



# Full wwPDB X-ray Structure Validation Report ⓘ

Sep 29, 2024 – 05:10 PM EDT

PDB ID : 3DTO  
Title : Crystal structure of the metal-dependent HD domain-containing hydrolase BH2835 from *Bacillus halodurans*, Northeast Structural Genomics Consortium Target BhR130.  
Authors : Forouhar, F.; Su, M.; Seetharaman, J.; Janjua, H.; Fang, Y.; Xiao, R.; Cunningham, K.; Ma, L.-C.; Owen, L.A.; Wang, D.; Tong, S.; Everett, J.K.; Acton, T.B.; Montelione, G.T.; Tong, L.; Hunt, J.F.; Northeast Structural Genomics Consortium (NESG)  
Deposited on : 2008-07-15  
Resolution : 3.30 Å(reported)

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

We welcome your comments at [validation@mail.wwpdb.org](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	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)

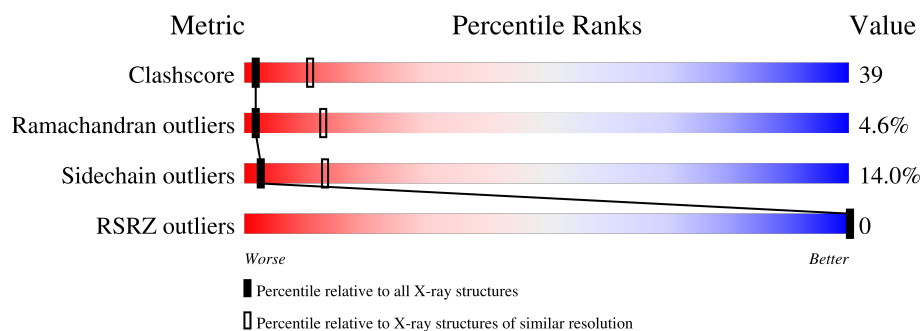
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.30 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	180529	1128 (3.32-3.28)
Ramachandran outliers	177936	1125 (3.32-3.28)
Sidechain outliers	177891	1124 (3.32-3.28)
RSRZ outliers	164620	1085 (3.32-3.28)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	223	
1	B	223	
1	C	223	
1	D	223	

Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.39

## 2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 6132 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 BH2835 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	189	Total	C	N	O	Se	0	0	0
			1533	975	263	287	8			
1	B	189	Total	C	N	O	Se	0	0	0
			1533	975	263	287	8			
1	C	189	Total	C	N	O	Se	0	0	0
			1533	975	263	287	8			
1	D	189	Total	C	N	O	Se	0	0	0
			1533	975	263	287	8			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	216	LEU	-	expression tag	UNP Q9K916
A	217	GLU	-	expression tag	UNP Q9K916
A	218	HIS	-	expression tag	UNP Q9K916
A	219	HIS	-	expression tag	UNP Q9K916
A	220	HIS	-	expression tag	UNP Q9K916
A	221	HIS	-	expression tag	UNP Q9K916
A	222	HIS	-	expression tag	UNP Q9K916
A	223	HIS	-	expression tag	UNP Q9K916
B	216	LEU	-	expression tag	UNP Q9K916
B	217	GLU	-	expression tag	UNP Q9K916
B	218	HIS	-	expression tag	UNP Q9K916
B	219	HIS	-	expression tag	UNP Q9K916
B	220	HIS	-	expression tag	UNP Q9K916
B	221	HIS	-	expression tag	UNP Q9K916
B	222	HIS	-	expression tag	UNP Q9K916
B	223	HIS	-	expression tag	UNP Q9K916
C	216	LEU	-	expression tag	UNP Q9K916
C	217	GLU	-	expression tag	UNP Q9K916
C	218	HIS	-	expression tag	UNP Q9K916
C	219	HIS	-	expression tag	UNP Q9K916
C	220	HIS	-	expression tag	UNP Q9K916

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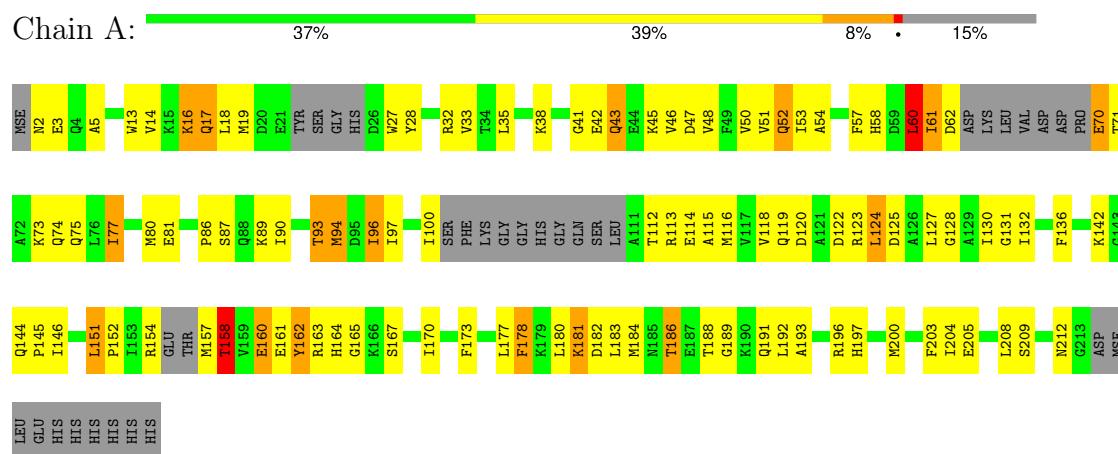
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Chain	Residue	Modelled	Actual	Comment	Reference
C	221	HIS	-	expression tag	UNP Q9K916
C	222	HIS	-	expression tag	UNP Q9K916
C	223	HIS	-	expression tag	UNP Q9K916
D	216	LEU	-	expression tag	UNP Q9K916
D	217	GLU	-	expression tag	UNP Q9K916
D	218	HIS	-	expression tag	UNP Q9K916
D	219	HIS	-	expression tag	UNP Q9K916
D	220	HIS	-	expression tag	UNP Q9K916
D	221	HIS	-	expression tag	UNP Q9K916
D	222	HIS	-	expression tag	UNP Q9K916
D	223	HIS	-	expression tag	UNP Q9K916

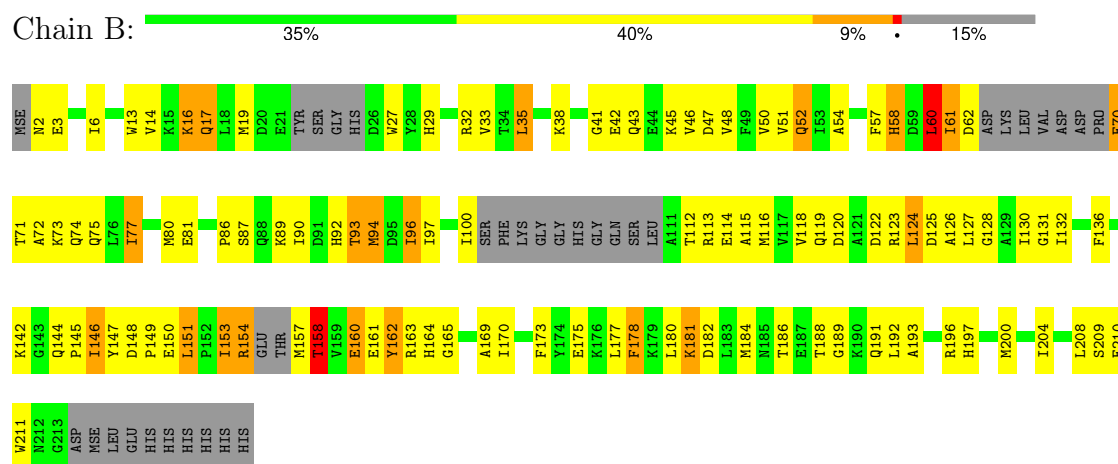
### 3 Residue-property plots

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.

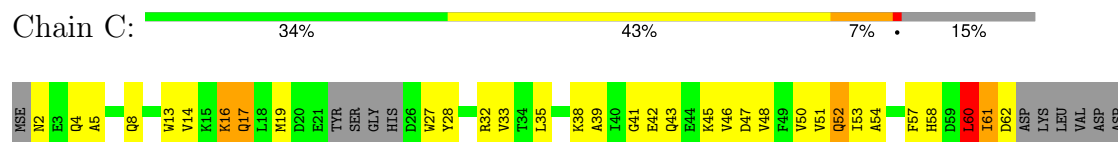
#### • Molecule 1: BH2835 protein

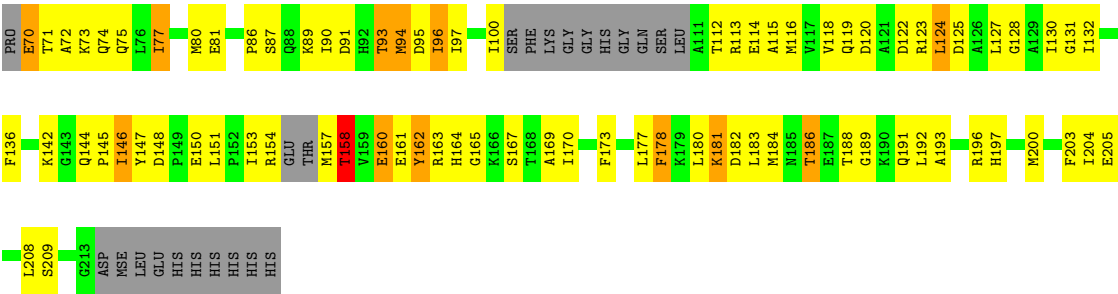


#### • Molecule 1: BH2835 protein

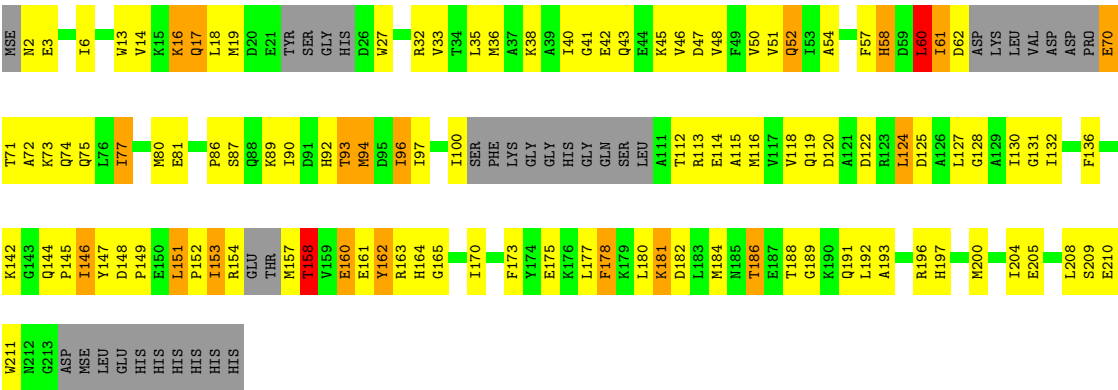


#### • Molecule 1: BH2835 protein





● Molecule 1: BH2835 protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	97.56Å 47.99Å 111.16Å 90.00° 89.99° 90.00°	Depositor
Resolution (Å)	19.76 – 3.30 19.76 – 3.32	Depositor EDS
% Data completeness (in resolution range)	76.3 (19.76-3.30) 87.4 (19.76-3.32)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.51 (at 3.31Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, $R_{free}$	0.248 , 0.292 0.283 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	71.4	Xtriage
Anisotropy	0.486	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 105.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.428 for h,-k,-l	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	6132	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.82% 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	0.51	0/1552	0.59	0/2076
1	B	0.51	0/1552	0.58	0/2076
1	C	0.51	0/1552	0.59	0/2076
1	D	0.53	0/1552	0.59	0/2076
All	All	0.52	0/6208	0.59	0/8304

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	1533	0	1512	117	0
1	B	1533	0	1512	117	0
1	C	1533	0	1512	122	0
1	D	1533	0	1512	118	0
All	All	6132	0	6048	469	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 39.

All (469) 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:D:181:LYS:HE3	1:D:181:LYS:H	1.23	1.02
1:B:181:LYS:HE3	1:B:181:LYS:H	1.22	1.02
1:C:181:LYS:HE3	1:C:181:LYS:H	1.21	1.01
1:B:186:THR:HG22	1:B:189:GLY:H	1.23	1.00
1:A:181:LYS:HE3	1:A:181:LYS:H	1.23	1.00
1:D:186:THR:HG22	1:D:189:GLY:H	1.23	0.98
1:C:186:THR:HG22	1:C:189:GLY:H	1.29	0.97
1:A:186:THR:HG22	1:A:189:GLY:H	1.28	0.96
1:B:146:ILE:HD13	1:B:146:ILE:H	1.30	0.95
1:C:113:ARG:HA	1:C:116:MSE:HE3	1.56	0.88
1:C:57:PHE:HA	1:C:60:LEU:HD21	1.56	0.88
1:D:113:ARG:HA	1:D:116:MSE:HE3	1.57	0.86
1:B:181:LYS:HE3	1:B:181:LYS:N	1.91	0.86
1:C:142:LYS:NZ	1:C:162:TYR:OH	2.10	0.85
1:A:57:PHE:HA	1:A:60:LEU:HD21	1.57	0.85
1:B:113:ARG:HA	1:B:116:MSE:HE3	1.57	0.85
1:C:181:LYS:HE3	1:C:181:LYS:N	1.90	0.85
1:D:181:LYS:HE3	1:D:181:LYS:N	1.92	0.85
1:A:181:LYS:HE3	1:A:181:LYS:N	1.92	0.84
1:D:146:ILE:HD13	1:D:146:ILE:H	1.43	0.84
1:A:113:ARG:HA	1:A:116:MSE:HE3	1.58	0.84
1:A:94:MSE:HA	1:A:97:ILE:CG1	2.09	0.83
1:C:94:MSE:HA	1:C:97:ILE:CG1	2.09	0.83
1:D:57:PHE:HA	1:D:60:LEU:HD21	1.61	0.81
1:B:57:PHE:HA	1:B:60:LEU:HD21	1.60	0.81
1:D:94:MSE:HA	1:D:97:ILE:CG1	2.10	0.81
1:C:181:LYS:H	1:C:181:LYS:CE	1.95	0.80
1:B:94:MSE:HA	1:B:97:ILE:CG1	2.11	0.80
1:D:181:LYS:H	1:D:181:LYS:CE	1.96	0.79
1:B:181:LYS:H	1:B:181:LYS:CE	1.95	0.78
1:C:94:MSE:HA	1:C:97:ILE:HG12	1.66	0.78
1:A:181:LYS:H	1:A:181:LYS:CE	1.97	0.77
1:A:154:ARG:NH1	1:A:167:SER:HB2	1.98	0.77
1:A:94:MSE:HA	1:A:97:ILE:HG12	1.65	0.76
1:B:80:MSE:HE3	1:B:90:ILE:HG13	1.68	0.75
1:D:94:MSE:HA	1:D:97:ILE:HG12	1.67	0.75
1:B:186:THR:HG22	1:B:189:GLY:N	2.00	0.74
1:B:94:MSE:HA	1:B:97:ILE:HG12	1.68	0.73
1:C:80:MSE:HE3	1:C:90:ILE:HG13	1.70	0.73
1:D:80:MSE:HE3	1:D:90:ILE:HG13	1.69	0.73
1:A:80:MSE:HE3	1:A:90:ILE:HG13	1.70	0.72
1:A:96:ILE:HG13	1:A:115:ALA:HA	1.72	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:146:ILE:HD12	1:C:146:ILE:H	1.55	0.72
1:C:96:ILE:HG13	1:C:115:ALA:HA	1.72	0.72
1:C:51:VAL:HG23	1:C:118:VAL:HG22	1.72	0.71
1:D:186:THR:HG22	1:D:189:GLY:N	2.00	0.71
1:A:186:THR:HG22	1:A:189:GLY:N	2.04	0.71
1:C:146:ILE:H	1:C:146:ILE:CD1	2.03	0.71
1:B:41:GLY:O	1:B:46:VAL:HG22	1.91	0.70
1:D:41:GLY:O	1:D:46:VAL:HG22	1.91	0.70
1:C:146:ILE:HD12	1:C:146:ILE:N	2.07	0.70
1:B:96:ILE:HG13	1:B:115:ALA:HA	1.73	0.69
1:C:186:THR:HG22	1:C:189:GLY:N	2.05	0.69
1:D:51:VAL:HG23	1:D:118:VAL:HG22	1.74	0.69
1:B:51:VAL:HG23	1:B:118:VAL:HG22	1.75	0.69
1:D:96:ILE:HG13	1:D:115:ALA:HA	1.73	0.69
1:A:100:ILE:HG13	1:A:119:GLN:HE21	1.57	0.68
1:A:177:LEU:O	1:A:180:LEU:HD12	1.92	0.68
1:B:144:GLN:HB2	1:B:145:PRO:HD2	1.75	0.68
1:C:177:LEU:O	1:C:180:LEU:HD12	1.93	0.68
1:A:128:GLY:N	1:A:200:MSE:HE3	2.09	0.68
1:A:146:ILE:HD12	1:A:146:ILE:N	2.09	0.68
1:A:151:LEU:HD22	1:A:152:PRO:HD2	1.75	0.68
1:D:74:GLN:O	1:D:77:ILE:HG22	1.95	0.67
1:D:144:GLN:HB2	1:D:145:PRO:HD2	1.77	0.67
1:A:144:GLN:HB2	1:A:145:PRO:HD2	1.77	0.67
1:D:128:GLY:N	1:D:200:MSE:HE3	2.10	0.67
1:C:120:ASP:O	1:C:124:LEU:HB2	1.95	0.67
1:B:128:GLY:N	1:B:200:MSE:HE3	2.09	0.67
1:C:144:GLN:HB2	1:C:145:PRO:HD2	1.75	0.67
1:B:154:ARG:HD3	1:B:175:GLU:OE1	1.94	0.66
1:A:51:VAL:HG23	1:A:118:VAL:HG22	1.76	0.66
1:C:57:PHE:HA	1:C:60:LEU:CD2	2.25	0.66
1:A:47:ASP:HB3	1:A:50:VAL:HG12	1.76	0.66
1:A:146:ILE:HD12	1:A:146:ILE:H	1.59	0.66
1:D:100:ILE:HG13	1:D:119:GLN:HE21	1.61	0.66
1:B:100:ILE:HG13	1:B:119:GLN:HE21	1.60	0.65
1:C:47:ASP:HB3	1:C:50:VAL:HG12	1.76	0.65
1:A:154:ARG:HH11	1:A:167:SER:HB2	1.59	0.65
1:B:74:GLN:O	1:B:77:ILE:HG22	1.96	0.65
1:A:41:GLY:O	1:A:46:VAL:HG22	1.96	0.65
1:B:47:ASP:HB3	1:B:50:VAL:HG12	1.77	0.65
1:D:47:ASP:HB3	1:D:50:VAL:HG12	1.77	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:128:GLY:H	1:B:200:MSE:HE3	1.62	0.64
1:C:100:ILE:HG13	1:C:119:GLN:HE21	1.61	0.64
1:A:57:PHE:HA	1:A:60:LEU:CD2	2.26	0.64
1:C:70:GLU:HG2	1:C:71:THR:HG22	1.79	0.64
1:B:57:PHE:HA	1:B:60:LEU:CD2	2.27	0.64
1:C:113:ARG:CA	1:C:116:MSE:HE3	2.28	0.64
1:D:130:ILE:HG13	1:D:131:GLY:N	2.12	0.64
1:A:120:ASP:O	1:A:124:LEU:HB2	1.98	0.63
1:A:163:ARG:HH11	1:A:163:ARG:HB3	1.63	0.63
1:C:41:GLY:O	1:C:46:VAL:HG22	1.99	0.63
1:D:128:GLY:H	1:D:200:MSE:HE3	1.64	0.63
1:C:128:GLY:N	1:C:200:MSE:HE3	2.14	0.63
1:D:154:ARG:HD3	1:D:175:GLU:OE1	1.99	0.63
1:A:70:GLU:HG2	1:A:71:THR:HG22	1.79	0.63
1:D:2:ASN:CG	1:D:3:GLU:N	2.52	0.63
1:A:142:LYS:NZ	1:A:162:TYR:OH	2.19	0.63
1:D:57:PHE:HA	1:D:60:LEU:CD2	2.29	0.63
1:B:120:ASP:O	1:B:124:LEU:HB2	1.98	0.62
1:A:128:GLY:H	1:A:200:MSE:HE3	1.63	0.62
1:B:177:LEU:O	1:B:180:LEU:HD12	2.00	0.62
1:C:163:ARG:HB3	1:C:163:ARG:HH11	1.63	0.62
1:D:177:LEU:O	1:D:180:LEU:HD12	1.98	0.62
1:D:2:ASN:OD1	1:D:3:GLU:N	2.32	0.62
1:D:120:ASP:O	1:D:124:LEU:HB2	1.99	0.62
1:A:16:LYS:HD3	1:A:17:GLN:N	2.15	0.62
1:B:70:GLU:HG2	1:B:71:THR:HG22	1.82	0.62
1:A:128:GLY:O	1:A:132:ILE:HG13	2.00	0.62
1:D:184:MSE:HE3	1:D:193:ALA:CB	2.29	0.62
1:D:70:GLU:HG2	1:D:71:THR:HG22	1.81	0.61
1:A:127:LEU:HD23	1:A:177:LEU:HD22	1.83	0.61
1:A:74:GLN:O	1:A:77:ILE:HG22	2.00	0.61
1:B:16:LYS:HD3	1:B:17:GLN:N	2.16	0.61
1:C:130:ILE:HG12	1:D:210:GLU:OE1	2.00	0.61
1:A:13:TRP:HD1	1:A:16:LYS:HD2	1.66	0.61
1:B:130:ILE:HG13	1:B:131:GLY:N	2.15	0.61
1:C:130:ILE:HG13	1:C:131:GLY:N	2.15	0.61
1:B:13:TRP:HD1	1:B:16:LYS:HD2	1.66	0.61
1:B:184:MSE:HE3	1:B:193:ALA:CB	2.30	0.61
1:D:127:LEU:HD23	1:D:177:LEU:HD22	1.83	0.61
1:B:127:LEU:HD23	1:B:177:LEU:HD22	1.83	0.61
1:A:136:PHE:CD1	1:A:146:ILE:HG23	2.35	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:128:GLY:O	1:C:132:ILE:HG13	2.01	0.61
1:D:113:ARG:CA	1:D:116:MSE:HE3	2.31	0.60
1:C:13:TRP:HD1	1:C:16:LYS:HD2	1.66	0.60
1:D:13:TRP:HD1	1:D:16:LYS:HD2	1.66	0.60
1:D:163:ARG:HH11	1:D:163:ARG:HB3	1.66	0.60
1:A:112:THR:HG22	1:A:113:ARG:N	2.16	0.60
1:A:130:ILE:HG13	1:A:131:GLY:N	2.16	0.60
1:C:16:LYS:HD3	1:C:17:GLN:N	2.16	0.60
1:A:181:LYS:HD3	1:A:197:HIS:CG	2.37	0.60
1:B:163:ARG:HH11	1:B:163:ARG:HB3	1.66	0.60
1:D:81:GLU:HG3	1:D:90:ILE:CD1	2.32	0.59
1:B:50:VAL:HG23	1:B:93:THR:HA	1.84	0.59
1:C:128:GLY:H	1:C:200:MSE:HE3	1.67	0.59
1:A:130:ILE:HG12	1:B:210:GLU:OE1	2.03	0.59
1:D:16:LYS:HD3	1:D:17:GLN:N	2.17	0.59
1:D:136:PHE:CD1	1:D:146:ILE:HG22	2.36	0.59
1:C:114:GLU:O	1:C:118:VAL:HG23	2.03	0.59
1:C:74:GLN:O	1:C:77:ILE:HG22	2.02	0.59
1:D:50:VAL:HG23	1:D:93:THR:HA	1.84	0.59
1:C:50:VAL:HG23	1:C:93:THR:HA	1.83	0.59
1:B:127:LEU:HD21	1:B:180:LEU:HD12	1.84	0.59
1:D:178:PHE:HA	1:D:181:LYS:NZ	2.18	0.59
1:C:112:THR:HG22	1:C:113:ARG:N	2.17	0.58
1:A:188:THR:O	1:A:191:GLN:HB3	2.03	0.58
1:C:51:VAL:HG23	1:C:118:VAL:CG2	2.33	0.58
1:D:181:LYS:HD3	1:D:197:HIS:CG	2.38	0.58
1:A:50:VAL:HG23	1:A:93:THR:HA	1.85	0.58
1:A:184:MSE:HE3	1:A:193:ALA:CB	2.34	0.58
1:B:178:PHE:HA	1:B:181:LYS:NZ	2.18	0.58
1:B:181:LYS:HD3	1:B:197:HIS:CG	2.37	0.58
1:C:127:LEU:HD23	1:C:177:LEU:HD22	1.86	0.58
1:A:81:GLU:HG3	1:A:90:ILE:CD1	2.34	0.58
1:B:154:ARG:HB2	1:B:175:GLU:OE1	2.03	0.58
1:B:113:ARG:CA	1:B:116:MSE:HE3	2.30	0.57
1:D:128:GLY:O	1:D:132:ILE:HG13	2.03	0.57
1:C:181:LYS:HD3	1:C:197:HIS:CG	2.39	0.57
1:B:188:THR:O	1:B:191:GLN:HB3	2.03	0.57
1:C:188:THR:O	1:C:191:GLN:HB3	2.04	0.57
1:B:136:PHE:CD1	1:B:146:ILE:HG22	2.39	0.57
1:D:186:THR:HG22	1:D:188:THR:N	2.20	0.57
1:B:181:LYS:HD3	1:B:197:HIS:ND1	2.20	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:113:ARG:CA	1:A:116:MSE:HE3	2.30	0.57
1:B:2:ASN:CG	1:B:3:GLU:N	2.58	0.57
1:B:81:GLU:HG3	1:B:90:ILE:CD1	2.35	0.57
1:D:70:GLU:HG2	1:D:71:THR:N	2.20	0.57
1:A:136:PHE:CE1	1:A:146:ILE:HG23	2.40	0.57
1:D:2:ASN:CG	1:D:3:GLU:H	2.08	0.57
1:A:114:GLU:O	1:A:118:VAL:HG23	2.05	0.56
1:A:51:VAL:HG23	1:A:118:VAL:CG2	2.35	0.56
1:D:13:TRP:O	1:D:16:LYS:HG3	2.05	0.56
1:D:51:VAL:HG23	1:D:118:VAL:CG2	2.35	0.56
1:D:188:THR:O	1:D:191:GLN:HB3	2.05	0.56
1:C:81:GLU:HG3	1:C:90:ILE:CD1	2.35	0.56
1:D:181:LYS:HD3	1:D:197:HIS:ND1	2.19	0.56
1:B:186:THR:HG22	1:B:188:THR:N	2.20	0.56
1:B:170:ILE:O	1:B:173:PHE:HB2	2.06	0.56
1:B:13:TRP:O	1:B:16:LYS:HG3	2.06	0.56
1:B:70:GLU:HG2	1:B:71:THR:N	2.20	0.56
1:B:184:MSE:HE3	1:B:193:ALA:HB2	1.87	0.56
1:C:184:MSE:HE3	1:C:193:ALA:CB	2.36	0.56
1:D:184:MSE:HE3	1:D:193:ALA:HB2	1.87	0.56
1:A:178:PHE:HA	1:A:181:LYS:NZ	2.21	0.55
1:B:142:LYS:NZ	1:B:162:TYR:OH	2.24	0.55
1:D:178:PHE:HA	1:D:181:LYS:HZ2	1.70	0.55
1:A:181:LYS:HD3	1:A:197:HIS:ND1	2.21	0.55
1:B:178:PHE:HA	1:B:181:LYS:HZ2	1.71	0.55
1:A:70:GLU:HG2	1:A:71:THR:N	2.22	0.55
1:C:33:VAL:HG23	1:C:122:ASP:HA	1.89	0.55
1:D:33:VAL:HG23	1:D:122:ASP:HA	1.89	0.55
1:C:70:GLU:HG2	1:C:71:THR:N	2.22	0.55
1:A:33:VAL:HG23	1:A:122:ASP:HA	1.89	0.55
1:D:114:GLU:O	1:D:118:VAL:HG23	2.07	0.54
1:D:170:ILE:O	1:D:173:PHE:HB2	2.08	0.54
1:B:51:VAL:HG23	1:B:118:VAL:CG2	2.37	0.54
1:A:146:ILE:H	1:A:146:ILE:CD1	2.20	0.54
1:B:112:THR:HG22	1:B:113:ARG:N	2.21	0.54
1:C:178:PHE:HA	1:C:181:LYS:NZ	2.22	0.54
1:A:170:ILE:O	1:A:173:PHE:HB2	2.07	0.54
1:A:127:LEU:HD21	1:A:180:LEU:HD12	1.90	0.54
1:D:154:ARG:HB2	1:D:175:GLU:OE1	2.08	0.54
1:C:204:ILE:C	1:C:204:ILE:HD12	2.28	0.54
1:D:42:GLU:O	1:D:45:LYS:HD2	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:204:ILE:C	1:A:204:ILE:HD12	2.28	0.53
1:D:96:ILE:HG12	1:D:100:ILE:HD13	1.91	0.53
1:B:114:GLU:O	1:B:118:VAL:HG23	2.08	0.53
1:C:93:THR:O	1:C:96:ILE:HG23	2.08	0.53
1:D:77:ILE:HD13	1:D:77:ILE:O	2.08	0.53
1:D:142:LYS:NZ	1:D:162:TYR:OH	2.23	0.53
1:D:204:ILE:HD12	1:D:204:ILE:C	2.29	0.53
1:B:204:ILE:HD12	1:B:204:ILE:C	2.29	0.53
1:A:13:TRP:O	1:A:16:LYS:HG3	2.08	0.53
1:C:170:ILE:O	1:C:173:PHE:HB2	2.09	0.53
1:C:147:TYR:HB2	1:C:170:ILE:CG1	2.39	0.53
1:D:112:THR:HG22	1:D:113:ARG:N	2.22	0.53
1:A:93:THR:O	1:A:96:ILE:HG23	2.08	0.52
1:D:148:ASP:HB3	1:D:151:LEU:HB2	1.91	0.52
1:B:128:GLY:O	1:B:132:ILE:HG13	2.09	0.52
1:A:47:ASP:HB3	1:A:50:VAL:CG1	2.39	0.52
1:B:33:VAL:HG23	1:B:122:ASP:HA	1.90	0.52
1:D:93:THR:O	1:D:96:ILE:HG23	2.09	0.52
1:D:74:GLN:HA	1:D:77:ILE:HG22	1.92	0.52
1:C:47:ASP:HB3	1:C:50:VAL:CG1	2.40	0.52
1:B:146:ILE:HD13	1:B:146:ILE:N	2.11	0.52
1:B:35:LEU:O	1:B:38:LYS:HB3	2.10	0.52
1:A:186:THR:HG22	1:A:188:THR:N	2.25	0.51
1:B:42:GLU:O	1:B:45:LYS:HD2	2.10	0.51
1:B:86:PRO:HG2	1:B:89:LYS:HB2	1.91	0.51
1:C:2:ASN:CG	1:C:5:ALA:HB3	2.31	0.51
1:C:38:LYS:O	1:C:41:GLY:N	2.43	0.51
1:D:127:LEU:HD21	1:D:180:LEU:HD12	1.91	0.51
1:D:47:ASP:HB3	1:D:50:VAL:CG1	2.40	0.51
1:D:86:PRO:HG2	1:D:89:LYS:HB2	1.91	0.51
1:A:74:GLN:HA	1:A:77:ILE:HG22	1.92	0.51
1:B:47:ASP:HB3	1:B:50:VAL:CG1	2.41	0.51
1:C:186:THR:HG22	1:C:188:THR:N	2.25	0.51
1:C:136:PHE:CD1	1:C:146:ILE:HG22	2.45	0.51
1:C:74:GLN:HA	1:C:77:ILE:HG22	1.92	0.51
1:D:146:ILE:H	1:D:146:ILE:CD1	2.11	0.51
1:D:151:LEU:HD13	1:D:152:PRO:HD2	1.92	0.51
1:D:181:LYS:H	1:D:181:LYS:CD	2.24	0.51
1:C:181:LYS:HD3	1:C:197:HIS:ND1	2.25	0.51
1:B:60:LEU:O	1:B:61:ILE:HG12	2.11	0.51
1:B:136:PHE:HD1	1:B:146:ILE:HG22	1.76	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:96:ILE:HG12	1:B:100:ILE:HD13	1.93	0.50
1:C:60:LEU:H	1:C:60:LEU:HD23	1.76	0.50
1:C:96:ILE:HG12	1:C:100:ILE:HD13	1.93	0.50
1:C:153:ILE:HG22	1:C:154:ARG:H	1.75	0.50
1:C:13:TRP:O	1:C:16:LYS:HG3	2.11	0.50
1:C:136:PHE:CE1	1:C:146:ILE:HG22	2.47	0.50
1:C:77:ILE:HD13	1:C:77:ILE:O	2.12	0.50
1:A:184:MSE:HE3	1:A:193:ALA:HB2	1.94	0.50
1:B:181:LYS:H	1:B:181:LYS:CD	2.24	0.50
1:B:93:THR:O	1:B:96:ILE:HG23	2.11	0.50
1:B:87:SER:HA	1:B:90:ILE:HG22	1.94	0.50
1:C:54:ALA:HB1	1:C:96:ILE:HD12	1.94	0.50
1:A:136:PHE:CE1	1:A:146:ILE:CG2	2.95	0.50
1:B:77:ILE:O	1:B:77:ILE:HD13	2.12	0.50
1:C:127:LEU:HD21	1:C:180:LEU:HD12	1.94	0.50
1:C:181:LYS:H	1:C:181:LYS:CD	2.25	0.50
1:A:77:ILE:HD13	1:A:77:ILE:O	2.12	0.49
1:D:60:LEU:O	1:D:61:ILE:HG12	2.12	0.49
1:C:178:PHE:HA	1:C:181:LYS:HZ2	1.76	0.49
1:D:157:MSE:O	1:D:158:THR:HG23	2.12	0.49
1:A:87:SER:HA	1:A:90:ILE:HG22	1.94	0.49
1:A:120:ASP:O	1:A:124:LEU:N	2.44	0.49
1:A:54:ALA:HB1	1:A:96:ILE:HD12	1.94	0.49
1:A:60:LEU:H	1:A:60:LEU:HD23	1.77	0.49
1:A:96:ILE:HG12	1:A:100:ILE:HD13	1.95	0.49
1:B:38:LYS:O	1:B:41:GLY:N	2.46	0.49
1:D:60:LEU:HD23	1:D:60:LEU:H	1.76	0.49
1:D:73:LYS:HD2	1:D:94:MSE:HE3	1.95	0.49
1:C:86:PRO:HG2	1:C:89:LYS:HB2	1.95	0.49
1:A:181:LYS:H	1:A:181:LYS:CD	2.26	0.49
1:B:60:LEU:HD23	1:B:60:LEU:H	1.76	0.49
1:B:73:LYS:HD2	1:B:94:MSE:HE3	1.95	0.49
1:C:54:ALA:HB1	1:C:96:ILE:CD1	2.43	0.49
1:D:35:LEU:O	1:D:38:LYS:HB3	2.13	0.49
1:B:177:LEU:O	1:B:181:LYS:NZ	2.43	0.49
1:C:13:TRP:CD1	1:C:16:LYS:HD2	2.48	0.49
1:C:60:LEU:O	1:C:61:ILE:HG12	2.13	0.49
1:A:146:ILE:N	1:A:146:ILE:CD1	2.76	0.49
1:B:74:GLN:HA	1:B:77:ILE:HG22	1.94	0.49
1:A:28:TYR:CE2	1:B:149:PRO:HG3	2.47	0.48
1:D:87:SER:HA	1:D:90:ILE:HG22	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:42:GLU:O	1:C:45:LYS:HD2	2.13	0.48
1:B:48:VAL:O	1:B:52:GLN:HG3	2.14	0.48
1:A:38:LYS:O	1:A:41:GLY:N	2.46	0.48
1:D:38:LYS:O	1:D:41:GLY:N	2.46	0.48
1:C:87:SER:HA	1:C:90:ILE:HG22	1.94	0.48
1:D:48:VAL:O	1:D:52:GLN:HG3	2.12	0.48
1:C:94:MSE:O	1:C:97:ILE:HB	2.14	0.48
1:C:120:ASP:O	1:C:124:LEU:N	2.41	0.48
1:A:54:ALA:HB1	1:A:96:ILE:CD1	2.44	0.48
1:D:33:VAL:CG2	1:D:122:ASP:HA	2.44	0.48
1:A:60:LEU:O	1:A:61:ILE:HG12	2.13	0.48
1:B:146:ILE:H	1:B:146:ILE:CD1	2.03	0.47
1:A:192:LEU:O	1:A:196:ARG:HG2	2.14	0.47
1:B:154:ARG:NH2	1:B:175:GLU:OE1	2.45	0.47
1:B:13:TRP:CD1	1:B:16:LYS:HD2	2.49	0.47
1:C:48:VAL:O	1:C:52:GLN:HG3	2.14	0.47
1:C:147:TYR:HB2	1:C:170:ILE:HG13	1.96	0.47
1:D:72:ALA:O	1:D:75:GLN:HB3	2.15	0.47
1:B:157:MSE:O	1:B:158:THR:HG23	2.14	0.47
1:A:86:PRO:HG2	1:A:89:LYS:HB2	1.96	0.47
1:A:2:ASN:HD21	1:A:5:ALA:HB3	1.79	0.47
1:A:112:THR:CG2	1:A:113:ARG:N	2.78	0.47
1:B:3:GLU:HA	1:B:6:ILE:HD12	1.97	0.47
1:A:94:MSE:O	1:A:97:ILE:HB	2.14	0.47
1:A:157:MSE:O	1:A:158:THR:HG23	2.15	0.46
1:D:177:LEU:O	1:D:181:LYS:NZ	2.46	0.46
1:A:2:ASN:O	1:A:3:GLU:HG2	2.14	0.46
1:A:42:GLU:O	1:A:45:LYS:HD2	2.14	0.46
1:B:33:VAL:CG2	1:B:122:ASP:HA	2.45	0.46
1:B:158:THR:O	1:C:73:LYS:NZ	2.48	0.46
1:A:178:PHE:HA	1:A:181:LYS:HZ2	1.78	0.46
1:B:54:ALA:HB1	1:B:96:ILE:HD12	1.98	0.46
1:C:57:PHE:HD1	1:C:60:LEU:HD11	1.81	0.46
1:C:184:MSE:HE3	1:C:193:ALA:HB2	1.96	0.46
1:D:54:ALA:HB1	1:D:96:ILE:CD1	2.46	0.46
1:B:120:ASP:O	1:B:124:LEU:N	2.43	0.46
1:B:54:ALA:HB1	1:B:96:ILE:CD1	2.45	0.46
1:B:148:ASP:OD1	1:B:150:GLU:HB2	2.16	0.46
1:C:112:THR:CG2	1:C:113:ARG:N	2.79	0.46
1:A:13:TRP:CD1	1:A:16:LYS:HD2	2.48	0.46
1:C:157:MSE:O	1:C:158:THR:HG23	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:2:ASN:ND2	1:C:5:ALA:HB3	2.30	0.45
1:D:13:TRP:CD1	1:D:16:LYS:HD2	2.48	0.45
1:A:54:ALA:HA	1:A:97:ILE:HD11	1.99	0.45
1:C:33:VAL:CG2	1:C:122:ASP:HA	2.46	0.45
1:D:130:ILE:HG13	1:D:131:GLY:H	1.81	0.45
1:A:73:LYS:HD2	1:A:94:MSE:HE3	1.98	0.45
1:A:87:SER:HA	1:A:90:ILE:CG2	2.47	0.45
1:C:192:LEU:O	1:C:196:ARG:HG2	2.16	0.45
1:C:87:SER:HA	1:C:90:ILE:CG2	2.47	0.45
1:C:124:LEU:HD11	1:C:192:LEU:HB2	1.98	0.45
1:C:4:GLN:O	1:C:4:GLN:HG2	2.17	0.45
1:C:4:GLN:O	1:C:8:GLN:OE1	2.34	0.45
1:D:54:ALA:HB1	1:D:96:ILE:HD12	1.98	0.45
1:A:48:VAL:O	1:A:52:GLN:HG3	2.16	0.45
1:B:13:TRP:HH2	1:B:75:GLN:HG2	1.82	0.45
1:C:208:LEU:HD12	1:C:209:SER:N	2.32	0.45
1:C:146:ILE:O	1:C:169:ALA:HB3	2.17	0.45
1:A:57:PHE:HD1	1:A:60:LEU:HD11	1.81	0.45
1:A:208:LEU:HD12	1:A:209:SER:N	2.32	0.45
1:D:208:LEU:O	1:D:211:TRP:HB3	2.17	0.45
1:C:35:LEU:O	1:C:38:LYS:HB3	2.17	0.44
1:C:148:ASP:OD1	1:C:150:GLU:HB2	2.16	0.44
1:C:177:LEU:O	1:C:181:LYS:NZ	2.46	0.44
1:A:33:VAL:CG2	1:A:122:ASP:HA	2.47	0.44
1:B:160:GLU:HA	1:B:163:ARG:NH1	2.33	0.44
1:B:208:LEU:O	1:B:211:TRP:HB3	2.17	0.44
1:C:120:ASP:OD1	1:C:186:THR:HB	2.17	0.44
1:A:35:LEU:O	1:A:38:LYS:HB3	2.17	0.44
1:B:54:ALA:HA	1:B:97:ILE:HD11	2.00	0.44
1:B:153:ILE:H	1:B:153:ILE:HG13	1.51	0.44
1:C:94:MSE:HA	1:C:97:ILE:CB	2.48	0.44
1:D:160:GLU:HA	1:D:163:ARG:NH1	2.33	0.44
1:A:94:MSE:HA	1:A:97:ILE:CB	2.48	0.44
1:C:183:LEU:HD22	1:C:183:LEU:N	2.33	0.44
1:D:61:ILE:HD13	1:D:61:ILE:HA	1.89	0.44
1:B:74:GLN:NE2	1:B:77:ILE:CG2	2.81	0.43
1:B:208:LEU:HD12	1:B:209:SER:N	2.33	0.43
1:A:203:PHE:CD1	1:A:203:PHE:C	2.92	0.43
1:C:203:PHE:CD1	1:C:203:PHE:C	2.92	0.43
1:D:42:GLU:H	1:D:42:GLU:HG2	1.68	0.43
1:C:54:ALA:HA	1:C:97:ILE:HD11	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:74:GLN:NE2	1:C:77:ILE:CG2	2.81	0.43
1:A:53:ILE:CG2	1:A:93:THR:HG21	2.49	0.43
1:B:72:ALA:O	1:B:75:GLN:HB3	2.18	0.43
1:D:13:TRP:HH2	1:D:75:GLN:HG2	1.83	0.43
1:D:192:LEU:O	1:D:196:ARG:HG2	2.18	0.43
1:A:183:LEU:N	1:A:183:LEU:HD22	2.34	0.43
1:B:32:ARG:HD2	1:B:125:ASP:OD1	2.19	0.43
1:C:53:ILE:CG2	1:C:93:THR:HG21	2.48	0.43
1:D:112:THR:CG2	1:D:113:ARG:N	2.82	0.43
1:B:147:TYR:HB2	1:B:170:ILE:HG13	2.00	0.43
1:C:13:TRP:CZ2	1:C:75:GLN:NE2	2.87	0.43
1:A:74:GLN:NE2	1:A:77:ILE:CG2	2.82	0.43
1:A:160:GLU:HA	1:A:163:ARG:NH1	2.33	0.43
1:A:177:LEU:O	1:A:181:LYS:NZ	2.47	0.43
1:C:14:VAL:HG13	1:C:60:LEU:HD22	2.01	0.43
1:C:163:ARG:O	1:C:163:ARG:HG2	2.17	0.43
1:B:87:SER:HA	1:B:90:ILE:CG2	2.49	0.43
1:B:112:THR:CG2	1:B:113:ARG:N	2.81	0.43
1:B:126:ALA:O	1:B:177:LEU:HD21	2.17	0.43
1:B:146:ILE:O	1:B:169:ALA:HB3	2.18	0.43
1:C:38:LYS:O	1:C:39:ALA:C	2.57	0.43
1:D:32:ARG:HD2	1:D:125:ASP:OD1	2.18	0.43
1:D:147:TYR:HB2	1:D:170:ILE:HG13	1.99	0.43
1:C:153:ILE:HG22	1:C:154:ARG:N	2.33	0.43
1:D:153:ILE:O	1:D:154:ARG:HG3	2.19	0.43
1:C:160:GLU:HA	1:C:163:ARG:NH1	2.34	0.42
1:C:204:ILE:HD12	1:C:205:GLU:N	2.34	0.42
1:A:53:ILE:HG22	1:A:93:THR:HG21	2.01	0.42
1:C:120:ASP:HA	1:C:123:ARG:HG2	2.01	0.42
1:D:74:GLN:HA	1:D:77:ILE:CG2	2.48	0.42
1:A:32:ARG:HD2	1:A:125:ASP:OD1	2.19	0.42
1:B:57:PHE:HD1	1:B:60:LEU:HD11	1.84	0.42
1:C:13:TRP:HH2	1:C:75:GLN:HG2	1.85	0.42
1:C:73:LYS:HD2	1:C:94:MSE:HE3	2.01	0.42
1:D:120:ASP:O	1:D:124:LEU:N	2.44	0.42
1:C:163:ARG:HB3	1:C:163:ARG:NH1	2.33	0.42
1:D:146:ILE:HD13	1:D:146:ILE:N	2.21	0.42
1:D:87:SER:HA	1:D:90:ILE:CG2	2.48	0.42
1:A:13:TRP:HH2	1:A:75:GLN:HG2	1.84	0.42
1:A:14:VAL:HG13	1:A:60:LEU:HD22	2.00	0.42
1:B:14:VAL:HG13	1:B:60:LEU:HD22	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:51:VAL:HG13	1:B:52:GLN:N	2.34	0.42
1:A:124:LEU:HD11	1:A:192:LEU:HB2	2.01	0.42
1:C:47:ASP:OD2	1:C:89:LYS:HE2	2.20	0.42
1:D:14:VAL:O	1:D:18:LEU:HD12	2.20	0.42
1:D:94:MSE:O	1:D:97:ILE:HB	2.19	0.42
1:D:204:ILE:HD12	1:D:205:GLU:N	2.34	0.42
1:D:208:LEU:HD12	1:D:209:SER:N	2.34	0.42
1:A:163:ARG:O	1:A:163:ARG:HG2	2.19	0.42
1:D:3:GLU:HA	1:D:6:ILE:HD12	2.01	0.42
1:B:74:GLN:HA	1:B:77:ILE:CG2	2.50	0.42
1:D:58:HIS:NE2	1:D:122:ASP:OD1	2.53	0.42
1:B:92:HIS:CE1	1:B:112:THR:HG21	2.55	0.42
1:C:91:ASP:O	1:C:95:ASP:HB2	2.20	0.42
1:B:29:HIS:O	1:B:33:VAL:HG12	2.19	0.41
1:B:58:HIS:NE2	1:B:122:ASP:OD1	2.53	0.41
1:B:192:LEU:O	1:B:196:ARG:HG2	2.19	0.41
1:C:32:ARG:HD2	1:C:125:ASP:OD1	2.19	0.41
1:C:154:ARG:NH1	1:C:167:SER:HB2	2.35	0.41
1:D:36:MSE:O	1:D:40:ILE:HG13	2.20	0.41
1:A:13:TRP:CZ2	1:A:75:GLN:NE2	2.88	0.41
1:B:16:LYS:HD3	1:B:17:GLN:H	1.85	0.41
1:B:94:MSE:O	1:B:97:ILE:HB	2.20	0.41
1:C:94:MSE:HA	1:C:97:ILE:HB	2.01	0.41
1:D:74:GLN:NE2	1:D:77:ILE:CG2	2.83	0.41
1:B:147:TYR:HB2	1:B:170:ILE:CG1	2.50	0.41
1:B:151:LEU:HA	1:B:151:LEU:HD22	1.77	0.41
1:A:74:GLN:HA	1:A:77:ILE:CG2	2.51	0.41
1:A:204:ILE:HD12	1:A:205:GLU:N	2.36	0.41
1:D:13:TRP:CZ2	1:D:75:GLN:NE2	2.88	0.41
1:D:57:PHE:HD1	1:D:60:LEU:HD11	1.85	0.41
1:D:112:THR:HG22	1:D:114:GLU:H	1.86	0.41
1:A:183:LEU:N	1:A:183:LEU:CD2	2.84	0.41
1:C:28:TYR:CE2	1:D:149:PRO:HG3	2.56	0.41
1:D:92:HIS:CE1	1:D:112:THR:HG21	2.55	0.41
1:D:163:ARG:HB3	1:D:163:ARG:NH1	2.34	0.41
1:C:183:LEU:N	1:C:183:LEU:CD2	2.83	0.41
1:D:80:MSE:CE	1:D:90:ILE:HG13	2.45	0.41
1:B:154:ARG:HB2	1:B:154:ARG:HH21	1.84	0.41
1:D:16:LYS:HD3	1:D:17:GLN:H	1.86	0.41
1:A:14:VAL:O	1:A:18:LEU:HD12	2.21	0.41
1:A:51:VAL:HG13	1:A:52:GLN:N	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:94:MSE:HA	1:A:97:ILE:HB	2.02	0.41
1:B:13:TRP:CZ2	1:B:75:GLN:NE2	2.89	0.41
1:B:120:ASP:HA	1:B:123:ARG:HG2	2.03	0.41
1:B:146:ILE:N	1:B:146:ILE:CD1	2.80	0.41
1:C:72:ALA:O	1:C:75:GLN:HB3	2.21	0.41
1:D:54:ALA:HA	1:D:97:ILE:HD11	2.02	0.41
1:A:47:ASP:OD2	1:A:89:LYS:HE2	2.20	0.41
1:D:92:HIS:HE1	1:D:112:THR:HG21	1.86	0.41
1:A:120:ASP:HA	1:A:123:ARG:HG2	2.03	0.40
1:B:61:ILE:HD13	1:B:61:ILE:HA	1.88	0.40
1:C:74:GLN:HA	1:C:77:ILE:CG2	2.51	0.40
1:A:42:GLU:H	1:A:42:GLU:HG2	1.68	0.40
1:A:178:PHE:HE1	1:A:200:MSE:HB2	1.87	0.40
1:C:51:VAL:HG13	1:C:52:GLN:N	2.37	0.40
1:A:43:GLN:OE1	1:A:43:GLN:HA	2.22	0.40
1:D:51:VAL:HG13	1:D:52:GLN:N	2.36	0.40
1:D:94:MSE:HA	1:D:97:ILE:CB	2.51	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	179/223 (80%)	146 (82%)	24 (13%)	9 (5%)	1	12
1	B	179/223 (80%)	144 (80%)	27 (15%)	8 (4%)	2	13
1	C	179/223 (80%)	146 (82%)	25 (14%)	8 (4%)	2	13
1	D	179/223 (80%)	147 (82%)	24 (13%)	8 (4%)	2	13
All	All	716/892 (80%)	583 (81%)	100 (14%)	33 (5%)	2	13

All (33) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	161	GLU
1	B	161	GLU
1	D	161	GLU
1	A	60	LEU
1	A	165	GLY
1	A	212	ASN
1	B	60	LEU
1	B	61	ILE
1	B	165	GLY
1	C	60	LEU
1	C	161	GLU
1	C	165	GLY
1	D	60	LEU
1	D	61	ILE
1	D	165	GLY
1	A	61	ILE
1	A	93	THR
1	A	158	THR
1	A	164	HIS
1	B	93	THR
1	B	158	THR
1	B	164	HIS
1	C	61	ILE
1	C	93	THR
1	C	158	THR
1	C	164	HIS
1	D	93	THR
1	D	158	THR
1	D	164	HIS
1	D	178	PHE
1	B	178	PHE
1	A	178	PHE
1	C	178	PHE

### 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	161/181 (89%)	140 (87%)	21 (13%)	3	14
1	B	161/181 (89%)	137 (85%)	24 (15%)	2	11
1	C	161/181 (89%)	139 (86%)	22 (14%)	3	13
1	D	161/181 (89%)	138 (86%)	23 (14%)	2	12
All	All	644/724 (89%)	554 (86%)	90 (14%)	3	13

All (90) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	LYS
1	A	17	GLN
1	A	19	MSE
1	A	27	TRP
1	A	43	GLN
1	A	52	GLN
1	A	58	HIS
1	A	60	LEU
1	A	62	ASP
1	A	70	GLU
1	A	77	ILE
1	A	94	MSE
1	A	96	ILE
1	A	124	LEU
1	A	151	LEU
1	A	158	THR
1	A	160	GLU
1	A	162	TYR
1	A	181	LYS
1	A	182	ASP
1	A	186	THR
1	B	16	LYS
1	B	17	GLN
1	B	19	MSE
1	B	27	TRP
1	B	35	LEU
1	B	43	GLN
1	B	52	GLN
1	B	58	HIS
1	B	60	LEU
1	B	62	ASP
1	B	70	GLU
1	B	77	ILE

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Mol	Chain	Res	Type
1	B	94	MSE
1	B	96	ILE
1	B	124	LEU
1	B	146	ILE
1	B	151	LEU
1	B	153	ILE
1	B	154	ARG
1	B	158	THR
1	B	160	GLU
1	B	162	TYR
1	B	181	LYS
1	B	182	ASP
1	C	16	LYS
1	C	17	GLN
1	C	19	MSE
1	C	27	TRP
1	C	43	GLN
1	C	52	GLN
1	C	58	HIS
1	C	60	LEU
1	C	62	ASP
1	C	70	GLU
1	C	77	ILE
1	C	94	MSE
1	C	96	ILE
1	C	124	LEU
1	C	146	ILE
1	C	151	LEU
1	C	158	THR
1	C	160	GLU
1	C	162	TYR
1	C	181	LYS
1	C	182	ASP
1	C	186	THR
1	D	16	LYS
1	D	17	GLN
1	D	19	MSE
1	D	27	TRP
1	D	43	GLN
1	D	52	GLN
1	D	58	HIS
1	D	60	LEU

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Mol	Chain	Res	Type
1	D	62	ASP
1	D	70	GLU
1	D	77	ILE
1	D	94	MSE
1	D	96	ILE
1	D	124	LEU
1	D	146	ILE
1	D	151	LEU
1	D	153	ILE
1	D	158	THR
1	D	160	GLU
1	D	162	TYR
1	D	181	LYS
1	D	182	ASP
1	D	186	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	GLN
1	A	75	GLN
1	A	88	GLN
1	A	92	HIS
1	A	119	GLN
1	A	144	GLN
1	A	171	ASN
1	A	212	ASN
1	B	17	GLN
1	B	75	GLN
1	B	88	GLN
1	B	92	HIS
1	B	119	GLN
1	B	144	GLN
1	B	171	ASN
1	C	8	GLN
1	C	17	GLN
1	C	75	GLN
1	C	88	GLN
1	C	92	HIS
1	C	119	GLN
1	C	144	GLN
1	C	171	ASN

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Mol	Chain	Res	Type
1	C	212	ASN
1	D	17	GLN
1	D	75	GLN
1	D	88	GLN
1	D	92	HIS
1	D	119	GLN
1	D	144	GLN
1	D	171	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

### 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

### 5.7 Other polymers [i](#)

There are no such residues in this entry.

### 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data [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	181/223 (81%)	-1.70	0 100 100	31, 65, 101, 118	0
1	B	181/223 (81%)	-1.68	0 100 100	29, 64, 93, 119	0
1	C	181/223 (81%)	-1.72	0 100 100	30, 65, 103, 118	0
1	D	181/223 (81%)	-1.63	0 100 100	26, 63, 94, 119	0
All	All	724/892 (81%)	-1.68	0 100 100	26, 64, 101, 119	0

There are no RSRZ outliers to report.

### 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 monosaccharides 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.