



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 05:59 AM UTC

PDB ID : 2BBZ / pdb\_00002bbz  
Title : Crystal Structure of MC159 Reveals Molecular Mechanism of DISC Assembly and vFLIP Inhibition  
Authors : Yang, J.K.; Wang, L.; Zheng, L.; Wan, F.; Ahmed, M.; Lenardo, M.J.; Wu, H.  
Deposited on : 2005-10-18  
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](mailto: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>.

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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

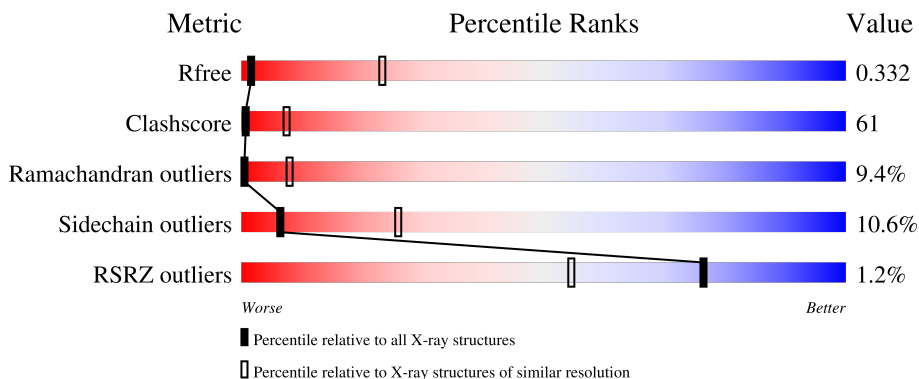
# 1 Overall quality at a glance

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  $\geq 3$ , 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	249	 2% 22% 40% 13% 24%
1	B	249	 2% 22% 41% 11% 24%
1	C	249	 22% 41% 12% 24%
1	D	249	 22% 41% 12% 24%

## 2 Entry composition i

There are 2 unique types of molecules in this entry. The entry contains 6016 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 Viral CASP8 and FADD-like apoptosis regulator.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	190	1491	937	256	287	11	0	0	0
1	B	190	1491	937	256	287	11	0	0	0
1	C	190	1491	937	256	287	11	0	0	0
1	D	190	1491	937	256	287	11	0	0	0

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	242	LEU	-	expression tag	UNP Q98325
A	243	GLU	-	expression tag	UNP Q98325
A	244	HIS	-	expression tag	UNP Q98325
A	245	HIS	-	expression tag	UNP Q98325
A	246	HIS	-	expression tag	UNP Q98325
A	247	HIS	-	expression tag	UNP Q98325
A	248	HIS	-	expression tag	UNP Q98325
A	249	HIS	-	expression tag	UNP Q98325
B	242	LEU	-	expression tag	UNP Q98325
B	243	GLU	-	expression tag	UNP Q98325
B	244	HIS	-	expression tag	UNP Q98325
B	245	HIS	-	expression tag	UNP Q98325
B	246	HIS	-	expression tag	UNP Q98325
B	247	HIS	-	expression tag	UNP Q98325
B	248	HIS	-	expression tag	UNP Q98325
B	249	HIS	-	expression tag	UNP Q98325
C	242	LEU	-	expression tag	UNP Q98325
C	243	GLU	-	expression tag	UNP Q98325
C	244	HIS	-	expression tag	UNP Q98325
C	245	HIS	-	expression tag	UNP Q98325
C	246	HIS	-	expression tag	UNP Q98325

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Chain	Residue	Modelled	Actual	Comment	Reference
C	247	HIS	-	expression tag	UNP Q98325
C	248	HIS	-	expression tag	UNP Q98325
C	249	HIS	-	expression tag	UNP Q98325
D	242	LEU	-	expression tag	UNP Q98325
D	243	GLU	-	expression tag	UNP Q98325
D	244	HIS	-	expression tag	UNP Q98325
D	245	HIS	-	expression tag	UNP Q98325
D	246	HIS	-	expression tag	UNP Q98325
D	247	HIS	-	expression tag	UNP Q98325
D	248	HIS	-	expression tag	UNP Q98325
D	249	HIS	-	expression tag	UNP Q98325

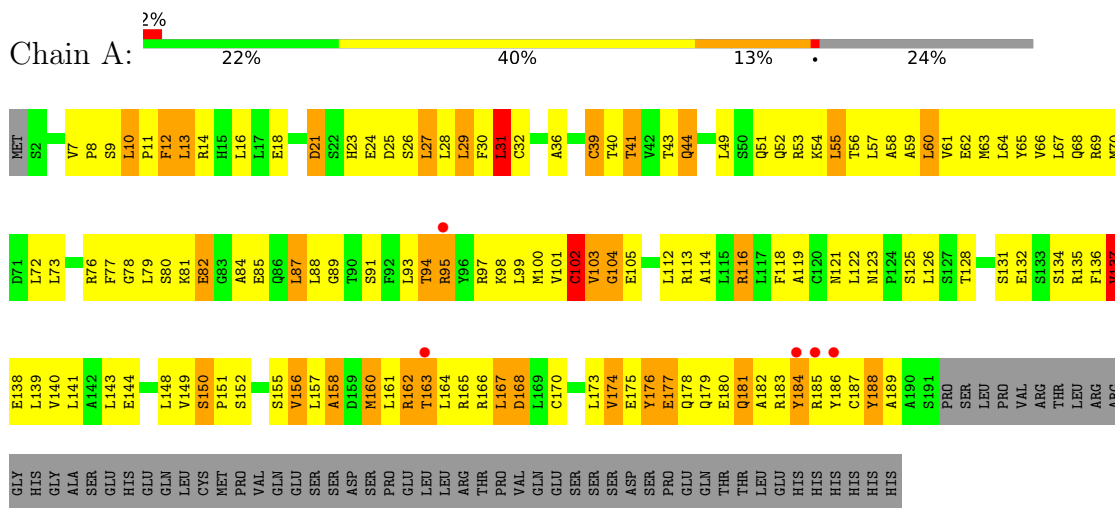
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	16	Total O 16 16	0	0
2	B	8	Total O 8 8	0	0
2	C	16	Total O 16 16	0	0
2	D	12	Total O 12 12	0	0

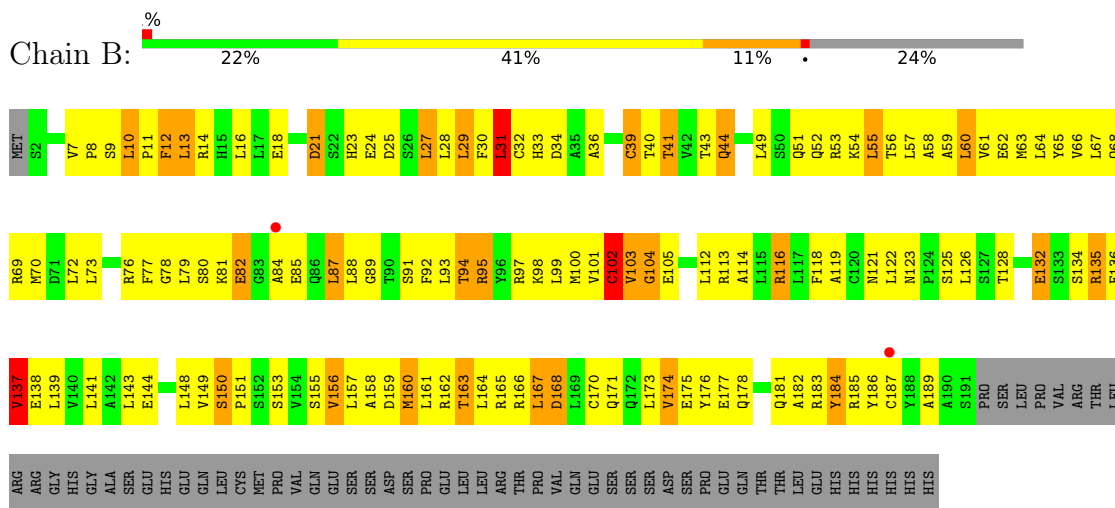
### 3 Residue-property plots

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.

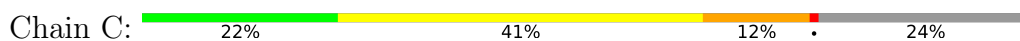
- Molecule 1: Viral CASP8 and FADD-like apoptosis regulator

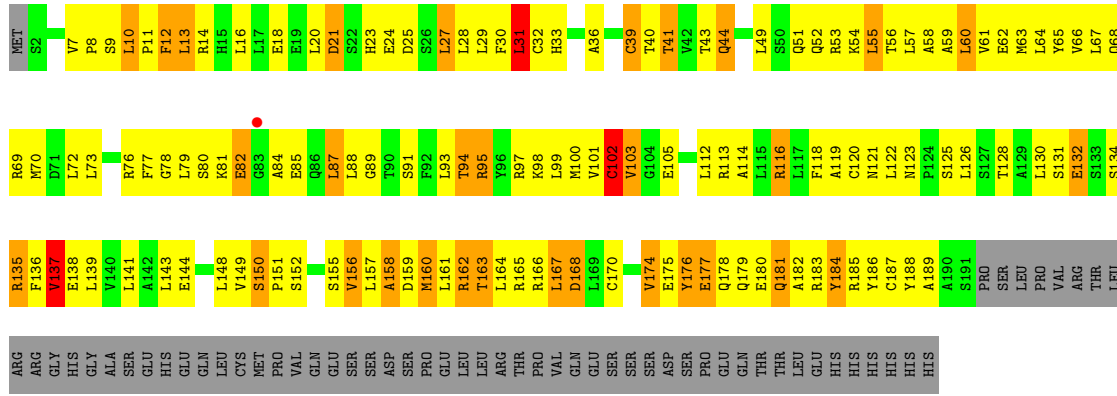


- Molecule 1: Viral CASP8 and FADD-like apoptosis regulator

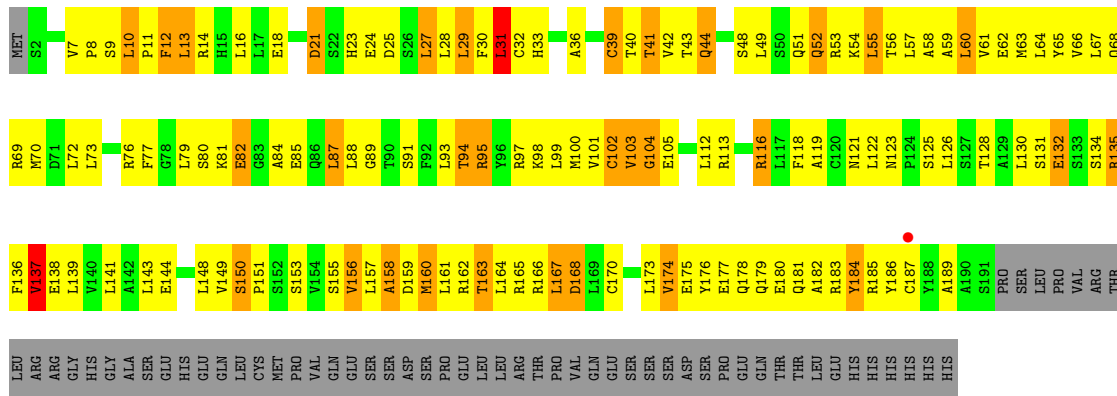
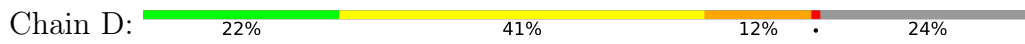


- Molecule 1: Viral CASP8 and FADD-like apoptosis regulator





• Molecule 1: Viral CASP8 and FADD-like apoptosis regulator



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	83.86Å 108.06Å 142.64Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 3.80 30.00 – 3.80	Depositor EDS
% Data completeness (in resolution range)	(Not available) (30.00-3.80) 93.5 (30.00-3.80)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.04 (at 3.75Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.325 , 0.372 0.294 , 0.332	Depositor DCC
$R_{free}$ test set	1361 reflections (10.24%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	88.7	Xtrriage
Anisotropy	0.551	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 110.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.88	EDS
Total number of atoms	6016	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	67.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.03% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	0.61	0/1510	1.05	9/2040 (0.4%)
1	B	0.52	0/1510	1.04	9/2040 (0.4%)
1	C	0.55	0/1510	1.04	10/2040 (0.5%)
1	D	0.48	0/1510	1.03	8/2040 (0.4%)
All	All	0.54	0/6040	1.04	36/8160 (0.4%)

There are no bond length outliers.

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	160	MET	N-CA-C	-11.11	99.85	113.41
1	C	160	MET	N-CA-C	-10.88	100.14	113.41
1	B	160	MET	N-CA-C	-10.88	100.14	113.41
1	D	160	MET	N-CA-C	-10.82	100.21	113.41
1	A	174	VAL	N-CA-C	-8.81	104.89	113.53
1	B	174	VAL	N-CA-C	-8.35	105.35	113.53
1	C	174	VAL	N-CA-C	-8.23	105.46	113.53
1	D	174	VAL	N-CA-C	-8.06	105.63	113.53
1	D	175	GLU	N-CA-C	-7.41	103.50	112.54
1	D	95	ARG	N-CA-C	-7.38	103.24	111.28
1	B	175	GLU	N-CA-C	-7.17	103.79	112.54
1	B	95	ARG	N-CA-C	-6.94	103.72	111.28
1	B	103	VAL	N-CA-C	-6.83	106.22	111.62
1	C	95	ARG	N-CA-C	-6.78	103.89	111.28
1	A	95	ARG	N-CA-C	-6.57	104.12	111.28
1	D	176	TYR	N-CA-C	-6.55	103.41	111.40
1	D	103	VAL	N-CA-C	-6.54	106.45	111.62
1	C	175	GLU	N-CA-C	-6.48	104.63	112.54
1	A	176	TYR	N-CA-C	-6.43	103.56	111.40
1	B	176	TYR	N-CA-C	-6.25	103.78	111.40
1	B	33	HIS	N-CA-C	-6.08	105.67	114.12
1	A	103	VAL	N-CA-C	-6.04	107.05	112.12
1	A	131	SER	N-CA-C	-5.96	99.98	109.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	176	TYR	N-CA-C	-5.92	104.18	111.40
1	A	175	GLU	N-CA-C	-5.89	105.35	112.54
1	D	33	HIS	N-CA-C	-5.83	105.94	114.39
1	C	103	VAL	N-CA-C	-5.79	107.05	111.62
1	C	114	ALA	N-CA-C	-5.70	104.68	111.69
1	C	131	SER	N-CA-C	-5.51	100.73	109.76
1	C	33	HIS	N-CA-C	-5.41	106.60	114.12
1	A	114	ALA	N-CA-C	-5.38	105.08	111.69
1	B	114	ALA	N-CA-C	-5.13	105.38	111.69
1	D	131	SER	N-CA-C	-5.12	101.36	109.76
1	A	162	ARG	N-CA-C	-5.11	105.69	112.34
1	C	162	ARG	N-CA-C	-5.09	105.72	112.34
1	B	92	PHE	N-CA-C	-5.04	106.72	112.92

There are no chirality outliers.

There are no planarity outliers.

## 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	1491	0	1518	220	0
1	B	1491	0	1518	180	0
1	C	1491	0	1518	212	1
1	D	1491	0	1518	187	1
2	A	16	0	0	17	0
2	B	8	0	0	6	0
2	C	16	0	0	12	0
2	D	12	0	0	7	0
All	All	6016	0	6072	735	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 61.

All (735) 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:181:GLN:HG2	1:C:165:ARG:HH22	1.17	1.01
1:D:177:GLU:HA	2:D:256:HOH:O	1.58	1.01
1:A:183:ARG:HG3	1:A:184:TYR:H	1.29	0.97
1:C:183:ARG:HG3	1:C:184:TYR:H	1.29	0.97
1:A:26:SER:HB3	2:A:259:HOH:O	1.64	0.97
1:D:183:ARG:HG3	1:D:184:TYR:H	1.32	0.95
1:A:167:LEU:CB	1:C:178:GLN:HG2	1.95	0.95
1:B:183:ARG:HG3	1:B:184:TYR:H	1.30	0.94
1:C:185:ARG:HD2	1:D:156:VAL:HA	1.48	0.94
1:A:165:ARG:HH12	1:C:181:GLN:CD	1.75	0.94
1:A:165:ARG:HH22	1:C:181:GLN:HG2	1.28	0.93
1:C:97:ARG:HA	2:C:251:HOH:O	1.67	0.93
1:A:167:LEU:HB2	1:C:178:GLN:HG2	1.48	0.92
1:C:27:LEU:HD12	1:C:137:VAL:HG21	1.54	0.89
1:A:167:LEU:HD22	1:A:167:LEU:H	1.39	0.88
1:C:186:TYR:CE2	1:D:159:ASP:HB2	2.08	0.88
1:C:44:GLN:HG2	2:C:255:HOH:O	1.72	0.87
1:A:178:GLN:OE1	1:C:168:ASP:HA	1.74	0.87
1:D:27:LEU:HD12	1:D:137:VAL:HG21	1.57	0.87
1:D:167:LEU:H	1:D:167:LEU:HD22	1.40	0.86
1:C:185:ARG:NE	1:D:156:VAL:HG22	1.90	0.86
1:C:167:LEU:HD22	1:C:167:LEU:H	1.38	0.86
1:A:9:SER:HB2	1:A:12:PHE:HD2	1.41	0.85
1:B:167:LEU:H	1:B:167:LEU:HD22	1.40	0.83
1:B:9:SER:HB2	1:B:12:PHE:HD2	1.43	0.83
1:D:9:SER:HB2	1:D:12:PHE:HD2	1.43	0.82
1:B:27:LEU:HD12	1:B:137:VAL:HG21	1.60	0.82
1:C:9:SER:HB2	1:C:12:PHE:HD2	1.43	0.82
1:A:168:ASP:HA	1:C:178:GLN:OE1	1.80	0.82
1:A:27:LEU:HD12	1:A:137:VAL:HG21	1.61	0.81
1:D:31:LEU:HD22	1:D:93:LEU:HD22	1.60	0.81
1:C:183:ARG:HG3	1:C:184:TYR:N	1.97	0.80
1:B:31:LEU:HD22	1:B:93:LEU:HD22	1.65	0.79
1:C:118:PHE:HA	2:C:260:HOH:O	1.82	0.78
1:B:183:ARG:HG3	1:B:184:TYR:N	1.98	0.78
1:C:31:LEU:HD22	1:C:93:LEU:HD22	1.66	0.77
1:B:53:ARG:O	1:B:53:ARG:HD3	1.84	0.77
1:A:183:ARG:HG3	1:A:184:TYR:N	1.96	0.77
1:D:53:ARG:O	1:D:53:ARG:HD3	1.85	0.76
1:D:183:ARG:HG3	1:D:184:TYR:N	1.98	0.76
1:C:53:ARG:O	1:C:53:ARG:HD3	1.85	0.76
1:C:163:THR:HG22	1:C:164:LEU:HD22	1.67	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:186:TYR:HE2	1:D:159:ASP:HB2	1.48	0.76
1:A:183:ARG:HG3	2:A:264:HOH:O	1.85	0.75
1:C:116:ARG:NH1	1:C:134:SER:O	2.19	0.75
1:C:187:CYS:C	1:C:189:ALA:H	1.94	0.75
1:C:28:LEU:HD11	1:C:63:MET:HB3	1.69	0.75
1:A:53:ARG:O	1:A:53:ARG:HD3	1.87	0.74
1:A:178:GLN:HG2	1:C:167:LEU:HB2	1.70	0.74
1:A:187:CYS:C	1:A:189:ALA:H	1.95	0.74
1:B:28:LEU:HD11	1:B:63:MET:HB3	1.70	0.74
1:B:103:VAL:C	1:B:105:GLU:H	1.94	0.74
1:C:185:ARG:CZ	1:D:156:VAL:HG22	2.18	0.74
2:A:250:HOH:O	1:B:118:PHE:CZ	2.39	0.73
1:B:187:CYS:C	1:B:189:ALA:H	1.95	0.73
1:A:181:GLN:CD	1:C:165:ARG:HH12	1.97	0.73
1:B:178:GLN:HG3	2:B:251:HOH:O	1.88	0.73
1:D:64:LEU:HD22	1:D:73:LEU:HG	1.71	0.73
1:B:163:THR:HG22	1:B:164:LEU:HD22	1.71	0.73
1:A:163:THR:HG22	1:A:164:LEU:HD22	1.71	0.72
1:C:62:GLU:HG2	1:C:88:LEU:HD23	1.71	0.72
1:D:187:CYS:C	1:D:189:ALA:H	1.95	0.72
1:B:41:THR:HG22	1:B:43:THR:H	1.56	0.71
1:B:174:VAL:HG11	1:C:163:THR:HA	1.73	0.71
1:A:31:LEU:HD22	1:A:93:LEU:HD22	1.71	0.71
1:C:41:THR:HG22	1:C:43:THR:H	1.56	0.71
1:A:132:GLU:C	2:A:263:HOH:O	2.33	0.71
1:A:163:THR:HA	1:D:174:VAL:HG11	1.73	0.71
1:D:103:VAL:C	1:D:105:GLU:H	1.98	0.71
1:D:28:LEU:HD11	1:D:63:MET:HB3	1.71	0.70
1:A:24:GLU:HB3	1:A:67:LEU:HD13	1.73	0.70
1:A:62:GLU:HG2	1:A:88:LEU:HD23	1.73	0.70
1:B:62:GLU:HG2	1:B:88:LEU:HD23	1.72	0.70
1:B:171:GLN:HB2	2:B:250:HOH:O	1.90	0.70
1:A:103:VAL:C	1:A:105:GLU:H	1.99	0.70
1:A:28:LEU:HD11	1:A:63:MET:HB3	1.74	0.70
1:C:103:VAL:C	1:C:105:GLU:H	1.98	0.70
1:A:10:LEU:HB3	1:A:14:ARG:NH2	2.07	0.70
1:A:165:ARG:NH2	1:C:181:GLN:HG2	2.04	0.69
1:D:41:THR:HG22	1:D:43:THR:H	1.57	0.69
1:C:183:ARG:O	1:C:185:ARG:N	2.25	0.69
1:A:64:LEU:HD22	1:A:73:LEU:HG	1.74	0.69
1:B:66:VAL:C	1:B:68:GLN:H	2.00	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:183:ARG:O	1:B:185:ARG:N	2.26	0.69
1:C:10:LEU:HB3	1:C:14:ARG:NH2	2.08	0.69
1:D:62:GLU:HG2	1:D:88:LEU:HD23	1.73	0.69
1:A:183:ARG:N	2:A:252:HOH:O	2.26	0.69
1:A:41:THR:HG22	1:A:43:THR:H	1.56	0.69
1:B:49:LEU:HB3	1:B:55:LEU:HB2	1.75	0.69
1:B:64:LEU:HD22	1:B:73:LEU:HG	1.75	0.68
1:D:24:GLU:HB3	1:D:67:LEU:HD13	1.75	0.68
1:B:116:ARG:NH1	1:B:134:SER:O	2.27	0.67
1:C:188:TYR:HB3	1:D:122:LEU:HA	1.77	0.67
1:B:178:GLN:CG	2:B:251:HOH:O	2.40	0.67
1:D:163:THR:HG22	1:D:164:LEU:HD22	1.75	0.67
1:B:10:LEU:HB3	1:B:14:ARG:NH2	2.09	0.67
1:C:64:LEU:HD22	1:C:73:LEU:HG	1.76	0.67
1:C:135:ARG:HD3	2:C:257:HOH:O	1.92	0.67
1:A:23:HIS:O	2:A:259:HOH:O	2.13	0.67
1:C:183:ARG:C	1:C:185:ARG:H	2.02	0.67
1:D:49:LEU:HB3	1:D:55:LEU:HB2	1.77	0.66
1:A:49:LEU:HB3	1:A:55:LEU:HB2	1.78	0.66
1:D:66:VAL:C	1:D:68:GLN:H	2.02	0.65
1:B:81:LYS:O	1:B:81:LYS:HD3	1.96	0.65
1:C:81:LYS:HD3	1:C:81:LYS:O	1.96	0.65
1:A:181:GLN:HG2	1:C:165:ARG:NH2	2.02	0.65
1:B:183:ARG:C	1:B:185:ARG:H	2.04	0.65
1:A:188:TYR:HB3	1:B:122:LEU:HA	1.78	0.65
1:D:116:ARG:NH1	1:D:134:SER:O	2.29	0.65
1:A:181:GLN:CG	1:C:165:ARG:HH22	2.02	0.65
1:A:183:ARG:C	1:A:185:ARG:H	2.04	0.65
1:D:112:LEU:HD22	2:D:258:HOH:O	1.95	0.65
1:C:149:VAL:O	1:C:150:SER:HB3	1.97	0.65
1:A:167:LEU:HB3	1:C:178:GLN:HG2	1.78	0.64
1:D:10:LEU:HB3	1:D:14:ARG:NH2	2.11	0.64
1:C:49:LEU:HB3	1:C:55:LEU:HB2	1.77	0.64
1:B:171:GLN:NE2	2:B:250:HOH:O	2.30	0.64
1:A:30:PHE:O	1:A:32:CYS:N	2.30	0.64
1:C:66:VAL:C	1:C:68:GLN:H	2.03	0.64
1:A:9:SER:HB2	1:A:12:PHE:CD2	2.29	0.64
1:A:168:ASP:CA	1:C:178:GLN:OE1	2.46	0.64
1:A:116:ARG:NH1	1:A:134:SER:O	2.30	0.63
1:D:101:VAL:C	1:D:103:VAL:H	2.06	0.63
1:B:30:PHE:O	1:B:32:CYS:N	2.32	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:66:VAL:C	1:A:68:GLN:H	2.05	0.63
1:A:70:MET:HE3	1:A:73:LEU:HD12	1.81	0.63
1:A:101:VAL:C	1:A:103:VAL:H	2.07	0.63
1:B:159:ASP:OD1	1:C:165:ARG:NH2	2.31	0.63
1:D:81:LYS:O	1:D:81:LYS:HD3	1.99	0.63
1:B:27:LEU:HD23	1:B:67:LEU:HD22	1.81	0.63
1:D:167:LEU:H	1:D:167:LEU:CD2	2.12	0.63
1:A:167:LEU:H	1:A:167:LEU:CD2	2.09	0.62
1:B:162:ARG:HD3	1:C:165:ARG:HB2	1.79	0.62
1:D:9:SER:HB2	1:D:12:PHE:CD2	2.30	0.62
1:A:57:LEU:O	1:A:61:VAL:HG23	2.00	0.62
1:B:24:GLU:HB3	1:B:67:LEU:HD13	1.80	0.62
1:D:32:CYS:HA	2:D:260:HOH:O	2.00	0.62
1:A:183:ARG:O	1:A:185:ARG:N	2.29	0.62
1:B:70:MET:HE3	1:B:73:LEU:HD12	1.81	0.62
1:D:183:ARG:C	1:D:185:ARG:H	2.07	0.62
1:D:94:THR:OG1	1:D:97:ARG:HG2	1.99	0.62
1:B:167:LEU:H	1:B:167:LEU:CD2	2.12	0.62
1:A:121:ASN:C	1:A:123:ASN:H	2.08	0.62
1:C:185:ARG:HG3	1:C:185:ARG:HH11	1.65	0.61
1:D:185:ARG:HG3	1:D:185:ARG:HH11	1.65	0.61
1:C:121:ASN:C	1:C:123:ASN:H	2.07	0.61
1:D:27:LEU:HD23	1:D:67:LEU:HD22	1.81	0.61
1:C:160:MET:O	1:C:163:THR:HB	2.01	0.61
1:A:183:ARG:CG	2:A:264:HOH:O	2.46	0.61
1:A:185:ARG:HD2	1:B:156:VAL:HA	1.83	0.61
1:A:178:GLN:HG2	1:C:167:LEU:CB	2.30	0.61
1:C:185:ARG:HD2	1:D:156:VAL:CA	2.28	0.61
1:B:57:LEU:O	1:B:61:VAL:HG23	2.00	0.60
1:C:185:ARG:CD	1:D:156:VAL:HG13	2.31	0.60
1:A:157:LEU:O	1:A:160:MET:HB3	2.01	0.60
1:A:185:ARG:CZ	1:B:156:VAL:HG22	2.31	0.60
1:B:149:VAL:O	1:B:150:SER:HB3	2.01	0.60
1:B:94:THR:OG1	1:B:97:ARG:HG2	2.02	0.60
1:B:185:ARG:HG3	1:B:185:ARG:HH11	1.66	0.60
1:D:149:VAL:O	1:D:150:SER:HB3	2.01	0.60
1:B:121:ASN:C	1:B:123:ASN:H	2.08	0.60
1:A:116:ARG:HG2	1:A:132:GLU:OE2	2.00	0.60
1:C:9:SER:HB2	1:C:12:PHE:CD2	2.31	0.60
1:D:121:ASN:C	1:D:123:ASN:H	2.09	0.60
1:A:178:GLN:OE1	1:C:168:ASP:CA	2.49	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:152:SER:HB3	1:D:153:SER:OG	2.02	0.60
1:C:157:LEU:O	1:C:160:MET:HB3	2.01	0.60
1:C:24:GLU:HB3	1:C:67:LEU:HD13	1.84	0.60
1:B:64:LEU:HD13	1:B:73:LEU:HD23	1.83	0.59
1:C:101:VAL:C	1:C:103:VAL:H	2.09	0.59
1:D:30:PHE:O	1:D:32:CYS:N	2.35	0.59
1:D:57:LEU:O	1:D:61:VAL:HG23	2.01	0.59
1:C:138:GLU:HG3	2:C:257:HOH:O	2.00	0.59
1:D:64:LEU:HD13	1:D:73:LEU:HD23	1.83	0.59
1:A:64:LEU:HD13	1:A:73:LEU:HD23	1.84	0.59
1:A:165:ARG:NH1	1:C:181:GLN:CD	2.54	0.59
1:B:9:SER:HB2	1:B:12:PHE:CD2	2.31	0.59
1:C:94:THR:OG1	1:C:97:ARG:HG2	2.03	0.59
1:C:178:GLN:NE2	2:C:262:HOH:O	2.35	0.59
1:C:189:ALA:HB1	1:D:118:PHE:CE1	2.37	0.59
1:A:160:MET:O	1:A:163:THR:HB	2.03	0.59
1:C:167:LEU:H	1:C:167:LEU:CD2	2.11	0.59
1:A:149:VAL:O	1:A:150:SER:HB3	2.01	0.59
1:A:94:THR:OG1	1:A:97:ARG:HG2	2.02	0.59
1:B:101:VAL:C	1:B:103:VAL:H	2.08	0.59
1:B:103:VAL:C	1:B:105:GLU:N	2.59	0.59
1:C:63:MET:O	1:C:67:LEU:HG	2.02	0.59
1:A:27:LEU:HD23	1:A:67:LEU:HD22	1.85	0.59
1:A:165:ARG:NH1	1:C:181:GLN:OE1	2.35	0.59
1:C:28:LEU:CD1	1:C:63:MET:HB3	2.33	0.59
1:D:181:GLN:HG3	1:D:185:ARG:HH12	1.68	0.59
1:C:64:LEU:HD13	1:C:73:LEU:HD23	1.85	0.58
1:A:89:GLY:C	1:A:91:SER:H	2.11	0.58
1:B:49:LEU:CB	1:B:55:LEU:HB2	2.33	0.58
1:B:185:ARG:O	1:B:189:ALA:HB2	2.03	0.58
1:C:139:LEU:HD23	1:C:139:LEU:O	2.03	0.58
1:A:185:ARG:HG3	1:A:185:ARG:HH11	1.68	0.58
1:B:59:ALA:HA	1:B:91:SER:HB3	1.85	0.58
1:B:183:ARG:CG	1:B:184:TYR:H	2.10	0.58
1:A:30:PHE:C	1:A:32:CYS:N	2.62	0.58
1:B:28:LEU:CD1	1:B:63:MET:HB3	2.33	0.58
1:D:183:ARG:O	1:D:185:ARG:N	2.34	0.58
1:A:81:LYS:O	1:A:81:LYS:HD3	2.04	0.58
1:A:99:LEU:O	1:A:100:MET:C	2.46	0.58
1:C:189:ALA:CB	1:D:118:PHE:CZ	2.87	0.58
1:A:168:ASP:N	1:C:178:GLN:OE1	2.37	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:184:TYR:O	1:A:185:ARG:HD3	2.04	0.58
1:C:30:PHE:O	1:C:32:CYS:N	2.36	0.58
1:A:58:ALA:HB3	2:A:261:HOH:O	2.03	0.57
1:D:28:LEU:CD1	1:D:63:MET:HB3	2.34	0.57
1:A:162:ARG:HB2	1:A:170:CYS:SG	2.44	0.57
2:C:253:HOH:O	1:D:153:SER:HA	2.03	0.57
1:D:59:ALA:HA	1:D:91:SER:HB3	1.85	0.57
1:C:70:MET:HE3	1:C:73:LEU:HD12	1.86	0.57
1:C:183:ARG:CG	1:C:184:TYR:H	2.10	0.57
1:A:183:ARG:CA	2:A:252:HOH:O	2.52	0.57
1:A:184:TYR:N	2:A:264:HOH:O	2.36	0.57
1:C:188:TYR:HB3	1:D:122:LEU:CB	2.35	0.57
1:C:116:ARG:HG2	1:C:132:GLU:OE2	2.04	0.57
1:D:183:ARG:CG	1:D:184:TYR:H	2.12	0.57
1:C:27:LEU:HD23	1:C:67:LEU:HD22	1.87	0.57
1:C:49:LEU:CB	1:C:55:LEU:HB2	2.35	0.57
1:D:157:LEU:O	1:D:160:MET:HB3	2.04	0.57
1:C:59:ALA:HA	1:C:91:SER:HB3	1.86	0.57
1:B:63:MET:O	1:B:67:LEU:HG	2.04	0.56
1:A:30:PHE:C	1:A:32:CYS:H	2.13	0.56
1:A:28:LEU:CD1	1:A:63:MET:HB3	2.34	0.56
1:C:30:PHE:HD1	1:C:141:LEU:HD11	1.71	0.56
1:B:30:PHE:C	1:B:32:CYS:H	2.12	0.56
1:C:103:VAL:C	1:C:105:GLU:N	2.62	0.56
1:D:30:PHE:HD1	1:D:141:LEU:HD11	1.70	0.56
1:A:165:ARG:HH22	1:C:181:GLN:CG	2.09	0.56
1:A:181:GLN:NE2	2:A:256:HOH:O	2.37	0.56
1:B:30:PHE:C	1:B:32:CYS:N	2.62	0.56
1:C:185:ARG:O	1:C:189:ALA:HB2	2.06	0.56
1:D:70:MET:HE3	1:D:73:LEU:HD12	1.86	0.56
1:B:54:LYS:O	1:B:56:THR:HG23	2.06	0.56
1:B:119:ALA:O	1:B:122:LEU:HD23	2.06	0.56
1:C:116:ARG:CZ	1:C:134:SER:O	2.53	0.56
1:D:49:LEU:CB	1:D:55:LEU:HB2	2.35	0.56
1:D:139:LEU:O	1:D:139:LEU:HD23	2.06	0.56
1:D:52:GLN:HA	2:D:252:HOH:O	2.06	0.56
1:D:185:ARG:O	1:D:189:ALA:HB2	2.06	0.56
1:A:49:LEU:CB	1:A:55:LEU:HB2	2.36	0.56
1:A:54:LYS:O	1:A:56:THR:N	2.39	0.56
1:B:144:GLU:HG3	1:B:149:VAL:HG23	1.88	0.56
1:C:113:ARG:HG2	1:C:113:ARG:HH11	1.71	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:170:CYS:O	1:B:174:VAL:HG23	2.06	0.56
1:C:57:LEU:O	1:C:61:VAL:HG23	2.06	0.56
1:B:116:ARG:HG2	1:B:132:GLU:OE2	2.05	0.55
1:D:116:ARG:HG2	1:D:132:GLU:OE2	2.06	0.55
2:A:250:HOH:O	1:B:118:PHE:HZ	1.83	0.55
1:D:66:VAL:C	1:D:68:GLN:N	2.64	0.55
1:D:103:VAL:C	1:D:105:GLU:N	2.62	0.55
1:C:16:LEU:HA	1:C:76:ARG:HH12	1.71	0.55
1:C:30:PHE:C	1:C:32:CYS:N	2.64	0.55
1:A:167:LEU:CB	1:C:178:GLN:CG	2.79	0.55
1:C:185:ARG:HD2	1:D:156:VAL:HG13	1.89	0.55
1:B:181:GLN:HG3	1:B:185:ARG:HH12	1.72	0.55
1:C:162:ARG:HB2	1:C:170:CYS:SG	2.47	0.55
1:D:63:MET:O	1:D:67:LEU:HG	2.06	0.55
1:A:185:ARG:O	1:A:189:ALA:HB2	2.07	0.55
1:B:139:LEU:O	1:B:139:LEU:HD23	2.06	0.55
1:C:166:ARG:O	1:C:167:LEU:C	2.49	0.55
1:B:113:ARG:HH11	1:B:113:ARG:HG2	1.72	0.55
1:D:39:CYS:HB3	1:D:44:GLN:HB3	1.88	0.55
1:A:30:PHE:CD1	1:A:137:VAL:HG12	2.42	0.54
1:B:99:LEU:O	1:B:100:MET:C	2.50	0.54
1:C:30:PHE:CE1	1:C:141:LEU:HD21	2.42	0.54
1:D:54:LYS:O	1:D:56:THR:N	2.39	0.54
1:A:161:LEU:C	1:A:163:THR:N	2.64	0.54
1:B:161:LEU:C	1:B:163:THR:N	2.65	0.54
1:D:30:PHE:CD1	1:D:137:VAL:HG12	2.42	0.54
1:A:155:SER:O	1:A:156:VAL:C	2.50	0.54
1:D:41:THR:O	1:D:44:GLN:HB2	2.07	0.54
1:B:166:ARG:O	1:B:167:LEU:C	2.51	0.54
1:C:89:GLY:C	1:C:91:SER:H	2.16	0.54
1:C:99:LEU:O	1:C:100:MET:C	2.51	0.54
1:D:10:LEU:HB2	1:D:11:PRO:HD3	1.89	0.54
1:B:54:LYS:O	1:B:56:THR:N	2.36	0.54
1:A:103:VAL:C	1:A:105:GLU:N	2.64	0.54
1:D:30:PHE:C	1:D:32:CYS:N	2.64	0.54
1:D:30:PHE:CE1	1:D:141:LEU:HD21	2.43	0.54
1:D:113:ARG:HG2	1:D:113:ARG:HH11	1.73	0.54
1:A:54:LYS:O	1:A:56:THR:HG23	2.07	0.54
1:A:57:LEU:C	1:A:57:LEU:HD13	2.33	0.53
1:A:152:SER:HB3	1:B:153:SER:OG	2.08	0.53
1:B:157:LEU:O	1:B:160:MET:HB3	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:122:LEU:HD23	1:B:122:LEU:H	1.73	0.53
1:C:40:THR:O	1:C:41:THR:C	2.50	0.53
1:C:66:VAL:C	1:C:68:GLN:N	2.66	0.53
1:D:187:CYS:C	1:D:189:ALA:N	2.65	0.53
1:B:39:CYS:HB3	1:B:44:GLN:HB3	1.90	0.53
1:C:158:ALA:C	1:C:160:MET:H	2.17	0.53
1:D:30:PHE:C	1:D:32:CYS:H	2.16	0.53
1:C:97:ARG:HD3	2:C:251:HOH:O	2.08	0.53
1:D:57:LEU:HD13	1:D:57:LEU:C	2.34	0.53
1:D:150:SER:HB2	1:D:151:PRO:HD2	1.90	0.53
1:B:41:THR:O	1:B:44:GLN:HB2	2.09	0.53
1:C:27:LEU:HD12	1:C:137:VAL:CG2	2.33	0.53
1:C:80:SER:HB3	1:C:82:GLU:HG3	1.90	0.53
1:D:166:ARG:O	1:D:167:LEU:C	2.51	0.53
1:A:59:ALA:HA	1:A:91:SER:HB3	1.91	0.53
1:A:63:MET:O	1:A:67:LEU:HG	2.08	0.53
1:A:144:GLU:HG3	1:A:149:VAL:HG23	1.89	0.53
1:A:166:ARG:O	1:A:167:LEU:C	2.52	0.53
1:C:186:TYR:CZ	1:D:159:ASP:HB2	2.44	0.53
1:A:10:LEU:HB2	1:A:11:PRO:HD3	1.90	0.53
1:C:30:PHE:C	1:C:32:CYS:H	2.15	0.53
1:C:155:SER:O	1:C:156:VAL:C	2.52	0.53
1:C:184:TYR:O	1:C:185:ARG:HD3	2.09	0.53
1:A:66:VAL:C	1:A:68:GLN:N	2.66	0.53
1:B:30:PHE:CD1	1:B:137:VAL:HG12	2.43	0.53
1:B:103:VAL:O	1:B:105:GLU:N	2.41	0.53
1:D:89:GLY:C	1:D:91:SER:H	2.17	0.53
1:A:39:CYS:HB3	1:A:44:GLN:HB3	1.90	0.52
1:A:113:ARG:HG2	1:A:113:ARG:HH11	1.75	0.52
1:C:39:CYS:HB3	1:C:44:GLN:HB3	1.91	0.52
1:A:16:LEU:HA	1:A:76:ARG:HH12	1.73	0.52
1:C:170:CYS:O	1:C:174:VAL:HG23	2.09	0.52
1:D:99:LEU:O	1:D:100:MET:C	2.51	0.52
1:D:126:LEU:C	1:D:128:THR:H	2.17	0.52
1:B:160:MET:O	1:B:163:THR:HB	2.09	0.52
1:C:16:LEU:C	1:C:18:GLU:H	2.18	0.52
1:B:84:ALA:O	1:B:87:LEU:HD23	2.09	0.52
1:B:89:GLY:C	1:B:91:SER:H	2.17	0.52
1:D:16:LEU:C	1:D:18:GLU:H	2.17	0.52
1:B:167:LEU:HD22	1:B:167:LEU:N	2.19	0.52
1:C:126:LEU:C	1:C:128:THR:H	2.18	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:166:ARG:HG2	1:B:166:ARG:HH11	1.75	0.52
1:C:188:TYR:HB3	1:D:122:LEU:CA	2.40	0.52
1:A:181:GLN:HG3	1:A:185:ARG:HH12	1.74	0.52
1:B:66:VAL:C	1:B:68:GLN:N	2.63	0.52
1:A:167:LEU:HB3	1:C:178:GLN:CG	2.40	0.52
1:B:116:ARG:CZ	1:B:134:SER:O	2.58	0.52
1:B:155:SER:O	1:B:156:VAL:C	2.52	0.52
1:C:181:GLN:HG3	1:C:185:ARG:HH12	1.74	0.52
1:C:189:ALA:HB1	1:D:118:PHE:CZ	2.45	0.52
1:B:81:LYS:HD3	1:B:81:LYS:C	2.34	0.52
1:B:177:GLU:O	1:B:181:GLN:HB2	2.10	0.52
1:C:161:LEU:C	1:C:163:THR:N	2.64	0.52
1:D:160:MET:O	1:D:163:THR:HB	2.10	0.52
1:D:81:LYS:HD3	1:D:81:LYS:C	2.35	0.51
1:D:122:LEU:HD11	1:D:148:LEU:HD13	1.92	0.51
1:D:16:LEU:HA	1:D:76:ARG:HH12	1.74	0.51
1:A:16:LEU:C	1:A:18:GLU:H	2.17	0.51
1:A:65:TYR:CE2	1:A:88:LEU:HD11	2.46	0.51
1:A:170:CYS:O	1:A:174:VAL:HG23	2.10	0.51
1:B:16:LEU:HA	1:B:76:ARG:HH12	1.74	0.51
1:D:13:LEU:C	1:D:13:LEU:HD23	2.35	0.51
1:C:144:GLU:HG3	1:C:149:VAL:HG23	1.92	0.51
1:D:161:LEU:C	1:D:163:THR:N	2.66	0.51
1:A:186:TYR:CE2	1:B:159:ASP:HB2	2.46	0.51
1:B:30:PHE:CE1	1:B:141:LEU:HD21	2.46	0.51
1:B:150:SER:HB2	1:B:151:PRO:HD2	1.93	0.51
1:D:84:ALA:O	1:D:87:LEU:HD23	2.11	0.51
1:A:80:SER:HB3	1:A:82:GLU:HG3	1.93	0.51
1:A:167:LEU:HD22	1:A:167:LEU:N	2.18	0.51
1:A:184:TYR:C	1:A:185:ARG:HG2	2.36	0.51
1:C:30:PHE:CD1	1:C:137:VAL:HG12	2.46	0.51
1:D:54:LYS:O	1:D:56:THR:HG23	2.11	0.51
1:B:10:LEU:HB2	1:B:11:PRO:HD3	1.92	0.51
1:C:41:THR:O	1:C:44:GLN:HB2	2.11	0.51
1:C:81:LYS:HD3	1:C:81:LYS:C	2.35	0.51
1:D:80:SER:HB3	1:D:82:GLU:HG3	1.92	0.51
1:D:177:GLU:HG3	2:D:256:HOH:O	2.11	0.51
1:B:16:LEU:C	1:B:18:GLU:H	2.19	0.51
1:B:40:THR:O	1:B:41:THR:C	2.54	0.51
1:B:186:TYR:CE2	1:D:178:GLN:NE2	2.79	0.51
1:C:150:SER:HB2	1:C:151:PRO:HD2	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:57:LEU:HD13	1:B:57:LEU:C	2.36	0.51
1:B:116:ARG:HD3	1:B:139:LEU:HD12	1.93	0.51
1:D:101:VAL:C	1:D:103:VAL:N	2.68	0.51
1:A:126:LEU:C	1:A:128:THR:H	2.18	0.50
1:B:122:LEU:HD23	1:B:122:LEU:N	2.25	0.50
1:A:13:LEU:HD23	1:A:13:LEU:C	2.36	0.50
1:B:126:LEU:C	1:B:128:THR:H	2.17	0.50
1:D:40:THR:O	1:D:41:THR:C	2.53	0.50
1:A:121:ASN:C	1:A:123:ASN:N	2.70	0.50
1:C:121:ASN:C	1:C:123:ASN:N	2.69	0.50
1:D:184:TYR:O	1:D:185:ARG:HD3	2.11	0.50
1:A:84:ALA:O	1:A:87:LEU:HD23	2.11	0.50
1:B:80:SER:HB3	1:B:82:GLU:HG3	1.93	0.50
1:D:170:CYS:O	1:D:174:VAL:HG23	2.12	0.50
1:C:103:VAL:O	1:C:105:GLU:N	2.45	0.50
1:D:162:ARG:HB2	1:D:170:CYS:SG	2.51	0.50
1:A:101:VAL:C	1:A:103:VAL:N	2.69	0.50
1:A:185:ARG:NE	1:B:156:VAL:HG22	2.26	0.50
1:B:184:TYR:O	1:B:185:ARG:HD3	2.12	0.50
1:A:81:LYS:HD3	1:A:81:LYS:C	2.37	0.50
1:A:187:CYS:C	1:A:189:ALA:N	2.63	0.50
1:C:164:LEU:HD22	1:C:164:LEU:N	2.26	0.50
1:C:187:CYS:C	1:C:189:ALA:N	2.63	0.50
1:A:40:THR:O	1:A:41:THR:C	2.54	0.49
1:B:13:LEU:C	1:B:13:LEU:HD23	2.37	0.49
1:A:185:ARG:CD	1:B:156:VAL:HG13	2.42	0.49
1:B:30:PHE:HD1	1:B:141:LEU:HD11	1.77	0.49
1:D:144:GLU:HG3	1:D:149:VAL:HG23	1.94	0.49
1:D:103:VAL:O	1:D:105:GLU:N	2.46	0.49
1:D:116:ARG:CZ	1:D:134:SER:O	2.60	0.49
1:B:101:VAL:C	1:B:103:VAL:N	2.70	0.49
1:C:54:LYS:O	1:C:56:THR:N	2.41	0.49
1:C:70:MET:SD	1:C:70:MET:N	2.84	0.49
1:D:69:ARG:HD3	1:D:72:LEU:HD12	1.94	0.49
1:A:70:MET:SD	1:A:70:MET:N	2.86	0.49
1:D:65:TYR:CE2	1:D:88:LEU:HD11	2.47	0.49
1:D:116:ARG:HD3	1:D:139:LEU:HD12	1.95	0.49
1:B:30:PHE:CG	1:B:137:VAL:HG12	2.47	0.49
1:B:121:ASN:C	1:B:123:ASN:N	2.69	0.49
1:C:139:LEU:HD23	1:C:139:LEU:C	2.38	0.49
1:A:165:ARG:HB2	1:D:162:ARG:HD3	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:160:MET:O	1:B:164:LEU:HD23	2.12	0.49
1:C:57:LEU:C	1:C:57:LEU:HD13	2.37	0.49
1:C:10:LEU:HB2	1:C:11:PRO:HD3	1.95	0.49
1:C:185:ARG:CD	1:D:156:VAL:HA	2.32	0.49
1:C:54:LYS:O	1:C:56:THR:HG23	2.13	0.48
1:C:84:ALA:O	1:C:87:LEU:HD23	2.12	0.48
1:D:101:VAL:O	1:D:103:VAL:N	2.46	0.48
1:D:155:SER:O	1:D:156:VAL:C	2.55	0.48
1:A:167:LEU:O	1:A:170:CYS:HB3	2.14	0.48
1:A:41:THR:O	1:A:44:GLN:HB2	2.12	0.48
1:B:155:SER:O	1:B:158:ALA:N	2.46	0.48
1:C:183:ARG:C	1:C:185:ARG:N	2.69	0.48
1:A:183:ARG:CG	1:A:184:TYR:H	2.11	0.48
1:C:179:GLN:O	1:C:183:ARG:HG2	2.14	0.48
1:C:182:ALA:O	1:C:183:ARG:C	2.57	0.48
1:D:30:PHE:CG	1:D:137:VAL:CG1	2.96	0.48
1:A:27:LEU:HD12	1:A:137:VAL:CG2	2.39	0.48
1:A:69:ARG:HD3	1:A:72:LEU:HD12	1.95	0.48
1:B:139:LEU:HD23	1:B:139:LEU:C	2.38	0.48
1:C:13:LEU:HD23	1:C:13:LEU:C	2.39	0.48
1:C:65:TYR:CE2	1:C:88:LEU:HD11	2.49	0.48
1:A:62:GLU:O	1:A:65:TYR:N	2.46	0.48
1:C:167:LEU:O	1:C:170:CYS:HB3	2.13	0.48
1:B:65:TYR:CE2	1:B:88:LEU:HD11	2.48	0.48
1:D:30:PHE:CG	1:D:137:VAL:HG12	2.49	0.48
1:A:122:LEU:N	1:A:122:LEU:HD23	2.28	0.47
1:A:182:ALA:O	1:A:183:ARG:C	2.57	0.47
1:B:29:LEU:HD12	1:B:29:LEU:O	2.14	0.47
1:C:122:LEU:N	1:C:122:LEU:HD23	2.28	0.47
1:A:30:PHE:CG	1:A:137:VAL:CG1	2.97	0.47
1:A:116:ARG:HD3	1:A:139:LEU:HD12	1.96	0.47
1:A:158:ALA:C	1:A:160:MET:H	2.22	0.47
1:B:10:LEU:HD23	2:B:255:HOH:O	2.14	0.47
1:C:30:PHE:CG	1:C:137:VAL:CG1	2.98	0.47
1:B:158:ALA:C	1:B:160:MET:H	2.23	0.47
1:A:122:LEU:HD23	1:A:122:LEU:H	1.79	0.47
1:A:136:PHE:O	1:A:138:GLU:N	2.46	0.47
1:C:122:LEU:HD23	1:C:122:LEU:H	1.79	0.47
1:D:112:LEU:HG	1:D:116:ARG:NH1	2.30	0.47
1:A:30:PHE:CG	1:A:137:VAL:HG12	2.50	0.47
1:A:139:LEU:HD23	1:A:139:LEU:O	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:51:GLN:O	1:B:53:ARG:N	2.48	0.47
1:C:176:TYR:C	1:C:178:GLN:N	2.71	0.47
1:D:143:LEU:HD22	1:D:148:LEU:HD12	1.97	0.47
1:B:13:LEU:HD13	2:B:255:HOH:O	2.15	0.47
1:B:112:LEU:HG	1:B:116:ARG:NH1	2.29	0.47
1:A:181:GLN:CD	1:C:165:ARG:NH1	2.71	0.47
1:A:185:ARG:HD2	1:B:156:VAL:HG13	1.96	0.47
1:B:51:GLN:C	1:B:53:ARG:N	2.72	0.47
1:C:51:GLN:C	1:C:53:ARG:N	2.72	0.47
1:C:122:LEU:HD11	1:C:148:LEU:HD13	1.96	0.47
1:D:167:LEU:HD22	1:D:167:LEU:N	2.19	0.47
1:D:168:ASP:OD1	1:D:168:ASP:N	2.47	0.47
1:A:30:PHE:CE1	1:A:141:LEU:HD21	2.50	0.47
1:A:167:LEU:N	2:A:258:HOH:O	2.48	0.47
1:C:10:LEU:HB3	1:C:14:ARG:HH22	1.80	0.47
1:D:27:LEU:HD23	1:D:67:LEU:CD2	2.45	0.47
1:A:69:ARG:CD	1:A:72:LEU:HD12	2.44	0.47
1:A:165:ARG:NH2	1:D:159:ASP:OD1	2.47	0.47
1:D:119:ALA:O	1:D:122:LEU:HD23	2.15	0.47
1:A:9:SER:H	1:A:12:PHE:HE2	1.63	0.47
1:B:27:LEU:HD12	1:B:137:VAL:CG2	2.39	0.47
1:B:30:PHE:CG	1:B:137:VAL:CG1	2.97	0.47
1:B:101:VAL:O	1:B:105:GLU:HG2	2.15	0.47
1:B:123:ASN:ND2	1:B:125:SER:H	2.13	0.47
1:A:30:PHE:HD1	1:A:141:LEU:HD11	1.81	0.46
1:B:51:GLN:C	1:B:53:ARG:H	2.24	0.46
1:C:119:ALA:O	1:C:122:LEU:HD23	2.15	0.46
1:D:48:SER:HB2	2:D:259:HOH:O	2.14	0.46
1:A:116:ARG:CZ	1:A:134:SER:O	2.63	0.46
1:B:182:ALA:O	1:B:183:ARG:C	2.58	0.46
1:C:8:PRO:HG3	1:C:57:LEU:HA	1.97	0.46
1:D:36:ALA:HB1	1:D:39:CYS:SG	2.55	0.46
1:D:158:ALA:C	1:D:160:MET:H	2.23	0.46
1:B:122:LEU:H	1:B:122:LEU:CD2	2.29	0.46
1:B:166:ARG:HG2	1:B:166:ARG:NH1	2.29	0.46
1:A:103:VAL:O	1:A:105:GLU:N	2.49	0.46
1:C:30:PHE:CG	1:C:137:VAL:HG12	2.50	0.46
1:D:177:GLU:O	1:D:181:GLN:HB2	2.16	0.46
1:B:7:VAL:HG21	1:B:53:ARG:HH22	1.80	0.46
1:B:122:LEU:HD11	1:B:148:LEU:HD13	1.97	0.46
1:D:27:LEU:HD12	1:D:137:VAL:CG2	2.37	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:183:ARG:C	1:A:185:ARG:N	2.71	0.46
1:D:182:ALA:O	1:D:183:ARG:C	2.58	0.46
1:A:180:GLU:C	1:A:182:ALA:H	2.23	0.46
1:A:183:ARG:HA	2:A:252:HOH:O	2.14	0.46
1:C:123:ASN:ND2	1:C:125:SER:H	2.14	0.46
1:C:164:LEU:N	1:C:164:LEU:CD2	2.79	0.46
1:D:10:LEU:H	1:D:11:PRO:CD	2.28	0.46
1:D:69:ARG:CD	1:D:72:LEU:HD12	2.46	0.46
1:D:122:LEU:HD23	1:D:122:LEU:N	2.31	0.46
1:D:139:LEU:HD23	1:D:139:LEU:C	2.39	0.46
1:A:143:LEU:HD22	1:A:148:LEU:HD12	1.97	0.46
1:A:164:LEU:HD22	1:A:164:LEU:N	2.31	0.46
1:C:51:GLN:C	1:C:53:ARG:H	2.24	0.46
1:C:116:ARG:HD3	1:C:139:LEU:HD12	1.97	0.46
1:A:7:VAL:HG21	1:A:53:ARG:HH22	1.81	0.46
1:D:121:ASN:C	1:D:123:ASN:N	2.71	0.46
1:B:27:LEU:HD23	1:B:67:LEU:CD2	2.45	0.46
1:C:177:GLU:O	1:C:181:GLN:HB2	2.15	0.46
1:D:51:GLN:C	1:D:53:ARG:H	2.24	0.46
1:D:99:LEU:HD21	1:D:173:LEU:HD12	1.98	0.46
1:A:8:PRO:HG3	1:A:57:LEU:HA	1.99	0.45
1:A:119:ALA:O	1:A:122:LEU:HD23	2.16	0.45
1:A:95:ARG:HG3	1:A:95:ARG:HH11	1.82	0.45
1:C:9:SER:H	1:C:12:PHE:HE2	1.64	0.45
1:D:160:MET:O	1:D:164:LEU:HD23	2.16	0.45
1:D:166:ARG:HH11	1:D:166:ARG:HG2	1.80	0.45
1:A:51:GLN:C	1:A:53:ARG:N	2.73	0.45
1:A:77:PHE:C	1:A:79:LEU:H	2.24	0.45
1:B:162:ARG:CD	1:C:165:ARG:HB2	2.46	0.45
1:C:10:LEU:H	1:C:11:PRO:CD	2.29	0.45
1:C:62:GLU:HB2	1:C:93:LEU:HD12	1.98	0.45
1:C:112:LEU:O	1:C:112:LEU:HD12	2.16	0.45
1:A:150:SER:HB2	1:A:151:PRO:HD2	1.98	0.45
1:C:157:LEU:O	1:C:158:ALA:C	2.58	0.45
1:C:168:ASP:OD1	1:C:168:ASP:N	2.48	0.45
1:B:10:LEU:H	1:B:11:PRO:CD	2.29	0.45
1:C:176:TYR:O	1:C:178:GLN:N	2.49	0.45
1:D:62:GLU:HB2	1:D:93:LEU:HD12	1.99	0.45
1:D:122:LEU:HD23	1:D:122:LEU:H	1.81	0.45
1:A:62:GLU:HB2	1:A:93:LEU:HD12	1.99	0.45
1:A:101:VAL:O	1:A:103:VAL:N	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:166:ARG:HG2	1:A:166:ARG:HH11	1.81	0.45
1:C:160:MET:O	1:C:164:LEU:HD23	2.17	0.45
1:D:174:VAL:HA	1:D:177:GLU:HB3	1.99	0.45
1:C:20:LEU:HB3	2:C:263:HOH:O	2.17	0.45
1:A:51:GLN:O	1:A:53:ARG:N	2.49	0.45
1:B:62:GLU:O	1:B:65:TYR:N	2.50	0.45
1:A:163:THR:C	1:A:165:ARG:H	2.23	0.45
1:B:95:ARG:HG3	1:B:95:ARG:HH11	1.82	0.45
1:B:99:LEU:O	1:B:102:CYS:N	2.50	0.45
1:B:135:ARG:HG3	1:B:135:ARG:HH11	1.82	0.45
1:A:10:LEU:HB3	1:A:14:ARG:HH22	1.80	0.45
1:C:101:VAL:C	1:C:103:VAL:N	2.72	0.45
1:C:189:ALA:HA	1:D:118:PHE:CZ	2.52	0.45
1:A:10:LEU:H	1:A:11:PRO:CD	2.29	0.44
1:A:123:ASN:ND2	1:A:125:SER:H	2.15	0.44
1:B:69:ARG:HD3	1:B:72:LEU:HD12	1.99	0.44
1:B:136:PHE:O	1:B:138:GLU:N	2.50	0.44
1:B:178:GLN:NE2	1:D:186:TYR:CE2	2.85	0.44
1:D:65:TYR:OH	1:D:81:LYS:HE3	2.17	0.44
1:B:116:ARG:C	1:B:118:PHE:H	2.26	0.44
1:B:143:LEU:HD22	1:B:148:LEU:HD12	1.99	0.44
1:C:163:THR:C	1:C:165:ARG:H	2.24	0.44
1:A:16:LEU:C	1:A:18:GLU:N	2.75	0.44
1:A:51:GLN:C	1:A:53:ARG:H	2.25	0.44
1:A:168:ASP:N	1:A:168:ASP:OD1	2.49	0.44
1:A:112:LEU:HG	1:A:116:ARG:NH1	2.30	0.44
1:C:16:LEU:C	1:C:18:GLU:N	2.75	0.44
1:C:77:PHE:C	1:C:79:LEU:H	2.24	0.44
1:D:51:GLN:O	1:D:53:ARG:N	2.50	0.44
1:D:58:ALA:O	1:D:61:VAL:N	2.51	0.44
1:D:135:ARG:HG3	1:D:135:ARG:HH11	1.83	0.44
1:B:29:LEU:HD12	1:B:29:LEU:C	2.43	0.44
1:C:51:GLN:O	1:C:53:ARG:N	2.50	0.44
1:D:118:PHE:CD2	1:D:160:MET:HE2	2.52	0.44
1:A:179:GLN:O	1:A:183:ARG:HG2	2.18	0.44
1:B:9:SER:H	1:B:12:PHE:HE2	1.65	0.44
1:C:69:ARG:HD3	1:C:72:LEU:HD12	1.99	0.44
1:C:188:TYR:O	1:D:122:LEU:HB3	2.18	0.44
1:A:62:GLU:O	1:A:63:MET:C	2.61	0.44
1:B:62:GLU:HB2	1:B:93:LEU:HD12	2.00	0.44
1:D:7:VAL:HG21	1:D:53:ARG:HH22	1.83	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:29:LEU:C	1:A:29:LEU:HD12	2.43	0.44
1:A:29:LEU:HD12	1:A:29:LEU:O	2.17	0.44
1:A:185:ARG:HG3	1:A:185:ARG:NH1	2.32	0.44
1:B:65:TYR:OH	1:B:81:LYS:HE3	2.18	0.44
1:B:70:MET:N	1:B:70:MET:SD	2.91	0.44
1:B:99:LEU:HD21	1:B:173:LEU:HD12	1.99	0.44
1:C:97:ARG:CA	2:C:251:HOH:O	2.45	0.44
1:C:161:LEU:C	1:C:163:THR:H	2.25	0.44
1:D:163:THR:C	1:D:165:ARG:H	2.26	0.44
1:D:185:ARG:HG3	1:D:185:ARG:NH1	2.30	0.44
1:D:70:MET:SD	1:D:70:MET:N	2.91	0.43
1:B:77:PHE:C	1:B:79:LEU:H	2.26	0.43
1:B:164:LEU:HD22	1:B:164:LEU:N	2.33	0.43
1:C:31:LEU:HD21	1:C:98:LYS:HG2	2.00	0.43
1:D:179:GLN:O	1:D:183:ARG:HG2	2.18	0.43
1:C:180:GLU:C	1:C:182:ALA:H	2.27	0.43
1:D:95:ARG:HG3	1:D:95:ARG:HH11	1.83	0.43
1:A:176:TYR:C	1:A:178:GLN:N	2.75	0.43
1:B:8:PRO:HG3	1:B:57:LEU:HA	2.00	0.43
1:B:185:ARG:HG3	1:B:185:ARG:NH1	2.32	0.43
1:C:136:PHE:O	1:C:138:GLU:N	2.52	0.43
1:D:29:LEU:HD12	1:D:29:LEU:O	2.18	0.43
1:D:51:GLN:C	1:D:53:ARG:N	2.72	0.43
1:C:143:LEU:HD22	1:C:148:LEU:HD12	1.99	0.43
1:D:8:PRO:HG3	1:D:57:LEU:HA	2.00	0.43
1:D:77:PHE:C	1:D:79:LEU:H	2.26	0.43
1:A:167:LEU:HD23	2:A:258:HOH:O	2.17	0.43
1:C:21:ASP:OD1	1:C:23:HIS:HB3	2.18	0.43
1:A:87:LEU:HD23	1:A:87:LEU:N	2.34	0.43
1:C:121:ASN:HB3	2:D:253:HOH:O	2.19	0.43
1:D:9:SER:H	1:D:12:PHE:HE2	1.65	0.43
1:D:31:LEU:CD2	1:D:93:LEU:HD22	2.41	0.43
1:A:122:LEU:HD11	1:A:148:LEU:HD13	2.01	0.43
1:A:155:SER:O	1:A:158:ALA:N	2.52	0.43
1:C:7:VAL:HG21	1:C:53:ARG:HH22	1.83	0.43
1:C:101:VAL:O	1:C:105:GLU:HG2	2.18	0.43
1:C:113:ARG:HG2	1:C:113:ARG:NH1	2.34	0.43
1:D:62:GLU:O	1:D:65:TYR:N	2.52	0.43
1:A:160:MET:O	1:A:164:LEU:HD23	2.19	0.43
1:B:16:LEU:C	1:B:18:GLU:N	2.77	0.43
1:B:36:ALA:HB1	1:B:39:CYS:SG	2.59	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:58:ALA:O	1:A:61:VAL:N	2.52	0.43
1:A:189:ALA:HB1	1:B:118:PHE:CE1	2.54	0.43
1:B:10:LEU:HB3	1:B:14:ARG:HH22	1.82	0.43
1:B:167:LEU:O	1:B:170:CYS:HB3	2.18	0.43
1:C:98:LYS:O	1:C:102:CYS:HB2	2.19	0.43
1:C:185:ARG:HG3	1:C:185:ARG:NH1	2.30	0.43
1:B:58:ALA:CB	1:B:87:LEU:HD12	2.49	0.42
1:B:168:ASP:OD1	1:B:168:ASP:N	2.51	0.42
1:B:177:GLU:O	1:B:181:GLN:CB	2.67	0.42
1:C:95:ARG:HH11	1:C:95:ARG:HG3	1.82	0.42
1:A:112:LEU:HD12	1:A:112:LEU:O	2.19	0.42
1:B:98:LYS:O	1:B:102:CYS:HB2	2.18	0.42
1:B:118:PHE:CD2	1:B:160:MET:HE2	2.54	0.42
1:C:58:ALA:CB	1:C:87:LEU:HD12	2.49	0.42
1:A:184:TYR:O	1:A:185:ARG:CD	2.67	0.42
1:B:119:ALA:C	1:B:121:ASN:H	2.27	0.42
1:D:57:LEU:O	1:D:60:LEU:HB3	2.20	0.42
1:D:123:ASN:ND2	1:D:125:SER:H	2.16	0.42
1:A:43:THR:O	1:A:44:GLN:C	2.61	0.42
1:C:62:GLU:O	1:C:65:TYR:N	2.53	0.42
1:D:16:LEU:C	1:D:18:GLU:N	2.74	0.42
1:A:36:ALA:HB1	1:A:39:CYS:SG	2.59	0.42
1:A:58:ALA:CB	1:A:87:LEU:HD12	2.49	0.42
1:C:36:ALA:HB1	1:C:39:CYS:SG	2.59	0.42
1:C:65:TYR:OH	1:C:81:LYS:HE3	2.19	0.42
1:D:29:LEU:HD12	1:D:29:LEU:C	2.44	0.42
1:D:116:ARG:HA	1:D:139:LEU:CD1	2.50	0.42
1:D:136:PHE:O	1:D:138:GLU:N	2.53	0.42
1:B:69:ARG:CD	1:B:72:LEU:HD12	2.50	0.42
1:C:116:ARG:C	1:C:118:PHE:H	2.28	0.42
1:A:13:LEU:HD23	1:A:14:ARG:N	2.35	0.42
1:A:139:LEU:HD23	1:A:139:LEU:C	2.44	0.42
1:A:157:LEU:O	1:A:158:ALA:C	2.61	0.42
1:D:58:ALA:CB	1:D:87:LEU:HD12	2.49	0.42
1:A:113:ARG:NE	2:A:253:HOH:O	2.48	0.42
1:A:176:TYR:O	1:A:178:GLN:N	2.53	0.42
1:D:10:LEU:N	1:D:11:PRO:CD	2.83	0.42
1:A:177:GLU:O	1:A:181:GLN:HB2	2.18	0.42
1:C:101:VAL:O	1:C:103:VAL:N	2.53	0.42
1:C:166:ARG:HG2	1:C:166:ARG:HH11	1.85	0.42
1:D:43:THR:O	1:D:44:GLN:C	2.63	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:166:ARG:HG2	1:D:166:ARG:NH1	2.35	0.42
1:D:184:TYR:C	1:D:185:ARG:HG2	2.43	0.42
1:A:24:GLU:O	1:A:28:LEU:HB2	2.19	0.42
1:A:99:LEU:HD21	1:A:173:LEU:HD12	2.02	0.42
1:D:116:ARG:C	1:D:118:PHE:H	2.28	0.42
1:D:143:LEU:CD2	1:D:148:LEU:HD12	2.50	0.42
1:A:21:ASP:OD1	1:A:23:HIS:HB3	2.20	0.41
1:A:116:ARG:C	1:A:118:PHE:H	2.27	0.41
1:B:157:LEU:O	1:B:158:ALA:C	2.63	0.41
1:B:183:ARG:C	1:B:185:ARG:N	2.71	0.41
1:C:69:ARG:CD	1:C:72:LEU:HD12	2.50	0.41
1:C:135:ARG:NH1	2:C:259:HOH:O	2.52	0.41
1:C:157:LEU:O	1:C:160:MET:CB	2.67	0.41
1:C:184:TYR:C	1:C:185:ARG:HG2	2.44	0.41
1:D:130:LEU:HD21	1:D:139:LEU:HA	2.02	0.41
1:A:40:THR:O	1:A:41:THR:O	2.38	0.41
1:A:155:SER:O	1:A:157:LEU:N	2.53	0.41
1:A:161:LEU:C	1:A:163:THR:H	2.26	0.41
1:D:62:GLU:O	1:D:63:MET:C	2.63	0.41
1:A:183:ARG:CD	2:A:264:HOH:O	2.68	0.41
1:B:57:LEU:O	1:B:60:LEU:HB3	2.20	0.41
1:B:104:GLY:CA	1:B:136:PHE:CD2	3.03	0.41
1:B:112:LEU:HD12	1:B:112:LEU:O	2.20	0.41
1:B:155:SER:O	1:B:157:LEU:N	2.53	0.41
1:C:57:LEU:O	1:C:60:LEU:HB3	2.19	0.41
1:D:31:LEU:HD21	1:D:98:LYS:HG2	2.02	0.41
1:A:98:LYS:O	1:A:102:CYS:HB2	2.21	0.41
1:A:122:LEU:H	1:A:122:LEU:CD2	2.33	0.41
1:B:21:ASP:OD1	1:B:23:HIS:HB3	2.20	0.41
1:A:57:LEU:O	1:A:60:LEU:HB3	2.21	0.41
1:A:138:GLU:C	1:A:140:VAL:N	2.78	0.41
1:B:62:GLU:O	1:B:63:MET:C	2.63	0.41
1:C:130:LEU:HD21	1:C:139:LEU:HA	2.02	0.41
1:A:143:LEU:CD2	1:A:148:LEU:HD12	2.51	0.41
1:C:167:LEU:HD22	1:C:167:LEU:N	2.18	0.41
1:D:104:GLY:CA	1:D:136:PHE:CD2	3.04	0.41
1:A:27:LEU:HD21	1:A:101:VAL:HG13	2.02	0.41
1:A:41:THR:HG22	1:A:43:THR:N	2.31	0.41
1:A:116:ARG:CD	1:A:139:LEU:HD12	2.50	0.41
1:A:136:PHE:C	1:A:138:GLU:N	2.78	0.41
1:B:10:LEU:N	1:B:11:PRO:CD	2.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:116:ARG:CD	1:B:139:LEU:HD12	2.51	0.41
1:B:161:LEU:C	1:B:163:THR:H	2.27	0.41
1:C:55:LEU:O	1:C:55:LEU:HG	2.20	0.41
1:C:118:PHE:CD2	1:C:160:MET:HE2	2.56	0.41
1:C:122:LEU:H	1:C:122:LEU:CD2	2.33	0.41
1:C:176:TYR:C	1:C:178:GLN:H	2.28	0.41
1:D:116:ARG:HA	1:D:139:LEU:HD11	2.02	0.41
1:D:161:LEU:C	1:D:163:THR:H	2.29	0.41
1:A:27:LEU:HD23	1:A:67:LEU:CD2	2.50	0.41
1:A:164:LEU:N	1:A:164:LEU:CD2	2.84	0.41
1:B:101:VAL:O	1:B:103:VAL:N	2.53	0.41
1:B:184:TYR:C	1:B:185:ARG:HG2	2.45	0.41
1:C:176:TYR:CZ	1:C:180:GLU:HG3	2.56	0.41
1:A:99:LEU:O	1:A:102:CYS:N	2.55	0.40
1:A:104:GLY:CA	1:A:136:PHE:CD2	3.04	0.40
1:B:34:ASP:CG	1:B:97:ARG:HH21	2.28	0.40
1:B:186:TYR:HD1	1:D:182:ALA:CB	2.35	0.40
1:A:165:ARG:HH12	1:C:181:GLN:CG	2.31	0.40
1:B:14:ARG:O	1:B:18:GLU:HG3	2.21	0.40
1:B:163:THR:C	1:B:165:ARG:H	2.28	0.40
1:D:24:GLU:O	1:D:28:LEU:HB2	2.21	0.40
1:D:180:GLU:C	1:D:182:ALA:H	2.29	0.40
1:A:69:ARG:NH1	1:A:69:ARG:HG2	2.36	0.40
1:A:186:TYR:HE2	1:B:159:ASP:HB2	1.85	0.40
1:C:43:THR:O	1:C:44:GLN:C	2.63	0.40
1:A:116:ARG:HB3	1:A:132:GLU:OE2	2.21	0.40
1:A:118:PHE:CD2	1:A:160:MET:HE2	2.56	0.40
1:A:166:ARG:HG2	1:A:166:ARG:NH1	2.36	0.40
1:B:113:ARG:HG2	1:B:113:ARG:NH1	2.35	0.40
2:C:252:HOH:O	1:D:118:PHE:HZ	2.04	0.40
1:D:21:ASP:OD1	1:D:23:HIS:HB3	2.21	0.40
1:D:41:THR:O	1:D:42:VAL:C	2.64	0.40
1:D:101:VAL:O	1:D:105:GLU:HG2	2.22	0.40
1:A:167:LEU:HB2	1:C:178:GLN:CG	2.34	0.40
1:C:27:LEU:HD21	1:C:101:VAL:HG13	2.04	0.40
1:C:119:ALA:C	1:C:121:ASN:H	2.29	0.40
1:D:164:LEU:HD22	1:D:164:LEU:N	2.36	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:C:123:ASN:ND2	1:D:51:GLN:OE1[4_546]	2.18	0.02

### 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	188/249 (76%)	123 (65%)	46 (24%)	19 (10%)	0	7
1	B	188/249 (76%)	121 (64%)	51 (27%)	16 (8%)	0	9
1	C	188/249 (76%)	123 (65%)	45 (24%)	20 (11%)	0	6
1	D	188/249 (76%)	123 (65%)	49 (26%)	16 (8%)	0	9
All	All	752/996 (76%)	490 (65%)	191 (25%)	71 (9%)	0	8

All (71) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	21	ASP
1	A	55	LEU
1	A	184	TYR
1	B	21	ASP
1	B	55	LEU
1	B	184	TYR
1	C	21	ASP
1	C	55	LEU
1	C	184	TYR
1	D	21	ASP
1	D	55	LEU
1	D	184	TYR
1	A	31	LEU
1	A	39	CYS
1	A	41	THR
1	A	87	LEU
1	A	137	VAL

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	B	31	LEU
1	B	39	CYS
1	B	87	LEU
1	C	31	LEU
1	C	39	CYS
1	C	87	LEU
1	D	31	LEU
1	D	39	CYS
1	D	87	LEU
1	D	102	CYS
1	A	52	GLN
1	A	102	CYS
1	A	177	GLU
1	B	41	THR
1	B	52	GLN
1	B	102	CYS
1	B	137	VAL
1	C	41	THR
1	C	52	GLN
1	C	102	CYS
1	C	137	VAL
1	C	150	SER
1	C	177	GLU
1	D	52	GLN
1	D	137	VAL
1	D	150	SER
1	A	181	GLN
1	B	150	SER
1	C	158	ALA
1	C	181	GLN
1	D	41	THR
1	A	150	SER
1	A	188	TYR
1	B	132	GLU
1	C	159	ASP
1	D	132	GLU
1	A	158	ALA
1	B	104	GLY
1	C	120	CYS
1	C	132	GLU
1	D	158	ALA
1	D	104	GLY

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Mol	Chain	Res	Type
1	A	156	VAL
1	B	156	VAL
1	C	78	GLY
1	D	156	VAL
1	A	10	LEU
1	A	78	GLY
1	A	104	GLY
1	B	10	LEU
1	B	78	GLY
1	C	10	LEU
1	C	156	VAL
1	D	10	LEU

### 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	170/226 (75%)	152 (89%)	18 (11%)	6	26
1	B	170/226 (75%)	152 (89%)	18 (11%)	6	26
1	C	170/226 (75%)	152 (89%)	18 (11%)	6	26
1	D	170/226 (75%)	152 (89%)	18 (11%)	6	26
All	All	680/904 (75%)	608 (89%)	72 (11%)	6	26

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

Mol	Chain	Res	Type
1	A	12	PHE
1	A	13	LEU
1	A	25	ASP
1	A	27	LEU
1	A	29	LEU
1	A	31	LEU
1	A	44	GLN
1	A	60	LEU
1	A	82	GLU

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	A	85	GLU
1	A	94	THR
1	A	102	CYS
1	A	116	ARG
1	A	135	ARG
1	A	137	VAL
1	A	163	THR
1	A	167	LEU
1	A	168	ASP
1	B	12	PHE
1	B	13	LEU
1	B	25	ASP
1	B	27	LEU
1	B	29	LEU
1	B	31	LEU
1	B	44	GLN
1	B	60	LEU
1	B	82	GLU
1	B	85	GLU
1	B	94	THR
1	B	102	CYS
1	B	116	ARG
1	B	135	ARG
1	B	137	VAL
1	B	163	THR
1	B	167	LEU
1	B	168	ASP
1	C	12	PHE
1	C	13	LEU
1	C	25	ASP
1	C	27	LEU
1	C	29	LEU
1	C	31	LEU
1	C	44	GLN
1	C	60	LEU
1	C	82	GLU
1	C	85	GLU
1	C	94	THR
1	C	102	CYS
1	C	116	ARG
1	C	135	ARG
1	C	137	VAL

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Mol	Chain	Res	Type
1	C	163	THR
1	C	167	LEU
1	C	168	ASP
1	D	12	PHE
1	D	13	LEU
1	D	25	ASP
1	D	27	LEU
1	D	29	LEU
1	D	31	LEU
1	D	44	GLN
1	D	60	LEU
1	D	82	GLU
1	D	85	GLU
1	D	94	THR
1	D	102	CYS
1	D	116	ARG
1	D	135	ARG
1	D	137	VAL
1	D	163	THR
1	D	167	LEU
1	D	168	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	HIS
1	A	23	HIS
1	A	123	ASN
1	A	171	GLN
1	A	172	GLN
1	B	15	HIS
1	B	23	HIS
1	B	123	ASN
1	B	171	GLN
1	B	172	GLN
1	B	178	GLN
1	C	15	HIS
1	C	123	ASN
1	C	171	GLN
1	C	172	GLN
1	D	15	HIS
1	D	23	HIS

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Mol	Chain	Res	Type
1	D	51	GLN
1	D	123	ASN
1	D	171	GLN
1	D	172	GLN
1	D	178	GLN
1	D	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 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 [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	190/249 (76%)	0.15	5 (2%) 57 39	8, 43, 127, 162	0
1	B	190/249 (76%)	0.08	2 (1%) 78 56	8, 70, 144, 190	0
1	C	190/249 (76%)	0.18	1 (0%) 87 71	8, 56, 140, 196	0
1	D	190/249 (76%)	0.21	1 (0%) 87 71	8, 86, 150, 185	0
All	All	760/996 (76%)	0.16	9 (1%) 76 54	8, 62, 141, 196	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	185	ARG	3.7
1	A	184	TYR	3.3
1	D	187	CYS	2.9
1	B	187	CYS	2.8
1	A	186	TYR	2.4
1	C	83	GLY	2.3
1	B	84	ALA	2.3
1	A	95	ARG	2.0
1	A	163	THR	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

There are no ligands in this entry.

## 6.5 Other polymers

There are no such residues in this entry.