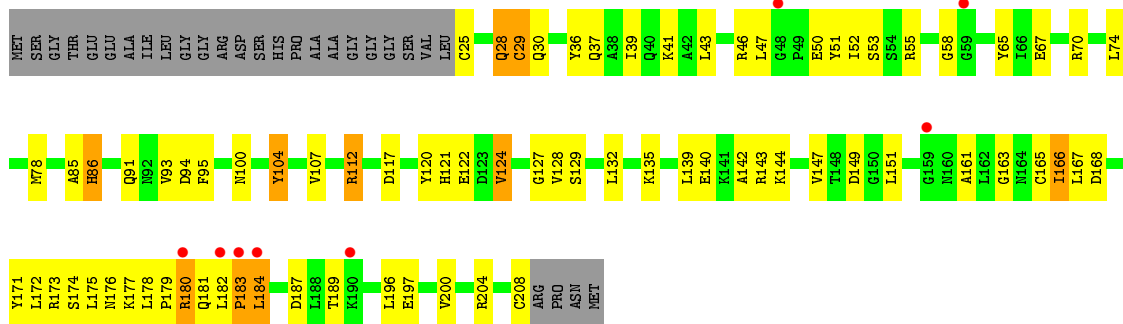


- Molecule 1: Rad52

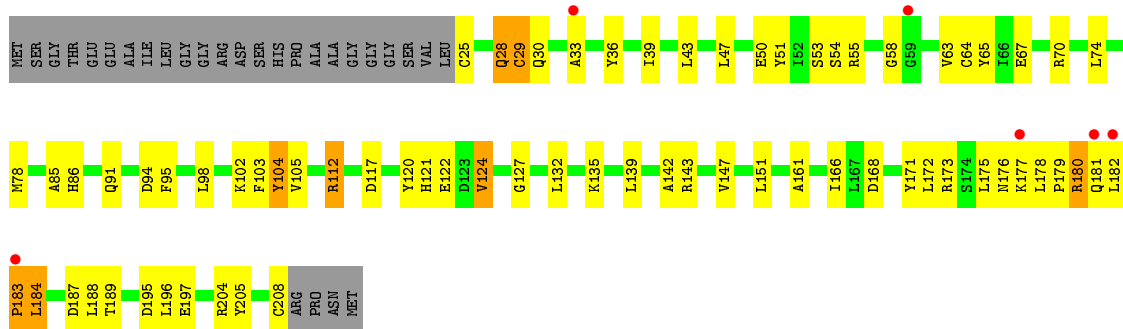


- Molecule 1: Rad52

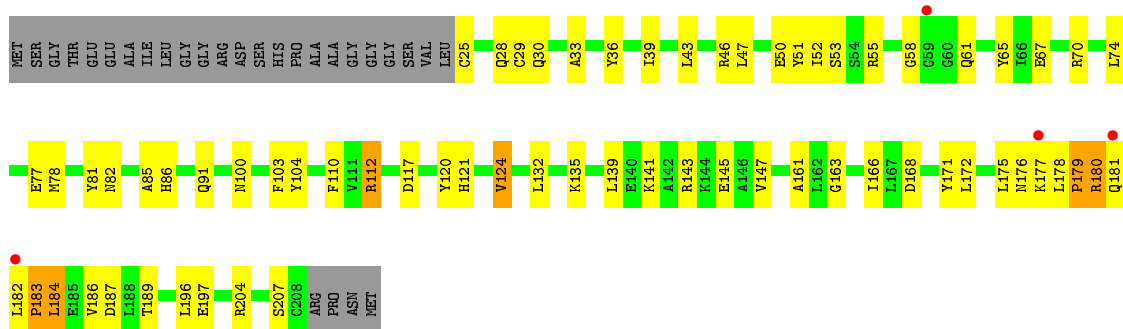




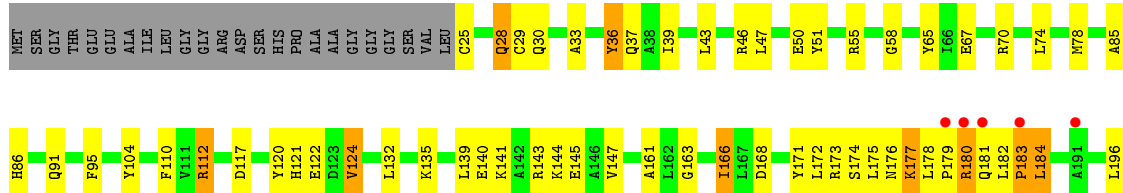
• Molecule 1: Rad52

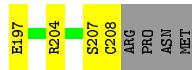


• Molecule 1: Rad52

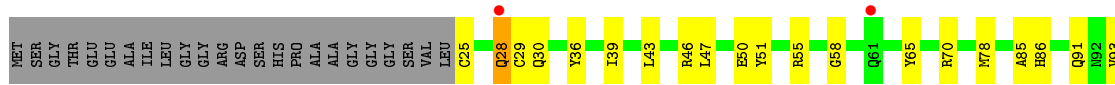


• Molecule 1: Rad52





- Molecule 1: Rad52



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	145.61Å 145.61Å 247.53Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.85 45.27 – 2.85	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-2.85) 99.3 (45.27-2.85)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	13.06 (at 2.86Å)	Xtrriage
Refinement program	CNS	Depositor
R, R_{free}	0.231 , 0.297 0.271 , 0.294	Depositor DCC
R_{free} test set	6303 reflections (10.11%)	wwPDB-VP
Wilson B-factor (Å ²)	32.5	Xtrriage
Anisotropy	0.084	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 76.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.28$, $\langle L^2 \rangle = 0.12$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.84	EDS
Total number of atoms	16032	wwPDB-VP
Average B, all atoms (Å ²)	81.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.31% 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	0.45	0/1474	0.63	0/1981
1	B	0.46	0/1474	0.63	1/1981 (0.1%)
1	C	0.42	0/1474	0.61	0/1981
1	D	0.42	0/1474	0.61	1/1981 (0.1%)
1	E	0.40	0/1474	0.60	0/1981
1	F	0.40	0/1474	0.59	0/1981
1	G	0.42	0/1474	0.62	0/1981
1	H	0.41	0/1474	0.61	0/1981
1	I	0.41	0/1474	0.62	0/1981
1	J	0.41	0/1474	0.62	0/1981
1	K	0.43	0/1474	0.61	0/1981
All	All	0.42	0/16214	0.61	2/21791 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	D	0	1
1	E	0	1
1	F	0	1
1	J	0	1
All	All	0	4

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	124	VAL	CB-CA-C	-5.46	101.03	111.40
1	D	124	VAL	CB-CA-C	-5.03	101.85	111.40

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	36	TYR	Sidechain
1	E	36	TYR	Sidechain
1	F	36	TYR	Sidechain
1	J	36	TYR	Sidechain

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1450	0	1433	74	0
1	B	1450	0	1433	74	0
1	C	1450	0	1433	76	0
1	D	1450	0	1433	83	1
1	E	1450	0	1433	87	1
1	F	1450	0	1433	81	0
1	G	1450	0	1433	87	1
1	H	1450	0	1433	71	0
1	I	1450	0	1433	72	0
1	J	1450	0	1433	77	0
1	K	1450	0	1433	69	0
2	A	9	0	0	2	0
2	B	12	0	0	0	0
2	C	9	0	0	0	0
2	D	6	0	0	0	0
2	E	3	0	0	0	0
2	F	2	0	0	0	0
2	G	8	0	0	0	0
2	H	10	0	0	0	0
2	I	7	0	0	0	0
2	J	9	0	0	0	0
2	K	7	0	0	0	0
All	All	16032	0	15763	681	2

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 21.

All (681) 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:G:182:LEU:HD12	1:G:183:PRO:HD2	1.43	0.99
1:A:86:HIS:HD2	1:B:121:HIS:HD2	1.09	0.98
1:G:182:LEU:CD1	1:G:183:PRO:HD2	1.96	0.95
1:A:121:HIS:HD2	1:K:86:HIS:HD2	1.14	0.93
1:B:86:HIS:HD2	1:C:121:HIS:HD2	1.14	0.93
1:J:86:HIS:HD2	1:K:121:HIS:HD2	1.15	0.89
1:B:147:VAL:HG21	1:C:124:VAL:CG2	2.04	0.88
1:G:182:LEU:HD12	1:G:183:PRO:CD	2.04	0.87
1:H:86:HIS:HD2	1:I:121:HIS:HD2	1.23	0.85
1:K:91:GLN:NE2	1:K:143:ARG:HE	1.75	0.85
1:I:86:HIS:HD2	1:J:121:HIS:HD2	1.23	0.85
1:J:147:VAL:HG21	1:K:124:VAL:CG2	2.09	0.82
1:F:86:HIS:CD2	1:G:121:HIS:HD2	1.97	0.82
1:A:39:ILE:HG23	1:A:78:MET:CE	2.10	0.82
1:A:86:HIS:CD2	1:B:121:HIS:HD2	1.96	0.82
1:E:39:ILE:HG23	1:E:78:MET:CE	2.09	0.81
1:F:86:HIS:HD2	1:G:121:HIS:HD2	1.24	0.81
1:A:86:HIS:HD2	1:B:121:HIS:CD2	1.97	0.80
1:B:86:HIS:HD2	1:C:121:HIS:CD2	2.00	0.79
1:G:86:HIS:CD2	1:H:121:HIS:HD2	2.00	0.79
1:C:39:ILE:HG23	1:C:78:MET:CE	2.13	0.78
1:E:141:LYS:HE2	1:E:145:GLU:OE2	1.83	0.78
1:D:147:VAL:HG21	1:E:124:VAL:CG2	2.13	0.78
1:D:91:GLN:NE2	1:D:143:ARG:HE	1.80	0.78
1:H:39:ILE:HG23	1:H:78:MET:CE	2.14	0.78
1:J:86:HIS:HD2	1:K:121:HIS:CD2	2.03	0.77
1:I:39:ILE:HG23	1:I:78:MET:CE	2.15	0.77
1:G:50:GLU:HG3	1:G:181:GLN:HE22	1.49	0.77
1:H:91:GLN:NE2	1:H:143:ARG:HE	1.83	0.76
1:K:39:ILE:HG23	1:K:78:MET:CE	2.15	0.76
1:F:39:ILE:HG23	1:F:78:MET:CE	2.16	0.75
1:C:86:HIS:HD2	1:D:121:HIS:HD2	1.33	0.75
1:A:121:HIS:CD2	1:K:86:HIS:HD2	2.02	0.75
1:I:147:VAL:HG21	1:J:124:VAL:CG2	2.17	0.75
1:F:50:GLU:HG3	1:F:181:GLN:HE22	1.52	0.74
1:D:86:HIS:CD2	1:E:121:HIS:HD2	2.05	0.74
1:F:86:HIS:HD2	1:G:121:HIS:CD2	2.05	0.74
1:C:86:HIS:CD2	1:D:121:HIS:HD2	2.05	0.74
1:I:182:LEU:HG	1:I:183:PRO:HD2	1.69	0.73
1:J:50:GLU:HG3	1:J:181:GLN:HE22	1.51	0.73
1:H:147:VAL:HG21	1:I:124:VAL:CG2	2.18	0.73
1:I:50:GLU:HG3	1:I:181:GLN:HE22	1.54	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:86:HIS:CD2	1:K:121:HIS:HD2	2.04	0.73
1:C:91:GLN:NE2	1:C:143:ARG:HE	1.87	0.73
1:D:39:ILE:HG23	1:D:78:MET:CE	2.19	0.73
1:G:147:VAL:HG21	1:H:124:VAL:CG2	2.17	0.73
1:H:86:HIS:CD2	1:I:121:HIS:HD2	2.07	0.72
1:H:28:GLN:HE21	1:H:28:GLN:HA	1.54	0.72
1:K:50:GLU:HG3	1:K:181:GLN:HE22	1.54	0.72
1:G:39:ILE:HG23	1:G:78:MET:CE	2.19	0.72
1:H:50:GLU:HG3	1:H:181:GLN:HE22	1.54	0.72
1:J:39:ILE:HG23	1:J:78:MET:CE	2.20	0.72
1:J:50:GLU:HG3	1:J:181:GLN:NE2	2.05	0.72
1:B:50:GLU:HG3	1:B:181:GLN:HE22	1.54	0.71
1:I:85:ALA:HB1	1:J:120:TYR:O	1.91	0.71
1:B:182:LEU:HG	1:B:183:PRO:HD2	1.72	0.70
1:I:91:GLN:NE2	1:I:143:ARG:HE	1.88	0.70
1:A:124:VAL:CG2	1:K:147:VAL:HG21	2.22	0.70
1:K:28:GLN:HE21	1:K:28:GLN:HA	1.54	0.70
1:E:50:GLU:HG3	1:E:181:GLN:HE22	1.55	0.70
1:F:204:ARG:HD2	1:H:36:TYR:OH	1.91	0.70
1:G:50:GLU:HG3	1:G:181:GLN:NE2	2.07	0.70
1:I:86:HIS:HD2	1:J:121:HIS:CD2	2.08	0.70
1:F:50:GLU:HG3	1:F:181:GLN:NE2	2.06	0.70
1:B:86:HIS:CD2	1:C:121:HIS:HD2	2.04	0.70
1:G:184:LEU:HD22	1:G:184:LEU:H	1.57	0.70
1:H:184:LEU:H	1:H:184:LEU:HD22	1.56	0.70
1:H:50:GLU:HG3	1:H:181:GLN:NE2	2.06	0.70
1:H:112:ARG:HG3	1:H:122:GLU:HB2	1.74	0.69
1:A:50:GLU:HG3	1:A:181:GLN:HE22	1.58	0.69
1:A:39:ILE:HG23	1:A:78:MET:HE1	1.75	0.69
1:J:28:GLN:HA	1:J:28:GLN:HE21	1.58	0.69
1:B:39:ILE:HG23	1:B:78:MET:CE	2.22	0.69
1:F:147:VAL:HG21	1:G:124:VAL:CG2	2.23	0.69
1:G:86:HIS:HD2	1:H:121:HIS:HD2	1.39	0.69
1:I:50:GLU:HG3	1:I:181:GLN:NE2	2.08	0.69
1:J:91:GLN:NE2	1:J:143:ARG:HE	1.91	0.69
1:E:28:GLN:HA	1:E:28:GLN:HE21	1.58	0.69
1:A:147:VAL:HG21	1:B:124:VAL:CG2	2.23	0.68
1:G:28:GLN:HA	1:G:28:GLN:HE21	1.58	0.68
1:I:184:LEU:H	1:I:184:LEU:HD22	1.58	0.68
1:J:184:LEU:H	1:J:184:LEU:HD22	1.58	0.68
1:C:86:HIS:HD2	1:D:121:HIS:CD2	2.11	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:70:ARG:HH11	1:D:70:ARG:HG3	1.57	0.68
1:E:50:GLU:HG3	1:E:181:GLN:NE2	2.08	0.68
1:F:184:LEU:HD22	1:F:184:LEU:H	1.57	0.68
1:F:25:CYS:HA	1:G:30:GLN:HG2	1.76	0.68
1:K:50:GLU:HG3	1:K:181:GLN:NE2	2.08	0.68
1:A:28:GLN:HE21	1:A:28:GLN:HA	1.59	0.67
1:B:50:GLU:HG3	1:B:181:GLN:NE2	2.09	0.67
1:D:25:CYS:HA	1:E:30:GLN:HG2	1.76	0.67
1:H:182:LEU:HG	1:H:183:PRO:HD2	1.76	0.67
1:A:121:HIS:HD2	1:K:86:HIS:CD2	2.05	0.67
1:K:184:LEU:H	1:K:184:LEU:HD22	1.60	0.67
1:A:184:LEU:H	1:A:184:LEU:HD22	1.59	0.67
1:G:86:HIS:CD2	1:H:121:HIS:CD2	2.83	0.67
1:C:147:VAL:HG21	1:D:124:VAL:CG2	2.24	0.66
1:E:204:ARG:HD2	1:G:36:TYR:OH	1.95	0.66
1:A:50:GLU:HG3	1:A:181:GLN:NE2	2.10	0.66
1:E:147:VAL:HG21	1:F:124:VAL:CG2	2.25	0.66
1:F:28:GLN:HE21	1:F:28:GLN:HA	1.58	0.66
1:E:25:CYS:HA	1:F:30:GLN:HG2	1.75	0.66
1:G:70:ARG:HH11	1:G:70:ARG:HG3	1.60	0.66
1:G:86:HIS:HD2	1:H:121:HIS:CD2	2.14	0.66
1:A:78:MET:O	1:K:204:ARG:NH2	2.29	0.66
1:E:112:ARG:HG3	1:E:122:GLU:HB2	1.78	0.66
1:H:208:CYS:HB2	1:J:33:ALA:HB2	1.78	0.66
1:C:184:LEU:H	1:C:184:LEU:HD22	1.60	0.66
1:B:28:GLN:HA	1:B:28:GLN:HE21	1.60	0.66
1:I:70:ARG:HG3	1:I:70:ARG:HH11	1.61	0.66
1:A:91:GLN:NE2	1:A:143:ARG:HE	1.93	0.66
2:A:213:HOH:O	1:B:121:HIS:HE1	1.79	0.65
1:D:184:LEU:HD22	1:D:184:LEU:H	1.61	0.65
1:F:182:LEU:HG	1:F:183:PRO:HD2	1.78	0.65
1:A:204:ARG:NH2	1:B:78:MET:O	2.29	0.65
1:D:50:GLU:HG3	1:D:181:GLN:HE22	1.59	0.65
1:F:91:GLN:NE2	1:F:143:ARG:HE	1.94	0.65
1:D:86:HIS:CD2	1:E:121:HIS:CD2	2.85	0.65
1:A:93:VAL:HA	1:A:107:VAL:HG22	1.79	0.65
1:B:91:GLN:NE2	1:B:143:ARG:HE	1.94	0.65
1:C:50:GLU:HG3	1:C:181:GLN:HE22	1.61	0.65
1:K:196:LEU:HD12	1:K:197:GLU:H	1.61	0.65
1:E:196:LEU:HD12	1:E:197:GLU:H	1.62	0.64
1:D:141:LYS:HE2	1:D:145:GLU:OE2	1.98	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:86:HIS:CD2	1:F:121:HIS:HD2	2.15	0.64
1:I:28:GLN:HE21	1:I:28:GLN:HA	1.63	0.64
1:B:70:ARG:HH11	1:B:70:ARG:HG3	1.63	0.64
1:E:91:GLN:NE2	1:E:143:ARG:HE	1.96	0.64
1:D:182:LEU:HG	1:D:183:PRO:HD2	1.80	0.64
1:B:184:LEU:H	1:B:184:LEU:HD22	1.63	0.63
1:K:182:LEU:HG	1:K:183:PRO:HD2	1.80	0.63
1:D:28:GLN:HA	1:D:28:GLN:HE21	1.64	0.63
1:E:184:LEU:H	1:E:184:LEU:HD22	1.63	0.63
1:J:182:LEU:HG	1:J:183:PRO:HD2	1.81	0.63
1:B:141:LYS:HE2	1:B:145:GLU:OE2	1.99	0.63
1:D:50:GLU:HG3	1:D:181:GLN:NE2	2.14	0.63
1:C:50:GLU:HG3	1:C:181:GLN:NE2	2.14	0.63
1:E:182:LEU:HG	1:E:183:PRO:HD2	1.81	0.63
1:G:85:ALA:HB1	1:H:120:TYR:O	1.97	0.62
1:H:85:ALA:HB1	1:I:120:TYR:O	1.99	0.62
1:D:147:VAL:HG21	1:E:124:VAL:HG22	1.81	0.62
1:H:70:ARG:HH11	1:H:70:ARG:HG3	1.63	0.62
1:J:70:ARG:HG3	1:J:70:ARG:HH11	1.64	0.62
1:B:204:ARG:HD2	1:D:36:TYR:OH	1.98	0.62
1:B:135:LYS:HG2	1:C:95:PHE:CE2	2.35	0.62
1:C:29:CYS:O	1:C:117:ASP:HA	2.00	0.62
1:I:29:CYS:O	1:I:117:ASP:HA	1.99	0.62
1:D:135:LYS:HG2	1:E:95:PHE:CE2	2.34	0.61
1:F:208:CYS:HB2	1:H:33:ALA:HB2	1.80	0.61
1:H:86:HIS:HD2	1:I:121:HIS:CD2	2.10	0.61
1:J:29:CYS:O	1:J:117:ASP:HA	2.00	0.61
1:I:86:HIS:CD2	1:J:121:HIS:HD2	2.11	0.61
1:A:124:VAL:HG22	1:K:147:VAL:HG21	1.80	0.61
1:F:86:HIS:CD2	1:G:121:HIS:CD2	2.83	0.61
1:G:25:CYS:HA	1:H:30:GLN:HG2	1.82	0.61
1:J:86:HIS:CD2	1:K:121:HIS:CD2	2.86	0.61
1:E:29:CYS:O	1:E:117:ASP:HA	2.00	0.61
1:G:29:CYS:O	1:G:117:ASP:HA	2.00	0.61
1:J:112:ARG:HG3	1:J:122:GLU:HB2	1.82	0.61
1:A:192:LYS:NZ	1:A:197:GLU:OE2	2.30	0.60
1:F:196:LEU:HD12	1:F:197:GLU:H	1.65	0.60
1:G:55:ARG:HD3	1:G:65:TYR:CE2	2.37	0.60
1:C:105:VAL:HG11	1:C:139:LEU:HD23	1.83	0.60
1:C:182:LEU:HG	1:C:183:PRO:HD2	1.82	0.60
1:J:25:CYS:HA	1:K:30:GLN:HG2	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:86:HIS:CD2	1:B:121:HIS:CD2	2.81	0.60
1:B:29:CYS:O	1:B:117:ASP:HA	2.01	0.60
1:G:112:ARG:HG3	1:G:122:GLU:HB2	1.82	0.60
1:A:30:GLN:HG2	1:K:25:CYS:HA	1.84	0.60
1:A:196:LEU:HD12	1:A:197:GLU:H	1.66	0.60
1:H:39:ILE:HG23	1:H:78:MET:HE3	1.84	0.60
1:B:36:TYR:OH	1:K:204:ARG:HD2	2.02	0.60
1:E:39:ILE:HG23	1:E:78:MET:HE3	1.84	0.59
1:G:43:LEU:HG	1:G:78:MET:CE	2.31	0.59
1:F:147:VAL:HG21	1:G:124:VAL:HG22	1.84	0.59
1:E:55:ARG:HD3	1:E:65:TYR:CE2	2.37	0.59
1:E:110:PHE:CE2	1:E:124:VAL:HG13	2.37	0.59
1:I:39:ILE:HG23	1:I:78:MET:HE1	1.84	0.59
1:A:29:CYS:O	1:A:117:ASP:HA	2.03	0.59
1:G:208:CYS:HB2	1:I:33:ALA:HB2	1.85	0.59
1:F:112:ARG:HG3	1:F:122:GLU:HB2	1.83	0.59
1:F:204:ARG:NH2	1:G:78:MET:O	2.36	0.59
1:G:204:ARG:HD2	1:I:36:TYR:OH	2.02	0.59
1:A:182:LEU:HG	1:A:183:PRO:HD2	1.83	0.59
1:F:135:LYS:HG2	1:G:95:PHE:CE2	2.38	0.59
1:B:147:VAL:HG21	1:C:124:VAL:HG22	1.86	0.58
1:J:147:VAL:HG21	1:K:124:VAL:HG22	1.83	0.58
1:A:47:LEU:HD12	1:A:51:TYR:CD1	2.38	0.58
1:G:182:LEU:CG	1:G:183:PRO:HD2	2.33	0.58
1:I:196:LEU:HD12	1:I:197:GLU:H	1.67	0.58
1:J:196:LEU:HD12	1:J:197:GLU:H	1.68	0.58
1:B:168:ASP:HB3	1:B:171:TYR:HB3	1.86	0.58
1:H:184:LEU:N	1:H:184:LEU:HD22	2.18	0.58
1:I:25:CYS:HA	1:J:30:GLN:HG2	1.84	0.58
1:D:55:ARG:HD3	1:D:65:TYR:CE2	2.39	0.58
1:K:70:ARG:HG3	1:K:70:ARG:HH11	1.67	0.58
1:E:70:ARG:HG3	1:E:70:ARG:HH11	1.67	0.57
1:F:39:ILE:HG23	1:F:78:MET:HE3	1.85	0.57
1:G:91:GLN:NE2	1:G:143:ARG:HE	2.01	0.57
1:H:29:CYS:O	1:H:117:ASP:HA	2.04	0.57
1:A:70:ARG:HH11	1:A:70:ARG:HG3	1.68	0.57
1:F:55:ARG:HD3	1:F:65:TYR:CE2	2.39	0.57
1:C:86:HIS:CD2	1:D:121:HIS:CD2	2.89	0.57
1:D:50:GLU:N	1:D:50:GLU:OE1	2.36	0.57
1:F:70:ARG:HG3	1:F:70:ARG:HH11	1.68	0.57
1:J:55:ARG:HD3	1:J:65:TYR:CE2	2.39	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:39:ILE:HG23	1:C:78:MET:HE1	1.87	0.57
1:D:112:ARG:HG3	1:D:122:GLU:HB2	1.87	0.57
1:C:208:CYS:HB2	1:E:33:ALA:HB2	1.87	0.57
1:F:29:CYS:O	1:F:117:ASP:HA	2.05	0.57
1:G:39:ILE:HG23	1:G:78:MET:HE3	1.87	0.57
1:C:28:GLN:HA	1:C:28:GLN:HE21	1.70	0.56
1:H:103:PHE:CE1	1:H:135:LYS:HB2	2.40	0.56
1:H:168:ASP:HB3	1:H:171:TYR:HB3	1.87	0.56
1:A:121:HIS:CD2	1:K:86:HIS:CD2	2.87	0.56
1:D:204:ARG:HD2	1:F:36:TYR:OH	2.04	0.56
1:H:196:LEU:HD12	1:H:197:GLU:H	1.69	0.56
1:B:86:HIS:CD2	1:C:121:HIS:CD2	2.85	0.56
1:C:184:LEU:N	1:C:184:LEU:HD22	2.21	0.56
1:D:196:LEU:HD12	1:D:197:GLU:H	1.71	0.56
1:H:43:LEU:HG	1:H:78:MET:HE1	1.88	0.56
1:C:93:VAL:HA	1:C:107:VAL:HG22	1.88	0.56
1:G:196:LEU:HD12	1:G:197:GLU:H	1.71	0.56
1:G:37:GLN:HG2	1:G:41:LYS:HE2	1.87	0.56
1:I:182:LEU:CG	1:I:183:PRO:HD2	2.36	0.56
1:A:147:VAL:HG21	1:B:124:VAL:HG22	1.86	0.56
1:C:147:VAL:HG21	1:D:124:VAL:HG22	1.88	0.56
1:E:85:ALA:HB1	1:F:120:TYR:O	2.06	0.56
1:I:184:LEU:N	1:I:184:LEU:HD22	2.21	0.56
1:A:55:ARG:HD3	1:A:65:TYR:CE2	2.41	0.55
1:B:196:LEU:HD12	1:B:197:GLU:H	1.69	0.55
1:H:43:LEU:HG	1:H:78:MET:CE	2.36	0.55
1:F:184:LEU:HD22	1:F:184:LEU:N	2.20	0.55
1:A:85:ALA:HB1	1:B:120:TYR:O	2.05	0.55
1:H:205:TYR:CE1	1:J:37:GLN:NE2	2.75	0.55
1:I:55:ARG:HD3	1:I:65:TYR:CE2	2.42	0.55
1:B:182:LEU:CG	1:B:183:PRO:HD2	2.37	0.55
1:A:143:ARG:NH1	1:B:94:ASP:OD1	2.40	0.55
1:G:135:LYS:HG2	1:H:95:PHE:CE2	2.42	0.55
1:J:184:LEU:N	1:J:184:LEU:HD22	2.21	0.55
1:A:112:ARG:HG3	1:A:122:GLU:HB2	1.88	0.55
1:H:25:CYS:HA	1:I:30:GLN:HG2	1.88	0.55
1:H:86:HIS:CD2	1:I:121:HIS:CD2	2.91	0.54
1:K:140:GLU:HG3	1:K:144:LYS:HE3	1.88	0.54
1:B:33:ALA:HB2	1:K:208:CYS:HB2	1.87	0.54
1:F:141:LYS:HE2	1:F:145:GLU:OE2	2.07	0.54
1:B:87:SER:HA	1:C:122:GLU:HB3	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:184:LEU:N	1:A:184:LEU:HD22	2.21	0.54
1:E:39:ILE:HG23	1:E:78:MET:HE1	1.89	0.54
1:J:143:ARG:NH1	1:K:94:ASP:OD1	2.41	0.54
1:K:112:ARG:HG3	1:K:122:GLU:HB2	1.90	0.54
1:E:86:HIS:HD2	1:F:121:HIS:HD2	1.56	0.54
1:E:140:GLU:HB2	1:F:126:TYR:CD2	2.42	0.54
1:F:187:ASP:OD1	1:F:189:THR:HG23	2.08	0.54
1:F:85:ALA:HB1	1:G:120:TYR:O	2.08	0.54
1:E:135:LYS:HG2	1:F:95:PHE:CE2	2.43	0.53
1:E:86:HIS:CD2	1:F:121:HIS:CD2	2.97	0.53
1:F:168:ASP:HB3	1:F:171:TYR:HB3	1.90	0.53
1:J:139:LEU:O	1:J:143:ARG:HG3	2.08	0.53
1:C:182:LEU:CG	1:C:183:PRO:HD2	2.39	0.53
1:D:184:LEU:HD22	1:D:184:LEU:N	2.23	0.53
1:E:86:HIS:HD2	1:F:121:HIS:CD2	2.27	0.53
1:G:184:LEU:HD22	1:G:184:LEU:N	2.21	0.53
1:C:182:LEU:CD1	1:C:183:PRO:HD2	2.38	0.53
1:C:85:ALA:HB1	1:D:120:TYR:O	2.08	0.53
1:H:28:GLN:HA	1:H:28:GLN:NE2	2.24	0.53
1:C:187:ASP:OD1	1:C:189:THR:HG23	2.08	0.53
1:A:91:GLN:HE22	1:A:143:ARG:HB3	1.75	0.52
1:A:91:GLN:HE21	1:A:143:ARG:HE	1.58	0.52
1:D:29:CYS:O	1:D:117:ASP:HA	2.09	0.52
1:D:86:HIS:HD2	1:E:121:HIS:CD2	2.25	0.52
1:F:67:GLU:O	1:F:70:ARG:HB2	2.10	0.52
1:B:184:LEU:N	1:B:184:LEU:HD22	2.23	0.52
1:G:143:ARG:NH1	1:H:94:ASP:OD1	2.42	0.52
1:K:141:LYS:HE2	1:K:145:GLU:OE2	2.10	0.52
1:D:103:PHE:CE1	1:D:135:LYS:HB2	2.44	0.52
1:D:86:HIS:CE1	1:D:151:LEU:HD22	2.44	0.52
1:G:43:LEU:HG	1:G:78:MET:HE1	1.91	0.52
1:I:86:HIS:CD2	1:J:121:HIS:CD2	2.92	0.52
1:C:112:ARG:HG3	1:C:122:GLU:HB2	1.90	0.52
1:G:168:ASP:HB3	1:G:171:TYR:HB3	1.92	0.52
1:D:179:PRO:O	1:D:181:GLN:N	2.43	0.52
1:I:43:LEU:HG	1:I:78:MET:CE	2.39	0.52
1:A:94:ASP:OD1	1:K:143:ARG:NH1	2.43	0.52
1:C:70:ARG:HH11	1:C:70:ARG:HG3	1.75	0.52
1:E:67:GLU:O	1:E:70:ARG:HB2	2.10	0.52
1:A:172:LEU:O	1:A:175:LEU:HB2	2.10	0.51
1:A:208:CYS:HB2	1:C:33:ALA:HB2	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:93:VAL:HA	1:B:107:VAL:HG22	1.92	0.51
1:D:105:VAL:HG11	1:D:139:LEU:HD23	1.92	0.51
1:E:43:LEU:HG	1:E:78:MET:HE1	1.92	0.51
1:F:135:LYS:HE3	1:G:95:PHE:CG	2.45	0.51
1:K:55:ARG:HD3	1:K:65:TYR:CE2	2.46	0.51
1:B:187:ASP:OD1	1:B:189:THR:HG23	2.10	0.51
1:G:50:GLU:CG	1:G:181:GLN:HE22	2.22	0.51
1:J:135:LYS:HG2	1:K:95:PHE:CE2	2.44	0.51
1:J:140:GLU:HG3	1:J:144:LYS:HE3	1.91	0.51
1:B:85:ALA:HB1	1:C:120:TYR:O	2.10	0.51
1:C:204:ARG:O	1:C:207:SER:HB3	2.10	0.51
1:D:39:ILE:HG23	1:D:78:MET:HE1	1.90	0.51
1:E:103:PHE:CE1	1:E:135:LYS:HB2	2.45	0.51
1:H:54:SER:HA	1:H:63:VAL:O	2.10	0.51
1:D:176:ASN:C	1:D:178:LEU:H	2.14	0.51
1:G:187:ASP:OD1	1:G:189:THR:HG23	2.10	0.51
1:E:204:ARG:CD	1:G:36:TYR:OH	2.58	0.51
1:I:172:LEU:O	1:I:175:LEU:HB2	2.11	0.51
1:B:105:VAL:HG11	1:B:139:LEU:HD23	1.92	0.51
1:E:182:LEU:CG	1:E:183:PRO:HD2	2.41	0.51
1:E:184:LEU:N	1:E:184:LEU:HD22	2.25	0.51
1:E:204:ARG:NH2	1:F:78:MET:O	2.43	0.51
1:B:143:ARG:NH1	1:C:94:ASP:OD1	2.43	0.51
1:C:55:ARG:HD3	1:C:65:TYR:CE2	2.46	0.50
1:D:140:GLU:HG3	1:D:144:LYS:HE3	1.93	0.50
1:F:46:ARG:HD3	1:F:171:TYR:CD1	2.47	0.50
1:G:52:ILE:HG22	1:G:53:SER:N	2.27	0.50
1:K:93:VAL:HA	1:K:107:VAL:HG22	1.93	0.50
1:H:182:LEU:CG	1:H:183:PRO:HD2	2.39	0.50
1:K:29:CYS:O	1:K:117:ASP:HA	2.11	0.50
1:A:36:TYR:OH	1:J:204:ARG:HD2	2.12	0.50
1:C:141:LYS:HE2	1:C:145:GLU:OE2	2.10	0.50
1:D:182:LEU:CG	1:D:183:PRO:HD2	2.41	0.50
1:D:43:LEU:HA	1:D:161:ALA:HB3	1.94	0.50
1:E:43:LEU:HG	1:E:78:MET:CE	2.42	0.50
1:H:55:ARG:HD3	1:H:65:TYR:CE2	2.47	0.50
1:I:176:ASN:C	1:I:178:LEU:H	2.15	0.50
1:E:135:LYS:HE3	1:F:95:PHE:CG	2.46	0.50
1:H:47:LEU:HD12	1:H:51:TYR:CD1	2.47	0.50
1:K:28:GLN:NE2	1:K:28:GLN:HA	2.24	0.50
1:A:163:GLY:O	1:A:166:ILE:HG22	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:163:GLY:O	1:D:166:ILE:HG22	2.12	0.50
1:E:179:PRO:O	1:E:181:GLN:N	2.45	0.50
1:I:135:LYS:HG2	1:J:95:PHE:CE2	2.46	0.50
1:K:163:GLY:O	1:K:166:ILE:HG22	2.12	0.50
1:J:204:ARG:NH2	1:K:78:MET:O	2.45	0.50
1:G:47:LEU:HD12	1:G:51:TYR:CD1	2.46	0.49
1:K:104:TYR:CD1	1:K:104:TYR:N	2.80	0.49
1:B:140:GLU:O	1:B:144:LYS:HG3	2.12	0.49
1:D:135:LYS:HE3	1:E:95:PHE:CD2	2.47	0.49
1:D:187:ASP:OD1	1:D:189:THR:HG23	2.12	0.49
1:G:173:ARG:HB3	1:G:173:ARG:HH11	1.77	0.49
1:H:53:SER:O	1:H:64:CYS:HA	2.12	0.49
1:K:182:LEU:CG	1:K:183:PRO:HD2	2.42	0.49
1:D:156:ARG:NH1	1:D:167:LEU:HD21	2.27	0.49
1:I:47:LEU:HD12	1:I:51:TYR:CD1	2.47	0.49
1:J:141:LYS:HE2	1:J:145:GLU:OE2	2.13	0.49
1:J:176:ASN:C	1:J:178:LEU:H	2.15	0.49
1:B:195:ASP:HB3	1:C:77:GLU:HG3	1.94	0.49
1:G:176:ASN:C	1:G:178:LEU:H	2.15	0.49
1:D:110:PHE:CE2	1:D:124:VAL:HG13	2.47	0.49
1:G:86:HIS:HA	1:G:112:ARG:O	2.12	0.49
1:E:143:ARG:NH1	1:F:94:ASP:OD1	2.45	0.49
1:G:28:GLN:NE2	1:G:28:GLN:HA	2.27	0.49
1:A:176:ASN:C	1:A:178:LEU:H	2.16	0.49
1:H:172:LEU:O	1:H:175:LEU:HB2	2.13	0.49
1:I:179:PRO:O	1:I:181:GLN:N	2.46	0.49
1:C:140:GLU:HB2	1:D:126:TYR:CD2	2.48	0.49
1:E:168:ASP:HB3	1:E:171:TYR:HB3	1.94	0.49
1:A:139:LEU:O	1:A:143:ARG:HG3	2.12	0.49
1:A:25:CYS:HA	1:B:30:GLN:HG2	1.94	0.49
1:E:163:GLY:O	1:E:166:ILE:HG22	2.13	0.49
1:G:179:PRO:O	1:G:181:GLN:N	2.46	0.49
1:F:143:ARG:NH1	1:G:94:ASP:OD1	2.46	0.49
1:H:139:LEU:O	1:H:143:ARG:HG3	2.12	0.49
1:J:173:ARG:HH11	1:J:173:ARG:HB3	1.78	0.49
1:C:131:GLY:O	1:C:132:LEU:O	2.31	0.48
1:I:91:GLN:HE21	1:I:143:ARG:HE	1.61	0.48
1:A:168:ASP:HB3	1:A:171:TYR:HB3	1.94	0.48
1:E:187:ASP:OD1	1:E:189:THR:HG23	2.12	0.48
1:H:204:ARG:HD2	1:J:36:TYR:OH	2.13	0.48
1:J:85:ALA:HB1	1:K:120:TYR:O	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:184:LEU:N	1:K:184:LEU:HD22	2.25	0.48
1:E:112:ARG:NH2	1:E:114:GLN:HE21	2.12	0.48
1:J:163:GLY:O	1:J:166:ILE:HG22	2.14	0.48
1:K:43:LEU:HA	1:K:161:ALA:HB3	1.95	0.48
1:B:112:ARG:HG3	1:B:122:GLU:HB2	1.96	0.48
1:C:168:ASP:HB3	1:C:171:TYR:HB3	1.96	0.48
1:F:176:ASN:C	1:F:178:LEU:H	2.17	0.48
1:G:104:TYR:CD1	1:G:104:TYR:N	2.81	0.48
1:H:187:ASP:OD1	1:H:189:THR:HG23	2.14	0.48
1:K:46:ARG:HD3	1:K:171:TYR:CD1	2.49	0.48
1:C:104:TYR:CD1	1:C:104:TYR:N	2.82	0.48
1:C:193:ARG:HD3	1:D:50:GLU:OE2	2.13	0.48
1:F:182:LEU:CG	1:F:183:PRO:HD2	2.41	0.48
1:H:147:VAL:HG21	1:I:124:VAL:HG22	1.95	0.48
1:B:39:ILE:HG23	1:B:78:MET:HE1	1.96	0.48
1:F:173:ARG:HB3	1:F:173:ARG:HH11	1.78	0.48
1:C:179:PRO:O	1:C:181:GLN:N	2.47	0.48
1:F:163:GLY:O	1:F:166:ILE:HG22	2.13	0.48
1:B:46:ARG:HD3	1:B:171:TYR:CD1	2.48	0.48
1:J:182:LEU:CG	1:J:183:PRO:HD2	2.43	0.48
1:C:176:ASN:C	1:C:178:LEU:H	2.17	0.47
1:B:25:CYS:HA	1:C:30:GLN:HG2	1.96	0.47
1:B:204:ARG:NH2	1:C:78:MET:O	2.47	0.47
1:D:70:ARG:NH1	1:D:70:ARG:HG3	2.28	0.47
1:G:67:GLU:O	1:G:70:ARG:HB2	2.13	0.47
1:F:179:PRO:O	1:F:181:GLN:N	2.47	0.47
1:A:110:PHE:CE2	1:A:124:VAL:HG13	2.49	0.47
1:A:50:GLU:CG	1:A:181:GLN:HE22	2.27	0.47
1:C:163:GLY:O	1:C:166:ILE:HG22	2.14	0.47
1:G:127:GLY:HA3	1:G:142:ALA:O	2.13	0.47
1:H:179:PRO:O	1:H:181:GLN:N	2.47	0.47
1:C:25:CYS:HA	1:D:30:GLN:HG2	1.96	0.47
1:G:43:LEU:HG	1:G:78:MET:HE2	1.95	0.47
1:H:104:TYR:N	1:H:104:TYR:CD1	2.80	0.47
1:D:135:LYS:HE3	1:E:95:PHE:CG	2.48	0.47
1:D:140:GLU:HB2	1:E:126:TYR:CD2	2.50	0.47
1:D:168:ASP:HB3	1:D:171:TYR:HB3	1.95	0.47
1:D:183:PRO:HB3	1:E:170:ASP:OD1	2.15	0.47
1:D:208:CYS:HB2	1:F:33:ALA:HB2	1.95	0.47
1:F:28:GLN:NE2	1:F:28:GLN:HA	2.27	0.47
1:E:104:TYR:CD1	1:E:104:TYR:N	2.83	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:151:LEU:HD12	1:G:151:LEU:O	2.14	0.47
1:I:74:LEU:HD13	1:I:161:ALA:O	2.14	0.47
1:F:104:TYR:N	1:F:104:TYR:CD1	2.83	0.47
1:I:186:VAL:HB	1:J:174:SER:OG	2.15	0.47
1:K:47:LEU:HD12	1:K:51:TYR:CD1	2.49	0.47
1:E:131:GLY:O	1:E:132:LEU:O	2.33	0.47
1:H:176:ASN:C	1:H:178:LEU:H	2.18	0.47
1:K:91:GLN:HE21	1:K:143:ARG:HE	1.58	0.47
1:A:71:VAL:HG12	1:A:155:LEU:HD13	1.97	0.47
1:D:50:GLU:CG	1:D:181:GLN:HE22	2.28	0.47
1:E:28:GLN:HA	1:E:28:GLN:NE2	2.26	0.47
1:K:192:LYS:NZ	1:K:197:GLU:OE2	2.35	0.47
1:K:39:ILE:HG23	1:K:78:MET:HE1	1.95	0.47
1:C:182:LEU:HD12	1:C:183:PRO:HD2	1.96	0.47
1:C:204:ARG:HD2	1:E:36:TYR:OH	2.15	0.47
1:K:172:LEU:O	1:K:175:LEU:HB2	2.15	0.47
1:D:134:SER:HB3	1:D:137:LEU:HB2	1.96	0.47
1:D:204:ARG:NH2	1:E:78:MET:O	2.48	0.47
1:D:204:ARG:O	1:D:207:SER:HB3	2.14	0.47
1:A:186:VAL:HB	1:B:174:SER:OG	2.15	0.46
1:E:127:GLY:HA3	1:E:142:ALA:O	2.15	0.46
1:G:140:GLU:HG3	1:G:144:LYS:HE3	1.97	0.46
1:H:127:GLY:HA3	1:H:142:ALA:O	2.15	0.46
1:H:195:ASP:HB3	1:I:77:GLU:HG3	1.96	0.46
1:A:204:ARG:O	1:A:207:SER:HB3	2.16	0.46
1:B:172:LEU:O	1:B:175:LEU:HB2	2.15	0.46
1:G:139:LEU:O	1:G:143:ARG:HG3	2.15	0.46
1:B:204:ARG:O	1:B:207:SER:HB3	2.16	0.46
1:E:105:VAL:HG11	1:E:139:LEU:HD23	1.96	0.46
1:A:50:GLU:N	1:A:50:GLU:OE1	2.40	0.46
1:C:182:LEU:HD12	1:C:183:PRO:CD	2.46	0.46
1:G:171:TYR:CE2	1:G:175:LEU:HD11	2.50	0.46
1:H:50:GLU:CG	1:H:181:GLN:HE22	2.25	0.46
1:H:86:HIS:HA	1:H:112:ARG:O	2.15	0.46
1:J:204:ARG:O	1:J:207:SER:HB3	2.16	0.46
1:A:120:TYR:O	1:K:85:ALA:HB1	2.15	0.46
1:E:176:ASN:C	1:E:178:LEU:H	2.18	0.46
1:J:43:LEU:HG	1:J:78:MET:CE	2.46	0.46
1:I:46:ARG:HD3	1:I:171:TYR:CD1	2.51	0.46
1:E:115:LEU:HD11	1:E:158:PHE:CE2	2.51	0.46
1:H:39:ILE:HG23	1:H:78:MET:HE1	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:143:ARG:NH1	1:D:94:ASP:OD1	2.49	0.46
1:E:151:LEU:HD12	1:E:151:LEU:O	2.15	0.46
1:G:147:VAL:HG21	1:H:124:VAL:HG22	1.97	0.46
1:K:179:PRO:O	1:K:181:GLN:N	2.48	0.46
1:K:187:ASP:OD1	1:K:189:THR:HG23	2.16	0.46
1:K:43:LEU:HG	1:K:78:MET:CE	2.45	0.46
1:A:165:CYS:O	1:A:167:LEU:N	2.49	0.46
1:B:50:GLU:CG	1:B:181:GLN:HE22	2.27	0.46
1:D:165:CYS:O	1:D:167:LEU:N	2.49	0.46
1:F:156:ARG:NH1	1:F:167:LEU:HD21	2.30	0.46
1:I:135:LYS:HE3	1:J:95:PHE:CG	2.51	0.46
1:C:46:ARG:HD3	1:C:171:TYR:CD1	2.51	0.46
1:D:131:GLY:O	1:D:132:LEU:O	2.34	0.46
1:J:173:ARG:NH1	1:J:173:ARG:CB	2.79	0.46
1:B:163:GLY:O	1:B:166:ILE:HG22	2.16	0.45
1:D:91:GLN:NE2	1:E:110:PHE:HZ	2.13	0.45
1:I:147:VAL:HG21	1:J:124:VAL:HG22	1.93	0.45
1:J:104:TYR:N	1:J:104:TYR:CD1	2.84	0.45
1:J:172:LEU:O	1:J:175:LEU:HB2	2.17	0.45
1:B:183:PRO:HB3	1:C:170:ASP:OD1	2.16	0.45
1:C:183:PRO:HB3	1:D:170:ASP:OD1	2.16	0.45
1:E:140:GLU:HG3	1:E:144:LYS:HE3	1.98	0.45
1:E:43:LEU:HA	1:E:161:ALA:HB3	1.98	0.45
1:I:104:TYR:CD1	1:I:104:TYR:N	2.84	0.45
1:J:173:ARG:O	1:J:177:LYS:HG3	2.16	0.45
1:A:33:ALA:HB2	1:J:208:CYS:HB2	1.98	0.45
1:A:140:GLU:HG3	1:A:144:LYS:HE3	1.99	0.45
1:C:173:ARG:HB3	1:C:173:ARG:HH11	1.82	0.45
1:F:43:LEU:HG	1:F:78:MET:CE	2.47	0.45
1:K:50:GLU:CG	1:K:181:GLN:HE22	2.26	0.45
1:D:112:ARG:NH2	1:D:114:GLN:HE21	2.14	0.45
1:D:46:ARG:HD3	1:D:171:TYR:CD1	2.51	0.45
1:G:93:VAL:HA	1:G:107:VAL:HG22	1.97	0.45
1:I:70:ARG:HG3	1:I:70:ARG:NH1	2.30	0.45
1:A:141:LYS:HE2	1:A:145:GLU:OE2	2.17	0.45
1:B:179:PRO:O	1:B:181:GLN:N	2.49	0.45
1:F:173:ARG:NH1	1:F:173:ARG:CB	2.80	0.45
1:G:173:ARG:NH1	1:G:173:ARG:CB	2.80	0.45
1:G:172:LEU:O	1:G:175:LEU:HB2	2.17	0.45
1:F:43:LEU:HA	1:F:161:ALA:HB3	1.97	0.45
1:H:70:ARG:NH1	1:H:70:ARG:HG3	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:196:LEU:HD12	1:C:197:GLU:H	1.81	0.45
1:J:173:ARG:HH11	1:J:173:ARG:CB	2.29	0.45
1:J:179:PRO:O	1:J:181:GLN:N	2.49	0.45
1:K:168:ASP:HB3	1:K:171:TYR:HB3	1.99	0.45
1:A:104:TYR:CD1	1:A:104:TYR:N	2.85	0.45
1:B:135:LYS:HE3	1:C:95:PHE:CG	2.52	0.45
1:E:28:GLN:CA	1:E:28:GLN:HE21	2.23	0.45
1:F:184:LEU:H	1:F:184:LEU:CD2	2.28	0.45
1:I:43:LEU:HG	1:I:78:MET:HE2	1.99	0.45
1:A:98:LEU:HD12	1:A:102:LYS:O	2.17	0.45
1:B:173:ARG:HB3	1:B:173:ARG:HH11	1.82	0.45
1:D:182:LEU:CD1	1:D:183:PRO:HD2	2.46	0.45
1:A:179:PRO:O	1:A:181:GLN:N	2.50	0.45
1:E:147:VAL:HG21	1:F:124:VAL:HG22	1.97	0.45
1:E:182:LEU:CD1	1:E:183:PRO:HD2	2.47	0.45
1:I:168:ASP:HB3	1:I:171:TYR:HB3	1.99	0.45
1:D:112:ARG:HH21	1:D:114:GLN:NE2	2.14	0.44
1:B:55:ARG:HD3	1:B:65:TYR:CE2	2.52	0.44
1:C:25:CYS:O	1:C:29:CYS:SG	2.67	0.44
1:D:172:LEU:O	1:D:175:LEU:HB2	2.17	0.44
1:J:39:ILE:HG23	1:J:78:MET:HE1	1.97	0.44
1:A:199:SER:HB3	2:A:218:HOH:O	2.16	0.44
1:D:104:TYR:N	1:D:104:TYR:CD1	2.86	0.44
1:C:204:ARG:NH2	1:D:78:MET:O	2.51	0.44
1:G:182:LEU:HD12	1:G:183:PRO:HD3	1.93	0.44
1:G:70:ARG:NH1	1:G:70:ARG:HG3	2.31	0.44
1:H:105:VAL:HG11	1:H:139:LEU:HD23	1.99	0.44
1:J:86:HIS:HA	1:J:112:ARG:O	2.17	0.44
1:D:173:ARG:HB3	1:D:173:ARG:HH11	1.82	0.44
1:E:135:LYS:HE3	1:F:95:PHE:CD2	2.53	0.44
1:I:139:LEU:O	1:I:143:ARG:HG3	2.16	0.44
1:H:67:GLU:O	1:H:70:ARG:HB2	2.18	0.44
1:B:104:TYR:N	1:B:104:TYR:CD1	2.85	0.44
1:E:172:LEU:O	1:E:175:LEU:HB2	2.17	0.44
1:I:28:GLN:NE2	1:I:28:GLN:HA	2.30	0.44
1:J:46:ARG:HD3	1:J:171:TYR:CD1	2.53	0.44
1:J:47:LEU:HD12	1:J:51:TYR:CD1	2.52	0.44
1:B:43:LEU:HG	1:B:78:MET:CE	2.48	0.44
1:G:184:LEU:H	1:G:184:LEU:CD2	2.28	0.44
1:A:182:LEU:CG	1:A:183:PRO:HD2	2.45	0.44
1:A:43:LEU:HA	1:A:161:ALA:HB3	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:163:GLY:O	1:G:166:ILE:HG22	2.17	0.44
1:I:141:LYS:HE2	1:I:145:GLU:OE2	2.18	0.44
1:J:74:LEU:HD13	1:J:161:ALA:O	2.18	0.44
1:H:182:LEU:CD1	1:H:183:PRO:HD2	2.48	0.43
1:I:81:TYR:CE1	1:I:82:ASN:HB3	2.53	0.43
1:D:112:ARG:NH2	1:D:114:GLN:NE2	2.66	0.43
1:F:186:VAL:HB	1:G:174:SER:OG	2.18	0.43
1:C:50:GLU:CG	1:C:181:GLN:HE22	2.31	0.43
1:G:182:LEU:HG	1:G:183:PRO:HD2	1.98	0.43
1:I:135:LYS:HE3	1:J:95:PHE:CD2	2.53	0.43
1:I:187:ASP:OD1	1:I:189:THR:HG23	2.18	0.43
1:C:184:LEU:CD2	1:C:184:LEU:H	2.30	0.43
1:E:173:ARG:HH11	1:E:173:ARG:HB3	1.84	0.43
1:F:53:SER:O	1:F:64:CYS:HA	2.19	0.43
1:H:184:LEU:H	1:H:184:LEU:CD2	2.27	0.43
1:I:50:GLU:CG	1:I:181:GLN:HE22	2.26	0.43
1:J:168:ASP:HB3	1:J:171:TYR:HB3	2.00	0.43
1:A:205:TYR:CE1	1:C:33:ALA:HB1	2.53	0.43
1:E:46:ARG:HD3	1:E:171:TYR:CD1	2.54	0.43
1:F:140:GLU:O	1:F:144:LYS:HG3	2.18	0.43
1:I:39:ILE:HG23	1:I:78:MET:HE3	1.98	0.43
1:J:43:LEU:HA	1:J:161:ALA:HB3	2.01	0.43
1:J:184:LEU:H	1:J:184:LEU:CD2	2.30	0.43
1:B:131:GLY:O	1:B:132:LEU:O	2.36	0.43
1:E:193:ARG:HB3	1:F:50:GLU:OE2	2.18	0.43
1:I:171:TYR:CE2	1:I:175:LEU:HD11	2.53	0.43
1:J:50:GLU:CG	1:J:181:GLN:HE22	2.23	0.43
1:K:158:PHE:HB2	1:K:162:LEU:HD12	2.00	0.43
1:J:67:GLU:O	1:J:70:ARG:HB2	2.19	0.43
1:K:176:ASN:C	1:K:178:LEU:H	2.21	0.43
1:A:193:ARG:HD3	1:B:50:GLU:OE2	2.19	0.43
1:E:98:LEU:HD12	1:E:102:LYS:O	2.19	0.43
1:E:86:HIS:CE1	1:E:151:LEU:HD22	2.53	0.43
1:H:151:LEU:O	1:H:151:LEU:HD12	2.18	0.43
1:H:43:LEU:HA	1:H:161:ALA:HB3	2.00	0.43
1:I:204:ARG:NH2	1:J:78:MET:O	2.51	0.43
1:D:55:ARG:HD3	1:D:65:TYR:CD2	2.54	0.43
1:G:46:ARG:HD3	1:G:171:TYR:CD1	2.53	0.43
1:I:103:PHE:CE1	1:I:135:LYS:HB2	2.54	0.43
1:I:184:LEU:CD2	1:I:184:LEU:H	2.29	0.43
1:I:61:GLN:N	1:I:61:GLN:OE1	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:173:ARG:HB3	1:K:173:ARG:HH11	1.83	0.43
1:C:173:ARG:NH1	1:C:173:ARG:CB	2.82	0.43
1:D:93:VAL:HA	1:D:107:VAL:HG22	2.01	0.43
1:F:127:GLY:HA3	1:F:142:ALA:O	2.19	0.43
1:G:168:ASP:HB3	1:G:171:TYR:CB	2.48	0.43
1:B:165:CYS:O	1:B:167:LEU:N	2.52	0.42
1:E:112:ARG:HH21	1:E:114:GLN:NE2	2.17	0.42
1:F:55:ARG:HD3	1:F:65:TYR:CD2	2.54	0.42
1:I:50:GLU:N	1:I:50:GLU:OE1	2.42	0.42
1:J:28:GLN:HA	1:J:28:GLN:NE2	2.30	0.42
1:J:147:VAL:CG2	1:K:124:VAL:CG2	2.90	0.42
1:F:72:ILE:HG22	1:F:76:ASN:ND2	2.35	0.42
1:G:86:HIS:CE1	1:G:151:LEU:HD22	2.55	0.42
1:B:47:LEU:HD12	1:B:51:TYR:CD1	2.54	0.42
1:D:195:ASP:HB3	1:E:77:GLU:HG3	2.01	0.42
1:C:52:ILE:HG22	1:C:53:SER:N	2.34	0.42
1:F:172:LEU:O	1:F:175:LEU:HB2	2.19	0.42
1:D:173:ARG:NH1	1:D:173:ARG:CB	2.82	0.42
1:E:171:TYR:CE2	1:E:175:LEU:HD11	2.55	0.42
1:F:131:GLY:O	1:F:132:LEU:O	2.37	0.42
1:F:50:GLU:CG	1:F:181:GLN:HE22	2.26	0.42
1:J:182:LEU:CD1	1:J:183:PRO:HD2	2.50	0.42
1:E:87:SER:HA	1:F:122:GLU:HB3	2.02	0.42
1:G:173:ARG:CB	1:G:173:ARG:HH11	2.32	0.42
1:A:105:VAL:HG11	1:A:139:LEU:HD23	2.01	0.42
1:B:193:ARG:HG3	1:B:193:ARG:HH11	1.84	0.42
1:G:128:VAL:HG12	1:G:129:SER:N	2.35	0.42
1:J:110:PHE:CE2	1:J:124:VAL:HG13	2.55	0.42
1:J:47:LEU:N	1:J:47:LEU:HD22	2.35	0.42
1:J:147:VAL:HG21	1:K:124:VAL:HG23	1.96	0.42
1:J:173:ARG:NH1	1:J:173:ARG:HB2	2.34	0.42
1:B:103:PHE:CE1	1:B:135:LYS:HB2	2.54	0.42
1:E:128:VAL:HG12	1:E:129:SER:N	2.35	0.42
1:A:95:PHE:CE2	1:K:135:LYS:HG2	2.55	0.41
1:F:86:HIS:CE1	1:F:151:LEU:HD22	2.55	0.41
1:J:70:ARG:HG3	1:J:70:ARG:NH1	2.32	0.41
1:K:182:LEU:CD1	1:K:183:PRO:HD2	2.49	0.41
1:A:127:GLY:HA3	1:A:142:ALA:O	2.20	0.41
1:A:184:LEU:H	1:A:184:LEU:CD2	2.31	0.41
1:F:43:LEU:HG	1:F:78:MET:HE1	2.02	0.41
1:G:175:LEU:O	1:G:178:LEU:HB2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:47:LEU:HD22	1:I:47:LEU:N	2.35	0.41
1:C:98:LEU:HD12	1:C:102:LYS:O	2.20	0.41
1:G:165:CYS:O	1:G:168:ASP:N	2.52	0.41
1:H:98:LEU:HD12	1:H:102:LYS:O	2.20	0.41
1:G:140:GLU:O	1:G:144:LYS:HG3	2.20	0.41
1:G:74:LEU:HD13	1:G:161:ALA:O	2.20	0.41
1:A:52:ILE:HG22	1:A:53:SER:N	2.35	0.41
1:D:67:GLU:O	1:D:70:ARG:HB2	2.21	0.41
1:H:168:ASP:HB3	1:H:171:TYR:CB	2.50	0.41
1:A:103:PHE:CE1	1:A:135:LYS:HB2	2.56	0.41
1:B:176:ASN:C	1:B:178:LEU:H	2.24	0.41
1:F:105:VAL:HG11	1:F:139:LEU:HD23	2.03	0.41
1:E:186:VAL:HB	1:F:174:SER:OG	2.21	0.41
1:H:74:LEU:HD13	1:H:161:ALA:O	2.21	0.41
1:A:195:ASP:HB3	1:B:77:GLU:HG3	2.02	0.41
1:D:165:CYS:O	1:D:168:ASP:N	2.52	0.41
1:E:204:ARG:O	1:E:207:SER:HB3	2.21	0.41
1:I:86:HIS:HA	1:I:112:ARG:O	2.20	0.41
1:C:110:PHE:CE2	1:C:124:VAL:HG13	2.55	0.41
1:C:54:SER:CB	1:C:62:LYS:HE3	2.51	0.41
1:D:127:GLY:HA3	1:D:142:ALA:O	2.21	0.41
1:D:140:GLU:O	1:D:144:LYS:HG3	2.21	0.41
1:E:53:SER:O	1:E:64:CYS:HA	2.21	0.41
1:F:140:GLU:HG3	1:F:144:LYS:HE3	2.02	0.41
1:G:55:ARG:HD3	1:G:65:TYR:CD2	2.56	0.41
1:G:85:ALA:CB	1:H:120:TYR:O	2.68	0.41
1:I:67:GLU:O	1:I:70:ARG:HB2	2.20	0.41
1:J:171:TYR:CE2	1:J:175:LEU:HD11	2.56	0.41
1:B:79:PHE:O	1:B:83:GLY:HA3	2.20	0.41
1:K:43:LEU:HA	1:K:161:ALA:CB	2.51	0.41
1:A:193:ARG:HH11	1:A:193:ARG:HG3	1.86	0.41
1:D:61:GLN:N	1:D:61:GLN:OE1	2.52	0.41
1:H:173:ARG:HH11	1:H:173:ARG:HB3	1.85	0.41
1:I:163:GLY:O	1:I:166:ILE:HG22	2.21	0.41
1:I:204:ARG:HD2	1:K:36:TYR:OH	2.21	0.41
1:C:103:PHE:CE1	1:C:135:LYS:HB2	2.56	0.41
1:F:204:ARG:O	1:F:207:SER:HB3	2.21	0.41
1:I:70:ARG:CG	1:I:70:ARG:HH11	2.31	0.41
1:B:28:GLN:HA	1:B:28:GLN:NE2	2.30	0.40
1:D:143:ARG:NH1	1:E:94:ASP:OD1	2.53	0.40
1:G:200:VAL:HG13	1:H:39:ILE:HG13	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:204:ARG:O	1:I:207:SER:HB3	2.20	0.40
1:J:25:CYS:SG	1:J:28:GLN:HG3	2.62	0.40
1:K:184:LEU:H	1:K:184:LEU:CD2	2.31	0.40
1:B:33:ALA:HB2	1:K:208:CYS:CB	2.51	0.40
1:C:86:HIS:CE1	1:C:151:LEU:HD22	2.56	0.40
1:F:47:LEU:HD12	1:F:51:TYR:CD1	2.56	0.40
1:I:52:ILE:HG22	1:I:53:SER:N	2.36	0.40
1:B:182:LEU:CD1	1:B:183:PRO:HD2	2.51	0.40
1:B:54:SER:HA	1:B:63:VAL:O	2.20	0.40
1:E:55:ARG:HD3	1:E:65:TYR:CD2	2.56	0.40
1:G:165:CYS:O	1:G:167:LEU:N	2.54	0.40
1:K:139:LEU:O	1:K:143:ARG:HG3	2.20	0.40
1:A:186:VAL:HG21	1:B:171:TYR:HA	2.02	0.40
1:F:182:LEU:CD1	1:F:183:PRO:HD2	2.51	0.40
1:F:144:LYS:NZ	1:G:149:ASP:OD2	2.38	0.40
1:H:188:LEU:HD12	1:I:178:LEU:HD11	2.02	0.40
1:A:182:LEU:CD1	1:A:183:PRO:HD2	2.52	0.40
1:C:127:GLY:HA3	1:C:142:ALA:O	2.22	0.40
1:C:135:LYS:HG2	1:D:95:PHE:CE2	2.56	0.40
1:C:180:ARG:CZ	1:C:181:GLN:O	2.69	0.40
1:E:70:ARG:NH1	1:E:70:ARG:HG3	2.35	0.40
1:I:110:PHE:CE2	1:I:124:VAL:HG13	2.57	0.40
1:K:173:ARG:NH1	1:K:173:ARG:CB	2.84	0.40
1:K:204:ARG:O	1:K:207:SER:HB3	2.21	0.40

All (2) 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:60:GLY:O	1:G:37:GLN:OE1[6_456]	2.14	0.06
1:E:185:GLU:OE2	1:E:185:GLU:OE2[8_666]	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	182/212 (86%)	165 (91%)	11 (6%)	6 (3%)	4	13
1	B	182/212 (86%)	166 (91%)	12 (7%)	4 (2%)	6	21
1	C	182/212 (86%)	165 (91%)	11 (6%)	6 (3%)	4	13
1	D	182/212 (86%)	166 (91%)	10 (6%)	6 (3%)	4	13
1	E	182/212 (86%)	163 (90%)	12 (7%)	7 (4%)	3	10
1	F	182/212 (86%)	158 (87%)	18 (10%)	6 (3%)	4	13
1	G	182/212 (86%)	162 (89%)	13 (7%)	7 (4%)	3	10
1	H	182/212 (86%)	162 (89%)	14 (8%)	6 (3%)	4	13
1	I	182/212 (86%)	165 (91%)	11 (6%)	6 (3%)	4	13
1	J	182/212 (86%)	167 (92%)	10 (6%)	5 (3%)	5	16
1	K	182/212 (86%)	164 (90%)	13 (7%)	5 (3%)	5	16
All	All	2002/2332 (86%)	1803 (90%)	135 (7%)	64 (3%)	4	13

All (64) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	132	LEU
1	B	132	LEU
1	C	132	LEU
1	D	132	LEU
1	E	132	LEU
1	F	132	LEU
1	G	132	LEU
1	H	132	LEU
1	I	132	LEU
1	J	132	LEU
1	K	132	LEU
1	A	180	ARG
1	D	166	ILE
1	D	180	ARG
1	E	180	ARG
1	I	180	ARG
1	A	166	ILE
1	A	177	LYS
1	C	180	ARG
1	D	177	LYS

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Mol	Chain	Res	Type
1	F	180	ARG
1	G	180	ARG
1	H	58	GLY
1	H	180	ARG
1	I	177	LYS
1	J	177	LYS
1	J	180	ARG
1	K	58	GLY
1	K	180	ARG
1	B	180	ARG
1	C	58	GLY
1	C	177	LYS
1	E	177	LYS
1	F	177	LYS
1	F	196	LEU
1	G	29	CYS
1	G	58	GLY
1	G	100	ASN
1	H	177	LYS
1	J	58	GLY
1	A	45	GLN
1	A	58	GLY
1	B	58	GLY
1	B	166	ILE
1	C	100	ASN
1	E	100	ASN
1	E	196	LEU
1	F	58	GLY
1	G	177	LYS
1	H	29	CYS
1	I	58	GLY
1	I	100	ASN
1	K	177	LYS
1	D	58	GLY
1	E	58	GLY
1	E	166	ILE
1	F	166	ILE
1	G	166	ILE
1	K	100	ASN
1	D	183	PRO
1	H	166	ILE
1	I	179	PRO

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Mol	Chain	Res	Type
1	J	166	ILE
1	C	183	PRO

5.3.2 Protein sidechains [i](#)

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	153/172 (89%)	147 (96%)	6 (4%)	32 63
1	B	153/172 (89%)	147 (96%)	6 (4%)	32 63
1	C	153/172 (89%)	146 (95%)	7 (5%)	27 56
1	D	153/172 (89%)	146 (95%)	7 (5%)	27 56
1	E	153/172 (89%)	145 (95%)	8 (5%)	23 51
1	F	153/172 (89%)	147 (96%)	6 (4%)	32 63
1	G	153/172 (89%)	145 (95%)	8 (5%)	23 51
1	H	153/172 (89%)	146 (95%)	7 (5%)	27 56
1	I	153/172 (89%)	148 (97%)	5 (3%)	38 68
1	J	153/172 (89%)	147 (96%)	6 (4%)	32 63
1	K	153/172 (89%)	147 (96%)	6 (4%)	32 63
All	All	1683/1892 (89%)	1611 (96%)	72 (4%)	29 59

All (72) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	GLN
1	A	112	ARG
1	A	124	VAL
1	A	180	ARG
1	A	183	PRO
1	A	184	LEU
1	B	28	GLN
1	B	112	ARG
1	B	124	VAL

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Mol	Chain	Res	Type
1	B	180	ARG
1	B	183	PRO
1	B	184	LEU
1	C	28	GLN
1	C	104	TYR
1	C	112	ARG
1	C	124	VAL
1	C	180	ARG
1	C	183	PRO
1	C	184	LEU
1	D	28	GLN
1	D	86	HIS
1	D	112	ARG
1	D	124	VAL
1	D	180	ARG
1	D	183	PRO
1	D	184	LEU
1	E	28	GLN
1	E	86	HIS
1	E	104	TYR
1	E	112	ARG
1	E	124	VAL
1	E	180	ARG
1	E	183	PRO
1	E	184	LEU
1	F	28	GLN
1	F	112	ARG
1	F	124	VAL
1	F	180	ARG
1	F	183	PRO
1	F	184	LEU
1	G	28	GLN
1	G	86	HIS
1	G	104	TYR
1	G	112	ARG
1	G	124	VAL
1	G	180	ARG
1	G	183	PRO
1	G	184	LEU
1	H	28	GLN
1	H	104	TYR
1	H	112	ARG

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Mol	Chain	Res	Type
1	H	124	VAL
1	H	180	ARG
1	H	183	PRO
1	H	184	LEU
1	I	112	ARG
1	I	124	VAL
1	I	180	ARG
1	I	183	PRO
1	I	184	LEU
1	J	28	GLN
1	J	112	ARG
1	J	124	VAL
1	J	180	ARG
1	J	183	PRO
1	J	184	LEU
1	K	28	GLN
1	K	112	ARG
1	K	124	VAL
1	K	180	ARG
1	K	183	PRO
1	K	184	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (85) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	28	GLN
1	A	37	GLN
1	A	45	GLN
1	A	86	HIS
1	A	91	GLN
1	A	114	GLN
1	A	121	HIS
1	A	181	GLN
1	B	28	GLN
1	B	37	GLN
1	B	86	HIS
1	B	91	GLN
1	B	121	HIS
1	B	181	GLN
1	C	28	GLN
1	C	37	GLN
1	C	45	GLN

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Mol	Chain	Res	Type
1	C	86	HIS
1	C	91	GLN
1	C	121	HIS
1	C	181	GLN
1	D	28	GLN
1	D	37	GLN
1	D	73	ASN
1	D	86	HIS
1	D	91	GLN
1	D	114	GLN
1	D	121	HIS
1	D	181	GLN
1	D	206	ASN
1	E	28	GLN
1	E	37	GLN
1	E	45	GLN
1	E	73	ASN
1	E	86	HIS
1	E	91	GLN
1	E	114	GLN
1	E	121	HIS
1	E	181	GLN
1	E	206	ASN
1	F	28	GLN
1	F	37	GLN
1	F	45	GLN
1	F	73	ASN
1	F	86	HIS
1	F	91	GLN
1	F	114	GLN
1	F	181	GLN
1	G	28	GLN
1	G	37	GLN
1	G	45	GLN
1	G	73	ASN
1	G	86	HIS
1	G	91	GLN
1	G	114	GLN
1	G	121	HIS
1	G	181	GLN
1	H	28	GLN
1	H	37	GLN

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Mol	Chain	Res	Type
1	H	45	GLN
1	H	86	HIS
1	H	91	GLN
1	H	121	HIS
1	H	181	GLN
1	I	28	GLN
1	I	37	GLN
1	I	45	GLN
1	I	86	HIS
1	I	91	GLN
1	I	121	HIS
1	I	181	GLN
1	J	28	GLN
1	J	37	GLN
1	J	45	GLN
1	J	86	HIS
1	J	91	GLN
1	J	121	HIS
1	J	181	GLN
1	K	28	GLN
1	K	37	GLN
1	K	45	GLN
1	K	86	HIS
1	K	91	GLN
1	K	121	HIS
1	K	181	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	184/212 (86%)	-0.25	2 (1%) 80 80	25, 60, 136, 198	0
1	B	184/212 (86%)	-0.13	5 (2%) 54 50	25, 56, 150, 200	0
1	C	184/212 (86%)	-0.14	6 (3%) 46 41	25, 61, 166, 200	0
1	D	184/212 (86%)	0.13	9 (4%) 29 25	28, 78, 156, 200	0
1	E	184/212 (86%)	0.55	23 (12%) 3 2	33, 85, 170, 200	0
1	F	184/212 (86%)	0.47	18 (9%) 7 5	38, 87, 163, 199	0
1	G	184/212 (86%)	0.18	8 (4%) 35 30	37, 82, 170, 200	0
1	H	184/212 (86%)	0.02	6 (3%) 46 41	30, 72, 149, 198	0
1	I	184/212 (86%)	-0.13	4 (2%) 62 59	31, 69, 139, 185	0
1	J	184/212 (86%)	-0.06	5 (2%) 54 50	31, 70, 158, 200	0
1	K	184/212 (86%)	0.09	15 (8%) 11 8	30, 65, 154, 182	0
All	All	2024/2332 (86%)	0.07	101 (4%) 28 24	25, 72, 160, 200	0

All (101) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	182	LEU	6.9
1	E	48	GLY	6.7
1	G	184	LEU	6.4
1	I	182	LEU	6.4
1	C	181	GLN	5.8
1	C	58	GLY	5.4
1	E	60	GLY	5.3
1	F	62	LYS	5.2
1	G	48	GLY	5.2
1	K	184	LEU	5.1
1	K	185	GLU	4.7
1	E	100	ASN	4.7

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Mol	Chain	Res	Type	RSRZ
1	H	182	LEU	4.7
1	K	182	LEU	4.6
1	E	181	GLN	4.5
1	B	180	ARG	4.4
1	G	59	GLY	4.4
1	E	59	GLY	4.3
1	J	181	GLN	4.3
1	E	61	GLN	4.3
1	K	61	GLN	4.3
1	D	187	ASP	4.2
1	B	61	GLN	4.1
1	K	183	PRO	4.1
1	F	59	GLY	3.9
1	F	183	PRO	3.9
1	E	183	PRO	3.8
1	F	182	LEU	3.8
1	J	180	ARG	3.6
1	B	182	LEU	3.6
1	K	207	SER	3.6
1	F	181	GLN	3.5
1	K	190	LYS	3.4
1	G	182	LEU	3.4
1	D	56	MET	3.4
1	F	176	ASN	3.4
1	H	59	GLY	3.3
1	D	179	PRO	3.3
1	J	183	PRO	3.2
1	C	184	LEU	3.2
1	E	179	PRO	3.2
1	F	179	PRO	3.2
1	E	99	ASN	3.1
1	H	177	LYS	3.1
1	E	193	ARG	3.1
1	F	97	ASP	3.0
1	E	180	ARG	3.0
1	G	180	ARG	3.0
1	J	191	ALA	2.9
1	E	47	LEU	2.9
1	E	184	LEU	2.9
1	K	28	GLN	2.9
1	K	181	GLN	2.9
1	C	62	LYS	2.9

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Mol	Chain	Res	Type	RSRZ
1	K	180	ARG	2.8
1	D	178	LEU	2.8
1	J	179	PRO	2.8
1	E	189	THR	2.8
1	F	133	LYS	2.8
1	A	180	ARG	2.7
1	F	48	GLY	2.7
1	K	178	LEU	2.7
1	F	57	ALA	2.7
1	E	58	GLY	2.7
1	F	198	PRO	2.6
1	E	185	GLU	2.6
1	E	104	TYR	2.6
1	I	177	LYS	2.6
1	F	184	LEU	2.5
1	F	180	ARG	2.5
1	G	190	LYS	2.5
1	E	177	LYS	2.4
1	E	172	LEU	2.4
1	E	25	CYS	2.4
1	E	178	LEU	2.4
1	D	59	GLY	2.3
1	F	170	ASP	2.3
1	K	189	THR	2.3
1	K	186	VAL	2.3
1	D	177	LYS	2.3
1	F	99	ASN	2.3
1	F	177	LYS	2.3
1	D	184	LEU	2.3
1	B	98	LEU	2.3
1	E	63	VAL	2.2
1	F	175	LEU	2.2
1	D	99	ASN	2.2
1	K	208	CYS	2.1
1	H	181	GLN	2.1
1	C	59	GLY	2.1
1	H	33	ALA	2.1
1	G	183	PRO	2.1
1	K	179	PRO	2.1
1	G	159	GLY	2.1
1	H	183	PRO	2.1
1	B	60	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	182	LEU	2.0
1	D	61	GLN	2.0
1	I	181	GLN	2.0
1	I	59	GLY	2.0
1	A	177	LYS	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.