



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:09 PM GMT

PDB ID : 1NR9  
Title : Crystal Structure of Escherichia coli 1262 (APC5008), Putative Isomerase  
Authors : Kim, Y.; Joachimiak, A.; Edwards, A.; Skarina, T.; Savchenko, A.; Midwest Center for Structural Genomics (MCSG)  
Deposited on : 2003-01-24  
Resolution : 2.70 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  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

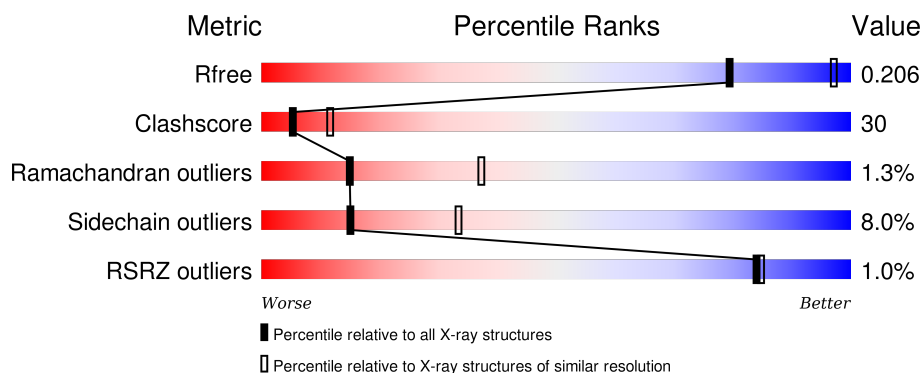
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.70 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))
$R_{free}$	91344	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (2.70-2.70)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	223	<div> <div>%</div> <div> <div></div> <div>53%</div> <div>39%</div> <div>• 5%</div> </div> </div>
1	B	223	<div> <div></div> <div> <div>48%</div> <div>42%</div> <div>5%</div> <div>6%</div> </div> </div>
1	C	223	<div> <div>%</div> <div> <div></div> <div>48%</div> <div>41%</div> <div>6%</div> <div>6%</div> </div> </div>
1	D	223	<div> <div></div> <div> <div>44%</div> <div>45%</div> <div>6%</div> <div>5%</div> </div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6697 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 Protein YCGM.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	211	Total	C	N	O	S	Se	0	0	0
			1606	1020	274	305	3	4			
1	B	210	Total	C	N	O	S	Se	0	0	0
			1599	1017	271	304	3	4			
1	C	210	Total	C	N	O	S	Se	0	0	0
			1602	1018	273	304	3	4			
1	D	211	Total	C	N	O	S	Se	0	0	0
			1609	1023	274	305	3	4			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	CLONING ARTIFACT	UNP P76004
A	0	HIS	-	CLONING ARTIFACT	UNP P76004
A	1	MSE	MET	MODIFIED RESIDUE	UNP P76004
A	33	MSE	MET	MODIFIED RESIDUE	UNP P76004
A	111	MSE	MET	MODIFIED RESIDUE	UNP P76004
A	165	MSE	MET	MODIFIED RESIDUE	UNP P76004
A	176	MSE	MET	MODIFIED RESIDUE	UNP P76004
A	220	GLY	-	CLONING ARTIFACT	UNP P76004
A	221	SER	-	CLONING ARTIFACT	UNP P76004
B	-1	GLY	-	CLONING ARTIFACT	UNP P76004
B	0	HIS	-	CLONING ARTIFACT	UNP P76004
B	1	MSE	MET	MODIFIED RESIDUE	UNP P76004
B	33	MSE	MET	MODIFIED RESIDUE	UNP P76004
B	111	MSE	MET	MODIFIED RESIDUE	UNP P76004
B	165	MSE	MET	MODIFIED RESIDUE	UNP P76004
B	176	MSE	MET	MODIFIED RESIDUE	UNP P76004
B	220	GLY	-	CLONING ARTIFACT	UNP P76004
B	221	SER	-	CLONING ARTIFACT	UNP P76004
C	-1	GLY	-	CLONING ARTIFACT	UNP P76004
C	0	HIS	-	CLONING ARTIFACT	UNP P76004
C	1	MSE	MET	MODIFIED RESIDUE	UNP P76004

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Chain	Residue	Modelled	Actual	Comment	Reference
C	33	MSE	MET	MODIFIED RESIDUE	UNP P76004
C	111	MSE	MET	MODIFIED RESIDUE	UNP P76004
C	165	MSE	MET	MODIFIED RESIDUE	UNP P76004
C	176	MSE	MET	MODIFIED RESIDUE	UNP P76004
C	220	GLY	-	CLONING ARTIFACT	UNP P76004
C	221	SER	-	CLONING ARTIFACT	UNP P76004
D	-1	GLY	-	CLONING ARTIFACT	UNP P76004
D	0	HIS	-	CLONING ARTIFACT	UNP P76004
D	1	MSE	MET	MODIFIED RESIDUE	UNP P76004
D	33	MSE	MET	MODIFIED RESIDUE	UNP P76004
D	111	MSE	MET	MODIFIED RESIDUE	UNP P76004
D	165	MSE	MET	MODIFIED RESIDUE	UNP P76004
D	176	MSE	MET	MODIFIED RESIDUE	UNP P76004
D	220	GLY	-	CLONING ARTIFACT	UNP P76004
D	221	SER	-	CLONING ARTIFACT	UNP P76004

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Mg 1 1	0	0
2	A	1	Total Mg 1 1	0	0
2	D	1	Total Mg 1 1	0	0
2	C	1	Total Mg 1 1	0	0

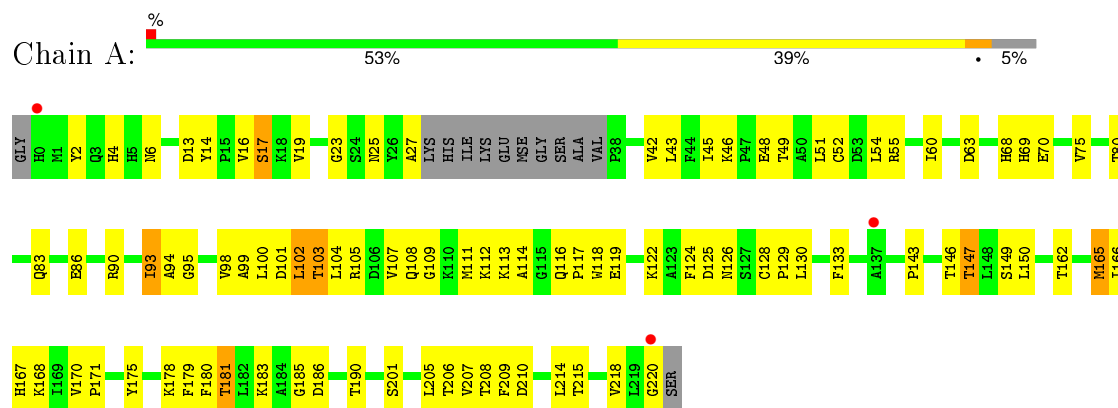
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	60	Total O 60 60	0	0
3	B	81	Total O 81 81	0	0
3	C	60	Total O 60 60	0	0
3	D	76	Total O 76 76	0	0

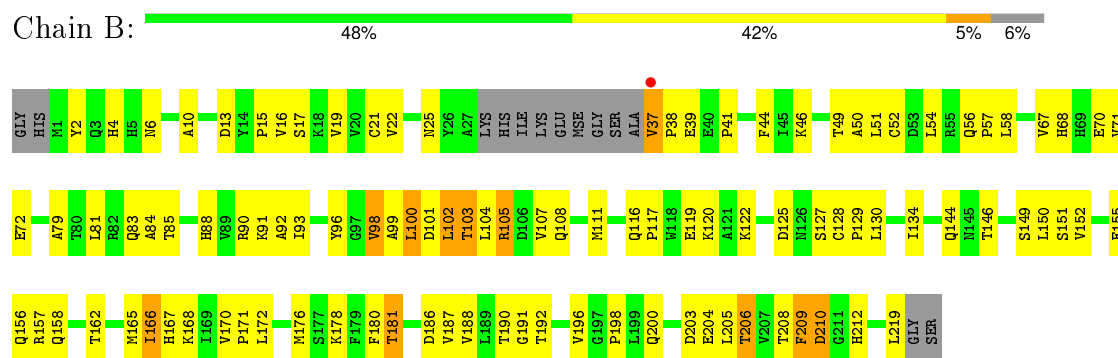
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

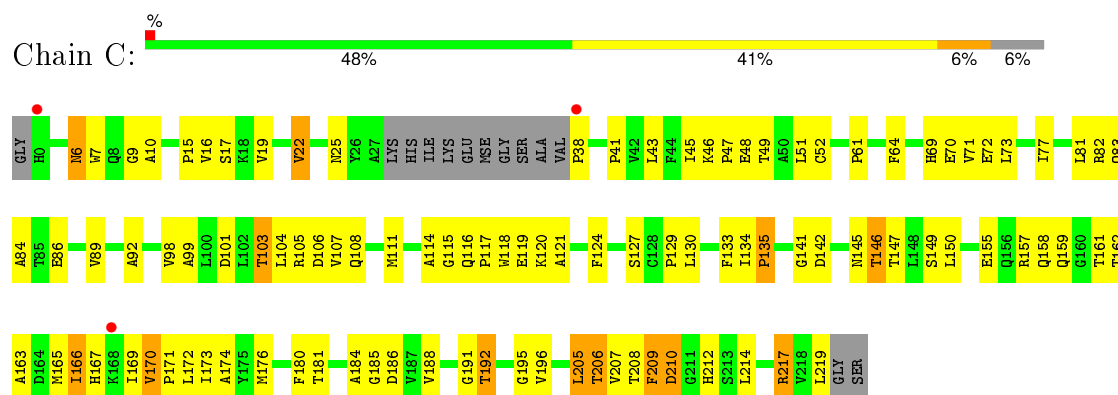
#### • Molecule 1: Protein YCGM



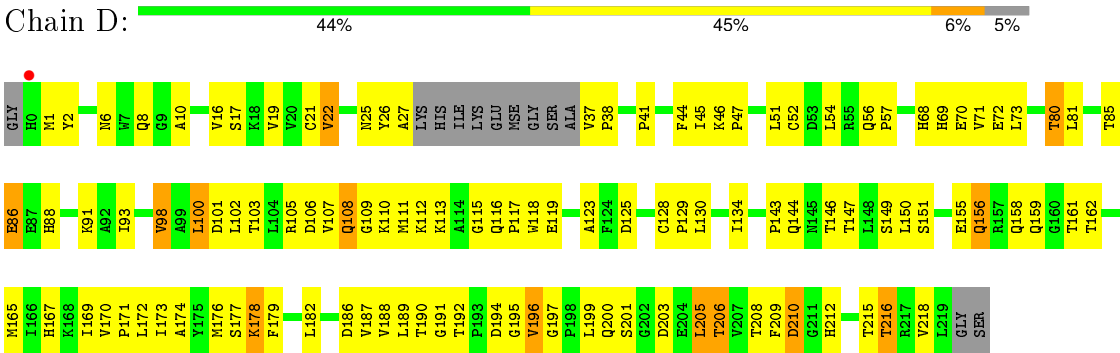
#### • Molecule 1: Protein YCGM



#### • Molecule 1: Protein YCGM



● Molecule 1: Protein YCGM



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	59.37Å 83.55Å 94.90Å 90.00° 93.08° 90.00°	Depositor
Resolution (Å)	47.38 – 2.70 47.38 – 2.67	Depositor EDS
% Data completeness (in resolution range)	97.0 (47.38-2.70) 97.9 (47.38-2.67)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.04 (at 2.65Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.206 , 0.275 0.212 , 0.206	Depositor DCC
$R_{free}$ test set	2546 reflections (10.07%)	DCC
Wilson B-factor (Å <sup>2</sup> )	36.2	Xtriage
Anisotropy	0.504	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 44.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 26030 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	6697	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.17% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG

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	0.40	0/1638	0.67	0/2220
1	B	0.39	0/1631	0.68	0/2214
1	C	0.39	0/1634	0.67	0/2215
1	D	0.39	0/1641	0.67	0/2226
All	All	0.39	0/6544	0.68	0/8875

There are no bond length outliers.

There are no bond angle outliers.

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	1606	0	1596	88	0
1	B	1599	0	1594	100	0
1	C	1602	0	1593	110	0
1	D	1609	0	1601	110	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	60	0	0	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	81	0	0	8	0
3	C	60	0	0	6	0
3	D	76	0	0	5	0
All	All	6697	0	6384	386	0

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 30.

All (386) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:146:THR:HG23	1:B:210:ASP:HB2	1.31	1.12
1:D:146:THR:HG23	1:D:210:ASP:HB2	1.29	1.10
1:C:173:ILE:HA	1:C:176:MSE:HE2	1.32	1.09
1:C:181:THR:HG23	1:D:125:ASP:OD2	1.55	1.06
1:D:173:ILE:HA	1:D:176:MSE:HE2	1.40	1.00
1:A:206:THR:HG22	1:A:215:THR:HG23	1.43	0.99
1:C:170:VAL:HG13	1:C:171:PRO:HD3	1.39	0.99
1:C:105:ARG:NH1	1:C:108:GLN:HE22	1.61	0.98
1:C:82:ARG:HD3	1:C:181:THR:HG22	1.46	0.97
1:C:161:THR:HG22	1:C:163:ALA:H	1.28	0.96
1:C:105:ARG:HH11	1:C:108:GLN:HE22	0.93	0.93
1:B:93:ILE:HD11	1:B:96:TYR:OH	1.69	0.91
1:D:146:THR:H	1:D:162:THR:HG22	1.34	0.91
1:D:73:LEU:HB3	1:D:190:THR:HG21	1.56	0.88
1:B:105:ARG:HH12	1:B:196:VAL:HG11	1.36	0.88
1:C:111:MSE:HE3	1:C:117:PRO:O	1.74	0.86
1:D:6:ASN:HD21	1:D:10:ALA:HB3	1.39	0.86
1:D:6:ASN:ND2	1:D:10:ALA:HB3	1.90	0.86
1:C:111:MSE:HE2	1:C:120:LYS:HB2	1.58	0.86
1:D:109:GLY:O	1:D:113:LYS:HD3	1.76	0.85
1:C:105:ARG:HH11	1:C:108:GLN:NE2	1.75	0.85
1:D:37:VAL:HG23	3:D:345:HOH:O	1.78	0.84
1:B:111:MSE:HE3	1:B:117:PRO:O	1.80	0.82
1:A:25:ASN:ND2	1:A:165:MSE:HG2	1.95	0.82
1:A:165:MSE:HE2	1:A:167:HIS:O	1.79	0.81
1:C:105:ARG:HH21	1:C:196:VAL:HG11	1.44	0.81
1:B:103:THR:HB	1:B:122:LYS:HD3	1.66	0.78
1:B:170:VAL:CG1	1:B:171:PRO:HD3	2.14	0.78
1:B:70:GLU:HG2	1:B:192:THR:HG21	1.65	0.77
1:C:161:THR:HG22	1:C:163:ALA:N	2.00	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:100:LEU:HD21	1:B:205:LEU:HD21	1.65	0.77
1:B:85:THR:H	1:B:88:HIS:HD2	1.32	0.77
1:C:82:ARG:HD3	1:C:181:THR:CG2	2.16	0.76
1:B:111:MSE:HE2	1:B:120:LYS:HB2	1.68	0.76
1:A:201:SER:OG	1:A:220:GLY:HA2	1.86	0.76
1:B:151:SER:HB2	1:B:206:THR:HG23	1.66	0.75
1:C:6:ASN:ND2	1:C:9:GLY:H	1.84	0.74
1:A:170:VAL:HG22	1:A:171:PRO:HD3	1.69	0.74
1:D:70:GLU:HB2	1:D:101:ASP:HB3	1.70	0.74
1:C:105:ARG:NH2	1:C:196:VAL:HG11	2.04	0.73
1:A:108:GLN:NE2	1:A:112:LYS:HE3	2.02	0.73
3:A:322:HOH:O	1:B:111:MSE:HE1	1.89	0.73
1:A:16:VAL:HG12	1:A:48:GLU:HG3	1.70	0.72
1:A:206:THR:HG22	1:A:215:THR:CG2	2.17	0.72
1:A:45:ILE:HG13	1:B:119:GLU:HG2	1.71	0.72
1:D:143:PRO:O	1:D:162:THR:HG21	1.90	0.71
1:A:93:ILE:HD12	1:A:95:GLY:H	1.55	0.71
1:D:56:GLN:O	1:D:216:THR:HG21	1.91	0.71
1:A:118:TRP:CE3	1:B:117:PRO:HG3	2.26	0.71
1:B:105:ARG:HE	1:B:108:GLN:NE2	1.88	0.70
1:A:147:THR:O	1:A:209:PHE:HA	1.92	0.70
1:B:105:ARG:NH1	1:B:196:VAL:HG11	2.05	0.70
1:A:99:ALA:HB2	1:A:130:LEU:CD1	2.22	0.69
1:D:205:LEU:HD12	1:D:205:LEU:O	1.93	0.69
1:D:146:THR:H	1:D:162:THR:CG2	2.06	0.69
1:A:68:HIS:CE1	1:A:105:ARG:HG3	2.26	0.69
1:B:105:ARG:HE	1:B:108:GLN:HE22	1.40	0.69
1:A:109:GLY:O	1:A:113:LYS:HD3	1.93	0.68
1:B:68:HIS:CE1	1:B:105:ARG:HG3	2.27	0.68
1:A:125:ASP:OD2	1:B:181:THR:HG22	1.93	0.68
1:B:22:VAL:CG2	1:B:172:LEU:HD21	2.23	0.68
1:B:107:VAL:O	1:B:111:MSE:HG2	1.94	0.68
1:B:98:VAL:HG22	1:B:209:PHE:CD2	2.29	0.67
1:B:68:HIS:ND1	1:B:105:ARG:HG3	2.10	0.67
1:B:151:SER:HB2	1:B:206:THR:CG2	2.24	0.67
1:C:101:ASP:O	1:C:103:THR:HG22	1.94	0.67
1:B:146:THR:HG23	1:B:210:ASP:CB	2.19	0.67
1:D:73:LEU:HB3	1:D:190:THR:CG2	2.25	0.67
1:B:57:PRO:HG2	3:B:372:HOH:O	1.94	0.66
1:D:107:VAL:O	1:D:111:MSE:HG2	1.94	0.66
1:D:173:ILE:CA	1:D:176:MSE:HE2	2.21	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:56:GLN:HG2	3:B:311:HOH:O	1.94	0.66
1:D:27:ALA:HB3	3:D:370:HOH:O	1.95	0.66
1:C:170:VAL:CG1	1:C:171:PRO:HD3	2.21	0.65
1:A:150:LEU:HD23	1:A:150:LEU:C	2.16	0.65
1:D:174:ALA:O	1:D:177:SER:HB3	1.96	0.65
1:C:145:ASN:HD22	1:C:161:THR:HG21	1.60	0.65
1:C:158:GLN:HG2	1:C:195:GLY:HA3	1.78	0.65
1:C:217:ARG:HH12	1:C:219:LEU:HD23	1.61	0.65
1:C:6:ASN:HD22	1:C:6:ASN:C	2.00	0.64
1:D:80:THR:HA	1:D:182:LEU:O	1.97	0.64
1:D:71:VAL:O	1:D:192:THR:HG23	1.97	0.64
1:B:134:ILE:HD11	1:B:212:HIS:CD2	2.33	0.64
1:A:68:HIS:ND1	1:A:105:ARG:HG3	2.13	0.64
1:D:173:ILE:HG12	1:D:176:MSE:HE2	1.79	0.63
1:C:70:GLU:HG2	1:C:196:VAL:HG22	1.80	0.63
1:B:168:LYS:O	1:B:171:PRO:HD2	1.98	0.63
1:C:71:VAL:O	1:C:192:THR:HG23	1.97	0.63
1:B:85:THR:H	1:B:88:HIS:CD2	2.16	0.63
1:D:22:VAL:HG23	1:D:189:LEU:O	1.99	0.62
1:B:170:VAL:HG12	1:B:171:PRO:HD3	1.81	0.62
1:D:170:VAL:HB	1:D:171:PRO:HD3	1.82	0.62
1:B:19:VAL:HA	1:B:187:VAL:O	2.00	0.62
1:B:22:VAL:HG21	1:B:172:LEU:HD21	1.82	0.62
1:B:81:LEU:HD11	1:B:92:ALA:CB	2.30	0.61
1:A:118:TRP:CD2	1:B:117:PRO:HG3	2.35	0.61
1:B:170:VAL:HG13	1:B:171:PRO:HD3	1.80	0.61
1:D:147:THR:O	1:D:209:PHE:HA	1.99	0.61
1:C:217:ARG:HG2	1:C:217:ARG:HH11	1.65	0.61
1:C:89:VAL:HB	1:C:170:VAL:HG23	1.83	0.61
1:D:128:CYS:SG	1:D:130:LEU:HD11	2.41	0.61
1:B:99:ALA:HB2	1:B:130:LEU:HD12	1.82	0.60
1:D:91:LYS:HD3	1:D:91:LYS:O	2.01	0.60
1:A:54:LEU:HD21	1:A:207:VAL:HG11	1.83	0.59
1:D:146:THR:N	1:D:162:THR:HG22	2.12	0.59
1:C:114:ALA:HB3	1:C:116:GLN:HG3	1.82	0.59
1:B:103:THR:HG21	1:B:122:LYS:HZ3	1.67	0.59
1:B:100:LEU:HB3	1:B:102:LEU:HD13	1.83	0.59
1:A:52:CYS:HB2	3:A:307:HOH:O	2.01	0.59
1:D:173:ILE:HG12	1:D:176:MSE:CE	2.32	0.59
1:C:118:TRP:CE3	1:D:117:PRO:HG3	2.37	0.59
1:D:37:VAL:HG12	1:D:38:PRO:HD3	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:49:THR:HG23	3:C:304:HOH:O	2.02	0.59
1:B:15:PRO:HG3	3:B:375:HOH:O	2.03	0.59
1:B:99:ALA:HB2	1:B:130:LEU:CD1	2.33	0.58
1:C:84:ALA:HA	3:C:339:HOH:O	2.02	0.58
1:B:149:SER:HA	1:B:158:GLN:O	2.03	0.58
1:A:98:VAL:HG13	1:A:209:PHE:CD2	2.39	0.58
1:A:99:ALA:HB2	1:A:130:LEU:HD12	1.86	0.58
1:B:91:LYS:HB3	3:B:323:HOH:O	2.03	0.58
1:D:176:MSE:HE1	1:D:188:VAL:HG11	1.85	0.58
1:D:37:VAL:N	1:D:38:PRO:HD2	2.18	0.58
1:D:206:THR:HB	1:D:215:THR:OG1	2.03	0.58
1:C:70:GLU:HA	1:C:192:THR:HG21	1.85	0.57
1:A:102:LEU:HD22	1:A:129:PRO:HD3	1.85	0.57
1:C:82:ARG:CD	1:C:181:THR:HG22	2.27	0.57
1:D:72:GLU:HG2	1:D:191:GLY:O	2.04	0.57
1:D:85:THR:H	1:D:88:HIS:CD2	2.22	0.56
1:C:6:ASN:HD21	1:C:9:GLY:H	1.52	0.56
1:D:200:GLN:HG2	1:D:203:ASP:OD2	2.05	0.56
1:A:107:VAL:O	1:A:111:MSE:HG2	2.05	0.56
1:C:206:THR:HA	1:C:214:LEU:O	2.06	0.56
1:A:93:ILE:HD12	1:A:95:GLY:N	2.21	0.56
1:C:104:LEU:HG	3:C:310:HOH:O	2.06	0.56
1:A:70:GLU:HB2	1:A:101:ASP:HB3	1.88	0.56
1:A:54:LEU:HD13	1:A:129:PRO:HB2	1.88	0.55
1:A:27:ALA:HB2	1:A:166:ILE:HD11	1.88	0.55
1:C:105:ARG:HH21	1:C:196:VAL:CG1	2.18	0.55
1:D:38:PRO:HG3	3:D:321:HOH:O	2.06	0.55
1:B:103:THR:HG21	1:B:122:LYS:NZ	2.21	0.55
1:A:100:LEU:HB3	1:A:102:LEU:HD13	1.89	0.55
1:B:90:ARG:HD2	3:B:369:HOH:O	2.07	0.55
1:B:156:GLN:OE1	1:B:156:GLN:HA	2.07	0.54
1:C:115:GLY:HA2	1:D:115:GLY:HA2	1.89	0.54
1:C:6:ASN:HD21	1:C:10:ALA:H	1.55	0.54
1:D:19:VAL:CG2	1:D:51:LEU:HD21	2.37	0.54
1:C:104:LEU:HD23	1:C:124:PHE:CE1	2.43	0.54
1:D:100:LEU:HB3	1:D:102:LEU:HD13	1.90	0.54
1:B:170:VAL:HG13	1:B:171:PRO:CD	2.38	0.54
1:C:6:ASN:ND2	1:C:6:ASN:C	2.60	0.54
1:D:162:THR:OG1	1:D:165:MSE:HE3	2.08	0.53
1:B:162:THR:HB	1:B:165:MSE:HE3	1.90	0.53
1:C:116:GLN:HB3	1:C:117:PRO:CD	2.39	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:25:ASN:ND2	1:D:165:MSE:HG2	2.23	0.53
1:D:105:ARG:NH1	1:D:108:GLN:HE22	2.07	0.53
1:D:98:VAL:HG22	1:D:209:PHE:CD2	2.44	0.53
1:D:47:PRO:HD3	1:D:123:ALA:HB1	1.90	0.53
1:C:176:MSE:HE1	1:C:188:VAL:HG21	1.89	0.53
1:C:104:LEU:HD12	3:C:333:HOH:O	2.07	0.53
1:C:52:CYS:SG	1:C:129:PRO:HA	2.48	0.53
1:D:172:LEU:O	1:D:176:MSE:HG3	2.10	0.52
1:B:71:VAL:H	1:B:158:GLN:HE22	1.57	0.52
1:A:98:VAL:CG1	1:A:209:PHE:HB3	2.39	0.52
1:D:17:SER:HB2	1:D:186:ASP:OD1	2.10	0.52
1:B:100:LEU:HD21	1:B:205:LEU:CD2	2.39	0.52
1:B:151:SER:O	1:B:205:LEU:HD12	2.09	0.52
1:C:105:ARG:HD3	1:C:108:GLN:NE2	2.25	0.52
1:C:69:HIS:HE1	1:C:157:ARG:HH11	1.56	0.52
1:A:55:ARG:HG3	1:A:214:LEU:HD13	1.91	0.52
1:B:105:ARG:HH12	1:B:196:VAL:CG1	2.15	0.52
1:C:45:ILE:HG13	1:D:119:GLU:HG2	1.92	0.52
1:A:49:THR:CG2	1:B:181:THR:HG23	2.40	0.51
1:A:17:SER:HB3	1:A:186:ASP:OD1	2.10	0.51
1:C:181:THR:CG2	1:D:125:ASP:OD2	2.43	0.51
1:C:98:VAL:HG22	1:C:209:PHE:CG	2.46	0.51
1:A:149:SER:OG	1:A:208:THR:HB	2.10	0.51
1:B:152:VAL:O	1:B:155:GLU:HG2	2.11	0.51
1:A:104:LEU:HG	1:A:124:PHE:CE1	2.46	0.51
1:A:183:LYS:HD2	3:B:332:HOH:O	2.09	0.51
1:D:69:HIS:HB3	1:D:199:LEU:HD11	1.91	0.51
1:D:86:GLU:HG3	1:D:170:VAL:HG12	1.93	0.51
1:D:85:THR:H	1:D:88:HIS:HD2	1.59	0.51
1:D:151:SER:HA	1:D:155:GLU:O	2.10	0.51
1:A:119:GLU:N	1:A:119:GLU:OE1	2.39	0.51
1:A:101:ASP:O	1:A:103:THR:HG22	2.11	0.50
1:A:103:THR:HB	1:A:122:LYS:HD3	1.93	0.50
1:D:21:CYS:HB2	1:D:44:PHE:CZ	2.46	0.50
1:B:168:LYS:C	1:B:171:PRO:HD2	2.32	0.50
1:C:117:PRO:HG3	1:D:118:TRP:CD2	2.46	0.50
1:A:114:ALA:CB	1:A:116:GLN:HG3	2.41	0.50
1:D:105:ARG:HH21	1:D:196:VAL:HG11	1.76	0.50
1:A:60:ILE:HD13	1:A:218:VAL:HG11	1.92	0.50
1:A:143:PRO:O	1:A:162:THR:HG21	2.12	0.50
1:C:145:ASN:ND2	1:C:161:THR:HG21	2.25	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:150:LEU:HD21	1:B:205:LEU:HD11	1.93	0.49
1:A:125:ASP:CG	1:B:181:THR:HG22	2.31	0.49
1:A:146:THR:HG23	1:A:210:ASP:CG	2.32	0.49
1:B:149:SER:OG	1:B:208:THR:HB	2.12	0.49
1:A:14:TYR:HB3	1:A:185:GLY:HA2	1.93	0.49
1:D:200:GLN:HA	3:D:322:HOH:O	2.11	0.49
1:A:42:VAL:HG22	1:B:116:GLN:OE1	2.12	0.49
1:D:37:VAL:N	1:D:38:PRO:CD	2.75	0.49
1:C:6:ASN:HA	1:C:133:PHE:O	2.13	0.49
1:B:72:GLU:HG2	1:B:191:GLY:O	2.13	0.49
1:C:51:LEU:HD22	1:C:130:LEU:HD22	1.94	0.49
1:A:175:TYR:O	1:A:178:LYS:HB2	2.13	0.49
1:B:93:ILE:CD1	1:B:96:TYR:OH	2.53	0.49
1:D:199:LEU:HD21	1:D:205:LEU:HD23	1.95	0.48
1:C:147:THR:O	1:C:209:PHE:HA	2.13	0.48
1:A:49:THR:HG21	1:B:181:THR:HG23	1.94	0.48
1:B:22:VAL:HB	1:B:190:THR:O	2.14	0.48
1:A:16:VAL:CG1	1:A:48:GLU:HG3	2.40	0.48
1:A:16:VAL:HG13	1:A:51:LEU:CD1	2.43	0.48
1:B:152:VAL:HG23	1:B:157:ARG:HG3	1.95	0.48
1:C:77:ILE:HG22	1:C:184:ALA:HA	1.95	0.48
1:D:69:HIS:HB3	1:D:199:LEU:CD1	2.44	0.48
1:D:1:MSE:HG2	1:D:2:TYR:N	2.27	0.48
1:C:142:ASP:O	1:C:146:THR:HB	2.14	0.48
1:C:161:THR:C	1:C:163:ALA:H	2.16	0.48
1:C:208:THR:HG23	1:C:212:HIS:O	2.14	0.48
1:C:47:PRO:HG2	1:C:49:THR:HG22	1.96	0.48
1:B:46:LYS:HD3	1:B:128:CYS:HB2	1.96	0.48
1:C:82:ARG:O	1:C:83:GLN:C	2.53	0.48
1:A:93:ILE:O	1:A:93:ILE:HG13	2.13	0.48
1:D:16:VAL:HG13	1:D:51:LEU:CD1	2.44	0.48
1:C:105:ARG:HD3	1:C:108:GLN:HE21	1.79	0.48
1:D:158:GLN:OE1	1:D:192:THR:HG22	2.14	0.48
1:C:176:MSE:HE1	1:C:188:VAL:CG2	2.42	0.47
1:C:38:PRO:HG3	1:C:166:ILE:HG21	1.96	0.47
1:D:8:GLN:NE2	3:D:312:HOH:O	2.48	0.47
1:C:70:GLU:HB3	1:C:192:THR:OG1	2.14	0.47
1:A:19:VAL:O	1:A:45:ILE:HG23	2.14	0.47
1:A:100:LEU:HB2	1:A:129:PRO:HD2	1.97	0.47
1:C:205:LEU:O	1:C:205:LEU:HD12	2.14	0.47
1:C:22:VAL:HG13	1:C:172:LEU:HD11	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:179:PHE:CD1	1:D:179:PHE:N	2.82	0.47
1:B:200:GLN:N	1:B:203:ASP:OD2	2.37	0.47
1:D:69:HIS:CE1	1:D:197:GLY:H	2.32	0.47
1:C:16:VAL:HG13	1:C:48:GLU:HG3	1.96	0.47
1:D:201:SER:HA	1:D:218:VAL:HG12	1.96	0.47
1:C:81:LEU:HD21	1:C:92:ALA:HB2	1.96	0.47
1:D:162:THR:OG1	1:D:165:MSE:CE	2.63	0.47
1:A:46:LYS:HD3	1:A:128:CYS:HB2	1.96	0.47
1:A:165:MSE:HE3	1:A:190:THR:HB	1.96	0.47
1:D:2:TYR:HB2	1:D:52:CYS:HA	1.96	0.47
1:C:19:VAL:HB	1:C:46:LYS:HB2	1.97	0.47
1:A:99:ALA:HB2	1:A:130:LEU:HD13	1.93	0.47
1:B:4:HIS:CD2	1:B:16:VAL:HG22	2.50	0.47
1:C:120:LYS:HG2	1:D:179:PHE:CD2	2.50	0.46
1:A:98:VAL:HG11	1:A:209:PHE:HB3	1.97	0.46
1:A:117:PRO:HB3	1:B:44:PHE:HB3	1.97	0.46
1:D:71:VAL:O	1:D:71:VAL:HG13	2.16	0.46
1:C:64:PHE:O	1:C:104:LEU:HD12	2.15	0.46
1:B:198:PRO:HG3	3:B:305:HOH:O	2.16	0.46
1:D:54:LEU:HD13	1:D:129:PRO:HB2	1.97	0.46
1:B:17:SER:HB3	1:B:186:ASP:OD1	2.16	0.46
1:B:84:ALA:HB1	1:B:88:HIS:HB2	1.97	0.46
1:C:117:PRO:HG3	1:D:118:TRP:CE3	2.50	0.46
1:A:168:LYS:C	1:A:171:PRO:HD2	2.36	0.46
1:C:116:GLN:HB3	1:C:117:PRO:HD2	1.98	0.46
1:B:150:LEU:C	1:B:150:LEU:HD23	2.37	0.46
1:A:93:ILE:C	1:A:93:ILE:HD12	2.36	0.45
1:B:172:LEU:HG	1:B:176:MSE:HE3	1.98	0.45
1:D:72:GLU:HG2	1:D:191:GLY:C	2.36	0.45
1:C:114:ALA:CB	1:C:116:GLN:HG3	2.46	0.45
1:C:149:SER:OG	1:C:208:THR:HB	2.16	0.45
1:A:25:ASN:HD21	1:A:165:MSE:CE	2.29	0.45
1:A:181:THR:HG23	1:B:49:THR:HG21	1.97	0.45
1:C:70:GLU:CA	1:C:192:THR:HG21	2.46	0.45
1:A:93:ILE:HD12	1:A:94:ALA:N	2.31	0.45
1:B:79:ALA:O	1:B:81:LEU:HD13	2.17	0.45
1:A:52:CYS:HB3	1:A:126:ASN:HD22	1.82	0.45
1:D:116:GLN:HB3	1:D:117:PRO:CD	2.47	0.45
1:A:75:VAL:HG13	1:A:93:ILE:HD13	1.98	0.45
1:B:67:VAL:HG22	1:B:104:LEU:HD23	1.98	0.45
1:C:209:PHE:CE2	1:C:212:HIS:HB2	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:108:GLN:O	1:D:112:LYS:HG3	2.17	0.45
1:D:144:GLN:C	1:D:162:THR:CG2	2.85	0.45
1:D:144:GLN:C	1:D:162:THR:HG23	2.37	0.45
1:D:128:CYS:SG	1:D:130:LEU:CD1	3.05	0.45
1:C:124:PHE:CD2	1:C:127:SER:HB3	2.52	0.45
1:A:83:GLN:NE2	1:A:178:LYS:O	2.50	0.44
1:C:17:SER:OG	1:C:186:ASP:OD1	2.32	0.44
1:A:90:ARG:HD2	3:A:350:HOH:O	2.17	0.44
1:D:22:VAL:HG22	1:D:172:LEU:HD21	1.98	0.44
1:B:57:PRO:HB2	1:B:219:LEU:HD21	1.99	0.44
1:D:156:GLN:NE2	1:D:159:GLN:OE1	2.43	0.44
1:C:141:GLY:HA3	3:C:352:HOH:O	2.17	0.44
1:C:149:SER:HA	1:C:158:GLN:O	2.18	0.44
1:C:41:PRO:HG3	1:C:167:HIS:CD2	2.52	0.44
1:C:72:GLU:HG2	1:C:191:GLY:O	2.18	0.44
1:D:70:GLU:CB	1:D:101:ASP:HB3	2.42	0.44
1:C:98:VAL:HG11	1:C:207:VAL:HG22	1.99	0.44
1:B:6:ASN:OD1	1:B:10:ALA:N	2.50	0.44
1:D:45:ILE:HG22	1:D:46:LYS:N	2.33	0.44
1:C:217:ARG:CG	1:C:217:ARG:HH11	2.31	0.44
1:A:86:GLU:HG3	3:A:360:HOH:O	2.18	0.44
1:D:149:SER:OG	1:D:208:THR:HB	2.18	0.44
1:B:166:ILE:HG22	1:B:167:HIS:N	2.33	0.44
1:D:113:LYS:N	1:D:113:LYS:CD	2.81	0.43
1:A:98:VAL:HG13	1:A:209:PHE:CG	2.53	0.43
1:C:41:PRO:HG3	1:C:167:HIS:CG	2.53	0.43
1:A:102:LEU:HD22	1:A:129:PRO:CD	2.48	0.43
1:D:134:ILE:HD11	1:D:212:HIS:CD2	2.53	0.43
1:A:4:HIS:NE2	1:A:185:GLY:O	2.43	0.43
1:C:107:VAL:HG12	1:C:121:ALA:HB2	1.99	0.43
1:D:16:VAL:HG13	1:D:51:LEU:HD12	2.00	0.43
1:B:83:GLN:OE1	1:B:178:LYS:HD3	2.19	0.43
1:D:56:GLN:HB3	1:D:57:PRO:HD2	2.00	0.43
1:A:170:VAL:CG2	1:A:171:PRO:HD3	2.43	0.43
1:D:19:VAL:HG21	1:D:51:LEU:HD21	2.00	0.43
1:C:7:TRP:CZ2	1:C:212:HIS:HB3	2.54	0.43
1:D:16:VAL:CG1	1:D:51:LEU:HD12	2.48	0.43
1:A:165:MSE:CE	1:A:190:THR:HB	2.48	0.43
1:B:206:THR:HG22	3:B:367:HOH:O	2.19	0.43
1:A:2:TYR:HB2	1:A:52:CYS:HA	2.00	0.43
1:D:173:ILE:HA	1:D:176:MSE:CE	2.30	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:73:LEU:HD22	1:C:162:THR:HG21	2.00	0.42
1:C:25:ASN:ND2	1:C:165:MSE:HG2	2.33	0.42
1:C:15:PRO:O	1:C:185:GLY:HA3	2.18	0.42
1:D:165:MSE:SE	1:D:190:THR:OG1	2.88	0.42
1:C:70:GLU:HB2	1:C:101:ASP:HB3	2.00	0.42
1:C:163:ALA:C	1:C:165:MSE:H	2.23	0.42
1:B:101:ASP:O	1:B:103:THR:HG22	2.19	0.42
1:D:26:TYR:O	1:D:27:ALA:C	2.57	0.42
1:D:68:HIS:CE1	1:D:105:ARG:HG3	2.53	0.42
1:A:178:LYS:HB3	1:A:178:LYS:HE2	1.83	0.42
1:A:83:GLN:NE2	1:A:178:LYS:HA	2.34	0.42
1:A:6:ASN:HA	1:A:133:PHE:O	2.19	0.42
1:C:6:ASN:HD21	1:C:9:GLY:N	2.15	0.42
1:B:98:VAL:HG13	1:B:209:PHE:HB2	2.01	0.42
1:D:68:HIS:ND1	1:D:105:ARG:HG3	2.35	0.42
1:B:52:CYS:SG	1:B:129:PRO:HA	2.59	0.42
1:D:176:MSE:HE1	1:D:188:VAL:CG1	2.50	0.42
1:B:144:GLN:CD	1:B:168:LYS:HD2	2.40	0.42
1:C:118:TRP:CD2	1:D:117:PRO:HG3	2.55	0.42
1:D:19:VAL:HA	1:D:187:VAL:O	2.20	0.42
1:C:147:THR:CG2	1:C:159:GLN:HE22	2.32	0.42
1:D:195:GLY:O	1:D:196:VAL:C	2.58	0.42
1:A:114:ALA:HB3	1:A:116:GLN:HG3	2.01	0.42
1:C:217:ARG:NH1	1:C:219:LEU:HD23	2.33	0.42
1:B:51:LEU:CD2	1:B:130:LEU:HD22	2.50	0.42
1:C:106:ASP:HB2	3:C:333:HOH:O	2.20	0.42
1:C:89:VAL:HG11	1:C:173:ILE:HB	2.03	0.41
1:C:105:ARG:NH1	1:C:108:GLN:NE2	2.45	0.41
1:A:179:PHE:CD1	1:A:179:PHE:N	2.87	0.41
1:D:21:CYS:HB3	1:D:189:LEU:HB2	2.02	0.41
1:D:41:PRO:HG3	1:D:167:HIS:CD2	2.54	0.41
1:D:106:ASP:O	1:D:110:LYS:HG3	2.20	0.41
1:B:2:TYR:OH	1:B:130:LEU:HD23	2.19	0.41
1:C:134:ILE:O	1:C:135:PRO:C	2.59	0.41
1:B:57:PRO:HB2	1:B:219:LEU:CD2	2.50	0.41
1:D:208:THR:HG22	1:D:209:PHE:N	2.36	0.41
1:A:181:THR:HG22	1:B:125:ASP:CG	2.41	0.41
1:C:119:GLU:OE2	1:D:44:PHE:HB2	2.21	0.41
1:C:209:PHE:O	1:C:209:PHE:CG	2.73	0.41
1:C:150:LEU:C	1:C:150:LEU:HD23	2.41	0.41
1:B:21:CYS:HB2	1:B:44:PHE:CZ	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:100:LEU:HB3	1:D:102:LEU:CD1	2.51	0.41
1:C:99:ALA:HB2	1:C:130:LEU:HD12	2.02	0.41
1:A:68:HIS:CG	1:A:105:ARG:CG	3.04	0.41
1:B:208:THR:HG23	1:B:212:HIS:O	2.20	0.41
1:D:98:VAL:HG13	1:D:209:PHE:HB3	2.02	0.41
1:B:38:PRO:O	1:B:39:GLU:C	2.59	0.41
1:D:178:LYS:HE3	1:D:178:LYS:HA	2.03	0.41
1:C:89:VAL:HG21	1:C:174:ALA:HB2	2.03	0.41
1:C:161:THR:C	1:C:163:ALA:N	2.74	0.41
1:B:41:PRO:HG3	1:B:167:HIS:CG	2.56	0.41
1:D:162:THR:O	1:D:165:MSE:HB2	2.22	0.40
1:A:69:HIS:CD2	1:A:150:LEU:HD13	2.57	0.40
1:C:119:GLU:N	1:C:119:GLU:OE1	2.47	0.40
1:A:25:ASN:ND2	1:A:165:MSE:CE	2.84	0.40
1:B:25:ASN:ND2	1:B:165:MSE:HG3	2.37	0.40
1:A:181:THR:HG22	1:B:125:ASP:OD2	2.21	0.40
1:C:86:GLU:HG3	1:C:170:VAL:HG22	2.02	0.40
1:C:142:ASP:HB3	1:C:145:ASN:HB3	2.04	0.40
1:B:71:VAL:H	1:B:158:GLN:NE2	2.19	0.40
1:C:61:PRO:HG2	1:C:104:LEU:HD21	2.03	0.40
1:A:116:GLN:HB3	1:A:117:PRO:CD	2.51	0.40
1:B:37:VAL:N	1:B:38:PRO:HD3	2.37	0.40
1:D:70:GLU:HB3	1:D:192:THR:OG1	2.22	0.40
1:A:25:ASN:HD21	1:A:165:MSE:HE2	1.86	0.40
1:B:100:LEU:HA	1:B:100:LEU:HD12	1.95	0.40
1:A:150:LEU:CD2	1:A:150:LEU:C	2.86	0.40
1:B:54:LEU:HD12	1:B:58:LEU:HD21	2.04	0.40
1:B:50:ALA:HB1	1:B:127:SER:OG	2.20	0.40

There are no symmetry-related clashes.

## 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	207/223 (93%)	192 (93%)	13 (6%)	2 (1%)	19	45
1	B	206/223 (92%)	191 (93%)	13 (6%)	2 (1%)	19	45
1	C	206/223 (92%)	184 (89%)	18 (9%)	4 (2%)	10	25
1	D	207/223 (93%)	190 (92%)	14 (7%)	3 (1%)	14	35
All	All	826/892 (93%)	757 (92%)	58 (7%)	11 (1%)	15	37

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	210	ASP
1	B	166	ILE
1	C	166	ILE
1	C	169	ILE
1	D	196	VAL
1	A	23	GLY
1	C	210	ASP
1	A	17	SER
1	C	155	GLU
1	D	210	ASP
1	D	169	ILE

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	174/178 (98%)	162 (93%)	12 (7%)	19	43
1	B	174/178 (98%)	161 (92%)	13 (8%)	17	38
1	C	174/178 (98%)	160 (92%)	14 (8%)	15	33
1	D	175/178 (98%)	158 (90%)	17 (10%)	10	23
All	All	697/712 (98%)	641 (92%)	56 (8%)	15	33

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	ASP
1	A	43	LEU
1	A	63	ASP
1	A	80	THR
1	A	93	ILE
1	A	102	LEU
1	A	103	THR
1	A	147	THR
1	A	165	MSE
1	A	180	PHE
1	A	181	THR
1	A	205	LEU
1	B	13	ASP
1	B	37	VAL
1	B	98	VAL
1	B	100	LEU
1	B	102	LEU
1	B	103	THR
1	B	105	ARG
1	B	180	PHE
1	B	181	THR
1	B	188	VAL
1	B	204	GLU
1	B	206	THR
1	B	209	PHE
1	C	6	ASN
1	C	22	VAL
1	C	43	LEU
1	C	103	THR
1	C	135	PRO
1	C	146	THR
1	C	170	VAL
1	C	180	PHE
1	C	192	THR
1	C	205	LEU
1	C	206	THR
1	C	209	PHE
1	C	210	ASP
1	C	217	ARG
1	D	22	VAL
1	D	80	THR
1	D	81	LEU
1	D	86	GLU

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Mol	Chain	Res	Type
1	D	93	ILE
1	D	98	VAL
1	D	100	LEU
1	D	103	THR
1	D	108	GLN
1	D	150	LEU
1	D	156	GLN
1	D	161	THR
1	D	178	LYS
1	D	194	ASP
1	D	205	LEU
1	D	206	THR
1	D	216	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	69	HIS
1	A	83	GLN
1	A	108	GLN
1	B	88	HIS
1	B	108	GLN
1	B	145	ASN
1	B	158	GLN
1	C	6	ASN
1	C	8	GLN
1	C	69	HIS
1	C	108	GLN
1	C	145	ASN
1	C	156	GLN
1	C	159	GLN
1	C	167	HIS
1	D	8	GLN
1	D	88	HIS
1	D	108	GLN
1	D	145	ASN

### 5.3.3 RNA ⓘ

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](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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	207/223 (92%)	-0.22	3 (1%) 78 77	13, 29, 53, 96	0
1	B	206/223 (92%)	-0.21	1 (0%) 91 93	14, 30, 56, 80	0
1	C	206/223 (92%)	-0.15	3 (1%) 76 76	15, 34, 60, 88	0
1	D	207/223 (92%)	-0.17	1 (0%) 91 93	16, 33, 61, 94	0
All	All	826/892 (92%)	-0.19	8 (0%) 84 85	13, 32, 59, 96	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	0	HIS	3.9
1	A	0	HIS	3.3
1	B	37	VAL	2.7
1	A	220	GLY	2.7
1	A	137	ALA	2.5
1	C	38	PRO	2.2
1	C	0	HIS	2.2
1	C	168	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	MG	D	304	1/1	0.91	0.20	1.85	24,24,24,24	0
2	MG	A	301	1/1	0.95	0.20	1.71	23,23,23,23	0
2	MG	B	302	1/1	0.94	0.21	1.24	23,23,23,23	0
2	MG	C	303	1/1	0.98	0.16	-0.05	22,22,22,22	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.