



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 10, 2021 – 06:11 PM EDT

PDB ID : 3ENO  
Title : Crystal structure of Pyrococcus furiosus Pcc1 in complex with Thermoplasma acidophilum Kae1  
Authors : Neculai, D.  
Deposited on : 2008-09-25  
Resolution : 3.02 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.

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

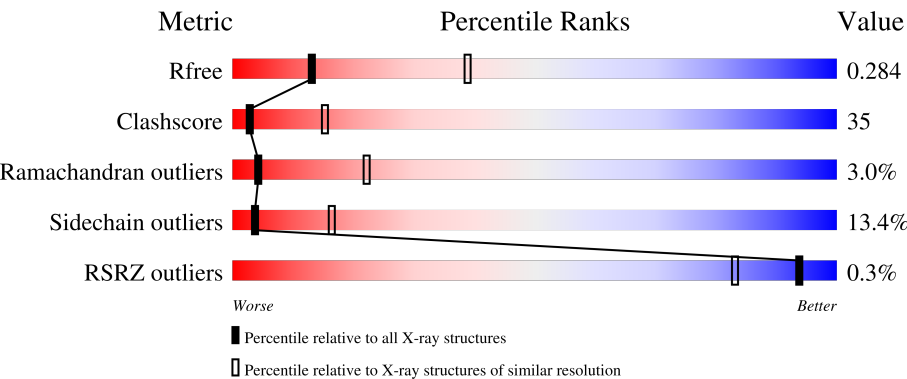
MolProbity : 4.02b-467  
Xtriage (Phenix) : 1.13  
EDS : 2.23.2  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.23.2

# 1 Overall quality at a glance i

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

The reported resolution of this entry is 3.02 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	2399 (3.04-3.00)
Clashscore	141614	2734 (3.04-3.00)
Ramachandran outliers	138981	2640 (3.04-3.00)
Sidechain outliers	138945	2643 (3.04-3.00)
RSRZ outliers	127900	2287 (3.04-3.00)

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	334	
1	B	334	
2	C	87	
2	D	87	
2	E	87	

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Mol	Chain	Length	Quality of chain
2	F	87	<div><div></div><div>24%</div><div>56%</div><div>10%</div><div>9%</div></div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7471 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 Putative O-sialoglycoprotein endopeptidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	329	Total	C	N	O	S	0	0	0
			2488	1571	432	466	19			
1	B	328	Total	C	N	O	S	0	0	0
			2480	1566	431	465	18			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	GLY	-	expression tag	UNP Q9HLA5
A	-3	ALA	-	expression tag	UNP Q9HLA5
A	-2	MET	-	expression tag	UNP Q9HLA5
A	-1	ASP	-	expression tag	UNP Q9HLA5
A	0	PRO	-	expression tag	UNP Q9HLA5
B	-4	GLY	-	expression tag	UNP Q9HLA5
B	-3	ALA	-	expression tag	UNP Q9HLA5
B	-2	MET	-	expression tag	UNP Q9HLA5
B	-1	ASP	-	expression tag	UNP Q9HLA5
B	0	PRO	-	expression tag	UNP Q9HLA5

- Molecule 2 is a protein called uncharacterized protein PF2011.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	77	Total	C	N	O	S	0	0	0
			624	402	104	117	1			
2	D	76	Total	C	N	O	S	0	0	0
			615	397	103	114	1			
2	E	77	Total	C	N	O	S	0	0	0
			624	402	104	117	1			
2	F	79	Total	C	N	O	S	0	0	0
			638	411	107	119	1			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-4	GLY	-	expression tag	UNP Q8TZI1
C	-3	ALA	-	expression tag	UNP Q8TZI1
C	-2	MET	-	expression tag	UNP Q8TZI1
C	-1	ASP	-	expression tag	UNP Q8TZI1
C	0	PRO	-	expression tag	UNP Q8TZI1
C	12	MET	ILE	engineered mutation	UNP Q8TZI1
D	-4	GLY	-	expression tag	UNP Q8TZI1
D	-3	ALA	-	expression tag	UNP Q8TZI1
D	-2	MET	-	expression tag	UNP Q8TZI1
D	-1	ASP	-	expression tag	UNP Q8TZI1
D	0	PRO	-	expression tag	UNP Q8TZI1
D	12	MET	ILE	engineered mutation	UNP Q8TZI1
E	-4	GLY	-	expression tag	UNP Q8TZI1
E	-3	ALA	-	expression tag	UNP Q8TZI1
E	-2	MET	-	expression tag	UNP Q8TZI1
E	-1	ASP	-	expression tag	UNP Q8TZI1
E	0	PRO	-	expression tag	UNP Q8TZI1
E	12	MET	ILE	engineered mutation	UNP Q8TZI1
F	-4	GLY	-	expression tag	UNP Q8TZI1
F	-3	ALA	-	expression tag	UNP Q8TZI1
F	-2	MET	-	expression tag	UNP Q8TZI1
F	-1	ASP	-	expression tag	UNP Q8TZI1
F	0	PRO	-	expression tag	UNP Q8TZI1
F	12	MET	ILE	engineered mutation	UNP Q8TZI1

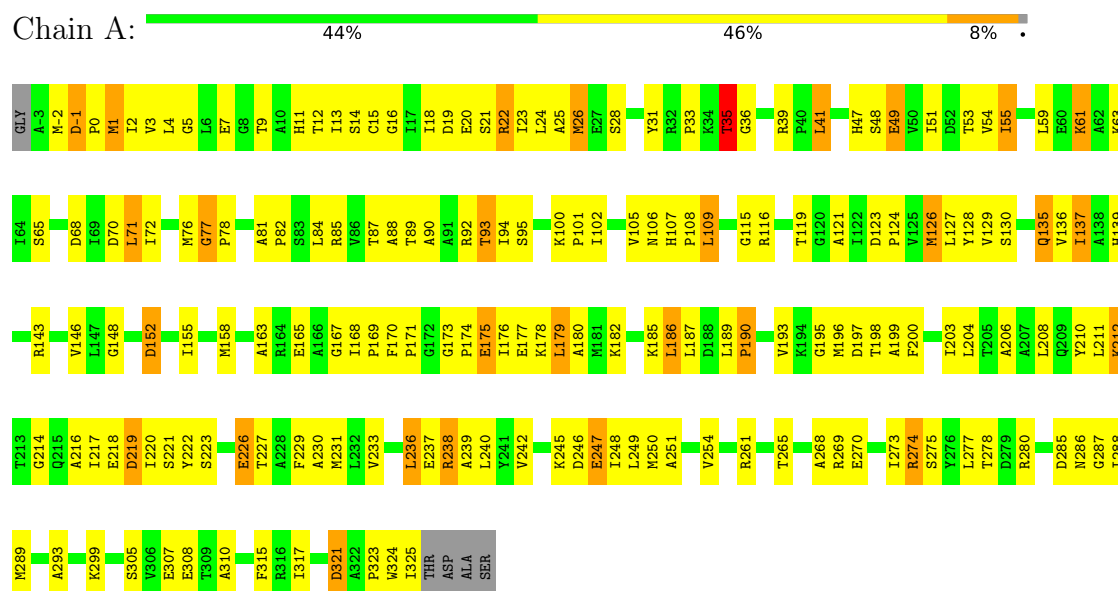
- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Mg 1 1	0	0
3	B	1	Total Mg 1 1	0	0

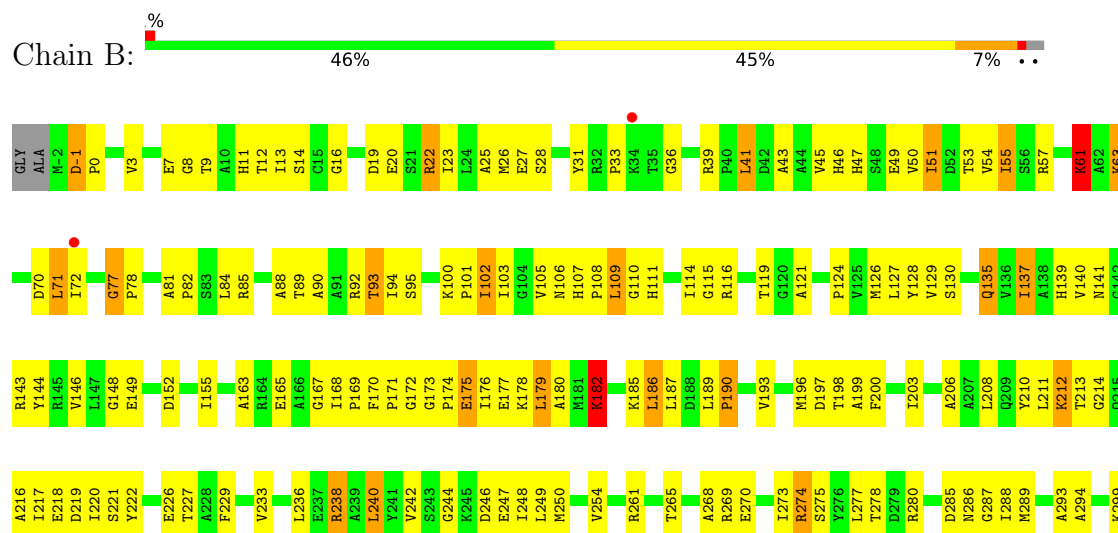
### 3 Residue-property plots [i](#)

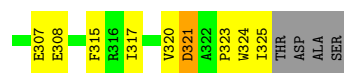
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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

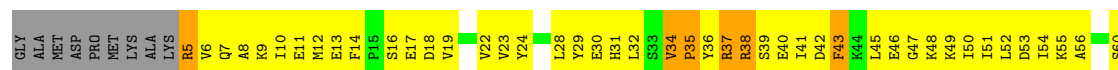


- Molecule 1: Putative O-sialoglycoprotein endopeptidase

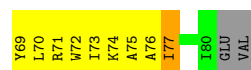




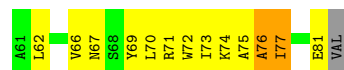
• Molecule 2: uncharacterized protein PF2011



• Molecule 2: uncharacterized protein PF2011



• Molecule 2: uncharacterized protein PF2011



• Molecule 2: uncharacterized protein PF2011



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	66.56Å 66.56Å 435.56Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	39.58 – 3.02 39.58 – 3.01	Depositor EDS
% Data completeness (in resolution range)	95.3 (39.58-3.02) 99.8 (39.58-3.01)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	0.05	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.63 (at 3.01Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, $R_{free}$	0.243 , 0.282 0.242 , 0.284	Depositor DCC
$R_{free}$ test set	1083 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	74.2	Xtriage
Anisotropy	0.454	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 59.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.438 for h,-h-k,-l	Xtriage
Reported twinning fraction	0.435 for h,-h-k,-l	Depositor
Outliers	0 of 21550 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	7471	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	84.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.39% 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](#)

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	0.26	0/2526	0.49	0/3416
1	B	0.26	0/2518	0.49	0/3406
2	C	0.25	0/634	0.46	0/855
2	D	0.26	0/625	0.46	0/843
2	E	0.25	0/634	0.45	0/855
2	F	0.25	0/648	0.46	0/873
All	All	0.26	0/7585	0.48	0/10248

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	2488	0	2559	190	0
1	B	2480	0	2547	174	0
2	C	624	0	642	60	0
2	D	615	0	636	36	0
2	E	624	0	642	45	0
2	F	638	0	660	54	0
3	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	1	0	0	0	0
All	All	7471	0	7686	532	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 35.

All (532) 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:128:TYR:HB3	1:B:135:GLN:HB2	1.37	1.07
1:B:26:MET:HB2	1:B:288:ILE:HD12	1.42	0.99
1:B:71:LEU:HD22	1:B:72:ILE:H	1.28	0.96
2:C:7:GLN:HE22	2:C:55:LYS:HE2	1.33	0.93
1:A:128:TYR:HB3	1:A:135:GLN:HB2	1.51	0.92
1:A:71:LEU:HD22	1:A:72:ILE:H	1.38	0.88
2:F:7:GLN:HE22	2:F:55:LYS:HE2	1.40	0.85
1:B:163:ALA:HB2	1:B:176:ILE:HD11	1.56	0.85
1:B:26:MET:HG3	1:B:288:ILE:HB	1.59	0.85
1:B:240:LEU:HD13	1:B:273:ILE:HD11	1.59	0.83
1:A:26:MET:HB2	1:A:288:ILE:HD12	1.59	0.83
1:A:208:LEU:HD13	2:E:35:PRO:HB3	1.59	0.83
1:B:169:PRO:HB2	1:B:171:PRO:HD2	1.58	0.83
1:B:286:ASN:ND2	1:B:288:ILE:HG22	1.95	0.82
1:B:23:ILE:HG23	1:B:23:ILE:O	1.80	0.82
1:A:127:LEU:HD23	1:A:250:MET:HG2	1.61	0.81
1:B:121:ALA:HB1	1:B:124:PRO:HG3	1.62	0.81
1:A:240:LEU:HD13	1:A:273:ILE:HD11	1.62	0.81
1:A:23:ILE:HG23	1:A:23:ILE:O	1.79	0.80
1:A:169:PRO:HB2	1:A:171:PRO:HD2	1.64	0.80
1:B:22:ARG:HD3	1:B:23:ILE:H	1.47	0.78
1:A:143:ARG:HG2	1:A:321:ASP:HA	1.64	0.78
1:B:286:ASN:ND2	1:B:289:MET:H	1.82	0.77
1:B:208:LEU:HD13	2:D:35:PRO:HB3	1.66	0.77
1:B:22:ARG:HA	1:B:22:ARG:NE	2.00	0.76
1:B:307:GLU:HA	2:F:29:TYR:CE2	2.21	0.76
1:A:121:ALA:HB1	1:A:124:PRO:HG3	1.67	0.76
2:E:7:GLN:HE22	2:E:55:LYS:HE2	1.50	0.76
1:B:-1:ASP:OD1	1:B:0:PRO:HD2	1.86	0.75
2:D:28:LEU:O	2:D:32:LEU:HG	1.86	0.75
1:B:11:HIS:HD2	1:B:31:TYR:HD2	1.34	0.75
1:A:163:ALA:HB2	1:A:176:ILE:HD11	1.67	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:137:ILE:HG22	1:B:146:VAL:HA	1.68	0.74
1:B:71:LEU:HD22	1:B:72:ILE:N	2.03	0.73
1:A:94:ILE:HG12	2:C:79:VAL:CG1	2.19	0.73
1:A:155:ILE:HD13	1:A:200:PHE:CE1	2.23	0.73
1:B:121:ALA:CB	1:B:124:PRO:HG3	2.20	0.72
1:A:196:MET:HG3	1:A:238:ARG:HG2	1.70	0.72
1:A:148:GLY:HA3	1:A:196:MET:O	1.90	0.71
1:A:286:ASN:ND2	1:A:289:MET:H	1.87	0.71
1:A:317:ILE:HD12	1:A:317:ILE:H	1.56	0.71
2:C:38:ARG:HE	2:E:81:GLU:C	1.94	0.71
2:F:31:HIS:HD2	2:F:43:PHE:HD2	1.40	0.70
1:A:11:HIS:HD2	1:A:31:TYR:HD2	1.39	0.70
1:A:22:ARG:HA	1:A:22:ARG:NE	2.05	0.70
2:F:67:ASN:HB3	2:F:71:ARG:NH2	2.05	0.70
2:C:67:ASN:HB3	2:C:71:ARG:NH2	2.07	0.70
1:A:26:MET:HG3	1:A:288:ILE:HB	1.73	0.70
1:B:143:ARG:HG2	1:B:321:ASP:HA	1.74	0.69
1:A:90:ALA:O	1:A:94:ILE:HG13	1.92	0.69
1:A:7:GLU:HB3	1:A:287:GLY:HA2	1.72	0.69
1:B:81:ALA:O	1:B:85:ARG:HG3	1.92	0.69
1:B:155:ILE:HD13	1:B:200:PHE:CE1	2.28	0.69
1:B:321:ASP:O	1:B:323:PRO:HD3	1.93	0.69
2:F:31:HIS:HD2	2:F:43:PHE:CD2	2.11	0.68
1:A:307:GLU:HA	2:C:29:TYR:CE2	2.27	0.68
1:A:109:LEU:HD13	1:A:293:ALA:HB1	1.75	0.68
1:B:180:ALA:HB1	1:B:222:TYR:HA	1.76	0.68
2:F:45:LEU:HD23	2:F:46:GLU:H	1.58	0.68
2:D:67:ASN:O	2:D:71:ARG:HB2	1.94	0.68
2:D:7:GLN:HE22	2:D:55:LYS:HE2	1.59	0.68
1:B:81:ALA:HB3	1:B:82:PRO:HD3	1.76	0.68
1:A:81:ALA:HB3	1:A:82:PRO:HD3	1.74	0.67
2:C:42:ASP:HB3	2:C:53:ASP:HB2	1.77	0.67
2:F:49:LYS:HB2	2:F:49:LYS:NZ	2.10	0.67
1:A:35:THR:HG23	1:A:36:GLY:H	1.58	0.67
1:A:180:ALA:HB1	1:A:222:TYR:HA	1.77	0.67
1:A:20:GLU:OE2	1:A:299:LYS:HE3	1.94	0.66
1:B:90:ALA:O	1:B:94:ILE:HG13	1.96	0.66
1:A:240:LEU:HD22	1:A:273:ILE:HD12	1.77	0.66
2:F:42:ASP:HB3	2:F:53:ASP:HB2	1.78	0.66
1:A:72:ILE:HD12	1:A:95:SER:HA	1.77	0.66
1:A:121:ALA:CB	1:A:124:PRO:HG3	2.26	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:71:LEU:CD2	1:B:72:ILE:H	2.06	0.66
1:B:31:TYR:O	1:B:33:PRO:HD3	1.95	0.66
1:B:143:ARG:HD3	1:B:321:ASP:OD2	1.96	0.65
2:F:31:HIS:CD2	2:F:43:PHE:HD2	2.13	0.65
2:F:5:ARG:O	2:F:5:ARG:HD3	1.95	0.65
2:C:38:ARG:NH1	2:C:64:GLY:HA3	2.11	0.65
1:A:106:ASN:OD1	1:A:108:PRO:HD2	1.97	0.65
1:B:25:ALA:CB	1:B:61:LYS:HB3	2.27	0.64
1:A:317:ILE:HD12	1:A:317:ILE:N	2.12	0.64
2:E:69:TYR:O	2:E:73:ILE:HG13	1.98	0.64
1:A:71:LEU:CD2	1:A:72:ILE:H	2.10	0.64
1:B:22:ARG:NH1	1:B:23:ILE:HG22	2.12	0.64
1:B:203:ILE:H	1:B:203:ILE:HD12	1.62	0.64
2:C:30:GLU:OE2	2:C:30:GLU:HA	1.98	0.64
2:D:49:LYS:HB2	2:D:49:LYS:NZ	2.13	0.64
1:B:7:GLU:HB3	1:B:287:GLY:HA2	1.80	0.64
1:B:89:THR:OG1	2:F:30:GLU:HG3	1.97	0.64
2:E:18:ASP:O	2:E:22:VAL:HG23	1.98	0.63
2:E:49:LYS:HB2	2:E:49:LYS:NZ	2.13	0.63
1:A:286:ASN:HD21	1:A:289:MET:HB3	1.62	0.63
1:A:22:ARG:HD3	1:A:23:ILE:H	1.63	0.63
1:A:71:LEU:HD22	1:A:72:ILE:N	2.12	0.63
1:A:240:LEU:HD21	1:A:248:ILE:CD1	2.28	0.63
2:E:6:VAL:HG21	2:E:62:LEU:HD22	1.79	0.63
1:A:94:ILE:HG12	2:C:79:VAL:HG11	1.79	0.63
1:A:186:LEU:HD22	1:A:187:LEU:O	1.98	0.63
1:B:22:ARG:HH11	1:B:23:ILE:HG22	1.64	0.63
1:B:148:GLY:HA3	1:B:196:MET:O	1.99	0.63
2:D:18:ASP:O	2:D:22:VAL:HG23	1.99	0.63
2:F:35:PRO:HG2	2:F:36:TYR:H	1.63	0.63
1:B:70:ASP:O	1:B:101:PRO:HD2	1.99	0.62
1:B:218:GLU:CD	1:B:218:GLU:H	2.02	0.62
2:E:55:LYS:NZ	2:E:55:LYS:HB3	2.14	0.62
1:A:137:ILE:HG22	1:A:146:VAL:HA	1.80	0.62
1:B:94:ILE:HG12	2:F:79:VAL:CG1	2.29	0.62
1:B:174:PRO:O	1:B:178:LYS:HG2	1.98	0.62
2:E:7:GLN:NE2	2:E:55:LYS:HE2	2.14	0.62
1:A:25:ALA:CB	1:A:61:LYS:HB3	2.28	0.62
1:B:26:MET:SD	1:B:286:ASN:HB2	2.39	0.62
2:D:6:VAL:HG21	2:D:62:LEU:HD22	1.82	0.62
2:E:31:HIS:CE1	2:E:37:ARG:HH12	2.18	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:174:PRO:O	1:A:178:LYS:HG2	2.00	0.61
1:B:240:LEU:HD21	1:B:248:ILE:CD1	2.29	0.61
1:A:286:ASN:ND2	1:A:288:ILE:HG22	2.15	0.61
1:B:88:ALA:HB1	1:B:92:ARG:NH2	2.15	0.61
2:D:8:ALA:HB3	2:D:54:ILE:HB	1.83	0.61
2:F:14:PHE:CD1	2:F:77:ILE:HD11	2.35	0.61
2:F:6:VAL:HG21	2:F:62:LEU:HD22	1.83	0.61
1:A:25:ALA:O	1:A:26:MET:HB2	2.01	0.61
1:B:286:ASN:HD22	1:B:288:ILE:HG22	1.66	0.61
2:C:38:ARG:HH12	2:C:64:GLY:HA3	1.65	0.61
2:F:55:LYS:HB3	2:F:55:LYS:NZ	2.16	0.61
2:C:49:LYS:HB2	2:C:49:LYS:NZ	2.16	0.60
1:B:129:VAL:HB	1:B:254:VAL:HB	1.83	0.60
2:F:38:ARG:O	2:F:39:SER:HB3	2.01	0.60
1:A:-1:ASP:OD1	1:A:0:PRO:HD2	2.01	0.60
1:A:240:LEU:HD11	1:A:248:ILE:HD11	1.83	0.60
2:C:31:HIS:HD2	2:C:43:PHE:HD2	1.47	0.60
1:A:94:ILE:HG12	2:C:79:VAL:HG13	1.83	0.60
1:B:240:LEU:HD22	1:B:273:ILE:HD12	1.83	0.60
1:A:143:ARG:HD3	1:A:321:ASP:OD2	2.01	0.60
1:B:107:HIS:HB3	1:B:108:PRO:HD3	1.82	0.59
1:B:317:ILE:HD12	1:B:317:ILE:N	2.16	0.59
1:A:31:TYR:O	1:A:33:PRO:HD3	2.02	0.59
1:B:26:MET:CG	1:B:288:ILE:HB	2.31	0.59
1:B:139:HIS:CE1	1:B:325:ILE:HD13	2.37	0.59
1:B:196:MET:HG3	1:B:238:ARG:HG2	1.84	0.59
2:E:8:ALA:HB3	2:E:54:ILE:HB	1.85	0.59
1:B:78:PRO:HG3	1:B:315:PHE:O	2.02	0.59
1:B:165:GLU:HG2	1:B:208:LEU:HD22	1.83	0.59
1:B:179:LEU:CD2	1:B:218:GLU:HG3	2.32	0.59
2:E:28:LEU:O	2:E:32:LEU:HG	2.03	0.58
1:A:88:ALA:HB1	1:A:92:ARG:NH2	2.18	0.58
2:C:5:ARG:HG2	2:C:7:GLN:HE21	1.69	0.58
1:B:20:GLU:OE2	1:B:299:LYS:HE3	2.03	0.58
1:B:212:LYS:C	1:B:214:GLY:H	2.06	0.58
1:B:16:GLY:HA3	1:B:288:ILE:HG13	1.83	0.58
1:B:23:ILE:O	1:B:23:ILE:CG2	2.51	0.58
2:C:16:SER:OG	2:C:19:VAL:HG23	2.03	0.58
1:A:268:ALA:HB2	1:A:275:SER:HB3	1.86	0.58
2:C:5:ARG:O	2:C:5:ARG:HD3	2.02	0.58
2:C:31:HIS:HD2	2:C:43:PHE:CD2	2.21	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:16:GLY:HA3	1:A:288:ILE:HG13	1.84	0.58
1:B:72:ILE:HD12	1:B:95:SER:HA	1.85	0.58
2:C:55:LYS:NZ	2:C:55:LYS:HB3	2.19	0.58
2:C:18:ASP:O	2:C:22:VAL:HG23	2.04	0.57
2:E:9:LYS:NZ	2:E:11:GLU:OE2	2.37	0.57
1:A:11:HIS:HD2	1:A:31:TYR:CD2	2.22	0.57
1:B:61:LYS:HB2	1:B:61:LYS:HZ2	1.70	0.57
1:A:22:ARG:NH1	1:A:23:ILE:HG22	2.20	0.57
1:B:317:ILE:HD12	1:B:317:ILE:H	1.69	0.57
2:C:5:ARG:HG2	2:C:7:GLN:NE2	2.19	0.57
2:F:16:SER:OG	2:F:19:VAL:HG23	2.03	0.57
1:A:70:ASP:O	1:A:101:PRO:HD2	2.03	0.57
2:F:28:LEU:O	2:F:32:LEU:HG	2.04	0.57
2:F:45:LEU:HD23	2:F:46:GLU:N	2.19	0.57
1:B:111:HIS:HB3	1:B:126:MET:HE1	1.85	0.57
2:E:11:GLU:HG2	2:E:51:ILE:HD12	1.86	0.57
1:B:25:ALA:HB2	1:B:61:LYS:HB3	1.87	0.57
2:D:11:GLU:HG2	2:D:51:ILE:HD12	1.87	0.57
1:A:12:THR:HG21	1:A:285:ASP:O	2.05	0.56
2:F:18:ASP:O	2:F:22:VAL:HG23	2.05	0.56
1:A:81:ALA:O	1:A:85:ARG:HG3	2.05	0.56
2:C:55:LYS:HB3	2:C:55:LYS:HZ2	1.70	0.56
2:D:55:LYS:NZ	2:D:55:LYS:HB3	2.21	0.56
2:C:31:HIS:CD2	2:C:43:PHE:HD2	2.22	0.56
1:A:165:GLU:HG2	1:A:208:LEU:HD22	1.87	0.56
1:B:94:ILE:HG12	2:F:79:VAL:HG13	1.88	0.56
1:B:185:LYS:HG2	1:B:186:LEU:N	2.21	0.56
2:C:31:HIS:CD2	2:C:43:PHE:HB2	2.41	0.56
1:B:51:ILE:O	1:B:55:ILE:HG23	2.06	0.56
2:D:13:GLU:HB2	2:D:49:LYS:HE3	1.87	0.56
2:F:17:GLU:HA	2:F:48:LYS:HB3	1.88	0.56
1:A:139:HIS:CE1	1:A:325:ILE:HD13	2.41	0.56
1:B:7:GLU:CB	1:B:287:GLY:HA2	2.36	0.56
1:B:186:LEU:HD22	1:B:187:LEU:O	2.07	0.55
1:A:126:MET:HE3	1:A:251:ALA:HB3	1.89	0.55
1:A:216:ALA:HB3	1:A:219:ASP:HB2	1.88	0.55
2:C:28:LEU:O	2:C:32:LEU:HG	2.06	0.55
2:E:72:TRP:O	2:E:75:ALA:HB3	2.07	0.55
2:E:49:LYS:HB2	2:E:49:LYS:HZ2	1.72	0.55
1:B:321:ASP:C	1:B:323:PRO:HD3	2.26	0.55
2:F:31:HIS:CE1	2:F:37:ARG:HH12	2.24	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:65:SER:HB3	1:A:68:ASP:OD2	2.06	0.55
1:B:250:MET:CE	1:B:277:LEU:HD22	2.36	0.55
1:A:89:THR:O	1:A:93:THR:HG22	2.06	0.55
1:A:185:LYS:HG2	1:A:186:LEU:N	2.22	0.55
1:A:179:LEU:CD2	1:A:218:GLU:HG3	2.36	0.55
1:B:77:GLY:HA2	1:B:315:PHE:HD1	1.71	0.55
2:D:7:GLN:NE2	2:D:55:LYS:HE2	2.21	0.55
2:F:14:PHE:CE1	2:F:77:ILE:HD11	2.42	0.55
1:A:1:MET:HB3	1:A:20:GLU:OE2	2.07	0.55
1:A:107:HIS:HB3	1:A:108:PRO:HD3	1.89	0.54
1:A:72:ILE:O	1:A:102:ILE:HG13	2.07	0.54
1:A:100:LYS:HB3	1:A:101:PRO:CD	2.37	0.54
1:A:18:ILE:HG22	1:A:23:ILE:HA	1.89	0.54
1:A:11:HIS:O	1:A:47:HIS:HE1	1.91	0.54
2:C:9:LYS:HA	2:C:52:LEU:O	2.07	0.54
2:E:67:ASN:O	2:E:71:ARG:HB2	2.07	0.54
1:A:26:MET:CG	1:A:288:ILE:HB	2.38	0.54
1:A:176:ILE:HG23	1:A:221:SER:HB2	1.89	0.53
1:A:240:LEU:HD21	1:A:248:ILE:HD12	1.90	0.53
1:A:218:GLU:CD	1:A:218:GLU:H	2.11	0.53
1:B:212:LYS:C	1:B:214:GLY:N	2.62	0.53
2:F:31:HIS:CD2	2:F:43:PHE:HB2	2.42	0.53
1:B:238:ARG:O	1:B:242:VAL:HG23	2.08	0.53
2:C:60:SER:HB2	2:D:74:LYS:HD2	1.89	0.53
2:E:66:VAL:HG13	2:E:70:LEU:HD12	1.91	0.53
1:A:25:ALA:HB1	1:A:61:LYS:HB3	1.89	0.53
1:A:286:ASN:ND2	1:A:289:MET:N	2.57	0.53
1:B:3:VAL:HG22	1:B:71:LEU:HB3	1.90	0.53
1:B:100:LYS:HB3	1:B:101:PRO:CD	2.39	0.53
2:E:5:ARG:O	2:E:5:ARG:HD3	2.09	0.53
2:D:24:TYR:O	2:D:28:LEU:HB2	2.09	0.53
1:B:268:ALA:HB2	1:B:275:SER:HB3	1.89	0.53
2:F:8:ALA:HB3	2:F:54:ILE:HB	1.90	0.53
2:F:43:PHE:HE1	2:F:50:ILE:HG21	1.74	0.53
1:B:250:MET:HE2	1:B:277:LEU:HD22	1.90	0.53
2:C:5:ARG:HD2	2:C:5:ARG:N	2.24	0.52
2:E:66:VAL:CG1	2:E:70:LEU:HD12	2.39	0.52
2:F:24:TYR:O	2:F:28:LEU:HB2	2.09	0.52
1:B:11:HIS:HD2	1:B:31:TYR:CD2	2.20	0.52
1:A:90:ALA:O	1:A:93:THR:HG23	2.09	0.52
1:A:196:MET:SD	1:A:239:ALA:HA	2.49	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:6:VAL:HG22	2:D:56:ALA:HB3	1.90	0.52
1:B:25:ALA:HB1	1:B:61:LYS:HB3	1.91	0.52
1:A:124:PRO:HA	1:A:247:GLU:HG3	1.90	0.52
1:A:93:THR:HG21	2:C:76:ALA:HA	1.90	0.52
1:B:43:ALA:O	1:B:46:HIS:HB3	2.10	0.52
1:B:127:LEU:HD23	1:B:250:MET:HG2	1.91	0.52
2:C:7:GLN:NE2	2:C:55:LYS:HE2	2.15	0.52
1:A:116:ARG:HH12	1:A:325:ILE:CG1	2.22	0.52
1:A:128:TYR:HD2	1:A:135:GLN:HG3	1.75	0.52
1:A:168:ILE:HG22	1:A:169:PRO:O	2.10	0.52
1:B:31:TYR:CE2	1:B:33:PRO:HG3	2.45	0.52
1:B:109:LEU:HD13	1:B:293:ALA:HB1	1.92	0.52
1:B:203:ILE:HD12	1:B:203:ILE:N	2.24	0.52
2:C:6:VAL:HG21	2:C:62:LEU:HD22	1.91	0.52
1:B:90:ALA:O	1:B:93:THR:HG23	2.10	0.51
1:B:105:VAL:HG12	1:B:106:ASN:N	2.24	0.51
1:B:240:LEU:HD11	1:B:248:ILE:HD11	1.91	0.51
2:F:37:ARG:HB3	2:F:39:SER:O	2.10	0.51
2:C:23:VAL:HG13	2:C:76:ALA:HB3	1.91	0.51
1:A:23:ILE:O	1:A:23:ILE:CG2	2.51	0.51
1:A:250:MET:HE2	1:A:277:LEU:HD22	1.91	0.51
1:B:187:LEU:HB3	1:B:210:TYR:HE1	1.75	0.51
1:B:189:LEU:HB2	1:B:227:THR:HG21	1.92	0.51
2:C:37:ARG:HB3	2:C:39:SER:O	2.10	0.51
1:B:307:GLU:HA	2:F:29:TYR:HE2	1.72	0.51
2:C:31:HIS:HD2	2:C:43:PHE:HB2	1.76	0.51
1:B:54:VAL:HG23	1:B:55:ILE:H	1.76	0.51
1:B:324:TRP:H	1:B:325:ILE:HD12	1.75	0.51
2:C:14:PHE:CD1	2:C:77:ILE:HD11	2.45	0.51
2:F:7:GLN:NE2	2:F:55:LYS:HE2	2.19	0.51
1:B:265:THR:O	1:B:268:ALA:HB3	2.11	0.51
2:D:72:TRP:O	2:D:75:ALA:HB3	2.11	0.51
1:A:182:LYS:O	1:A:182:LYS:HG3	2.11	0.51
1:B:94:ILE:HG12	2:F:79:VAL:HG11	1.93	0.51
1:A:78:PRO:HD2	1:A:107:HIS:HB3	1.93	0.50
1:B:115:GLY:O	1:B:119:THR:HG23	2.10	0.50
2:E:6:VAL:HG22	2:E:56:ALA:HB3	1.93	0.50
1:A:261:ARG:NH1	1:A:280:ARG:HG2	2.26	0.50
1:A:7:GLU:CB	1:A:287:GLY:HA2	2.41	0.50
1:A:195:GLY:O	1:A:238:ARG:NH1	2.45	0.50
1:A:321:ASP:C	1:A:323:PRO:HD3	2.32	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:116:ARG:HH12	1:B:325:ILE:CG1	2.24	0.50
2:D:5:ARG:O	2:D:5:ARG:HD3	2.11	0.50
1:B:25:ALA:O	1:B:26:MET:HB2	2.11	0.50
1:A:229:PHE:O	1:A:233:VAL:HG23	2.11	0.50
1:A:321:ASP:O	1:A:323:PRO:HD3	2.11	0.50
1:A:317:ILE:H	1:A:317:ILE:CD1	2.24	0.50
2:C:8:ALA:HB3	2:C:54:ILE:HB	1.93	0.50
1:B:12:THR:HG21	1:B:285:ASP:O	2.12	0.49
1:B:54:VAL:HG23	1:B:55:ILE:N	2.27	0.49
2:C:45:LEU:HD23	2:C:46:GLU:N	2.26	0.49
1:A:22:ARG:HH11	1:A:23:ILE:HG22	1.76	0.49
1:A:115:GLY:O	1:A:119:THR:HG23	2.12	0.49
2:D:9:LYS:NZ	2:D:11:GLU:OE2	2.45	0.49
2:F:18:ASP:O	2:F:21:LYS:HB2	2.12	0.49
1:A:25:ALA:HB2	1:A:61:LYS:HB3	1.92	0.49
1:A:208:LEU:CD1	2:E:35:PRO:HB3	2.37	0.49
1:A:274:ARG:HE	1:A:274:ARG:N	2.11	0.49
2:D:66:VAL:CG1	2:D:70:LEU:HD12	2.42	0.49
1:A:152:ASP:O	2:E:29:TYR:CE2	2.66	0.49
1:B:182:LYS:O	1:B:182:LYS:HG3	2.13	0.49
1:B:261:ARG:NH1	1:B:280:ARG:HG2	2.27	0.49
2:F:13:GLU:HB2	2:F:49:LYS:HE3	1.95	0.49
1:B:168:ILE:HG22	1:B:169:PRO:O	2.13	0.49
1:A:23:ILE:HG13	1:A:288:ILE:CD1	2.43	0.48
1:B:89:THR:O	1:B:93:THR:HG22	2.13	0.48
1:A:212:LYS:C	1:A:214:GLY:H	2.15	0.48
2:F:37:ARG:C	2:F:39:SER:H	2.15	0.48
1:A:187:LEU:HB3	1:A:210:TYR:HE1	1.78	0.48
1:B:36:GLY:HA2	1:B:170:PHE:CD2	2.48	0.48
2:F:43:PHE:HE1	2:F:50:ILE:CG2	2.27	0.48
1:A:61:LYS:HZ2	1:A:61:LYS:HB2	1.79	0.48
2:C:17:GLU:HA	2:C:48:LYS:HB3	1.95	0.48
1:A:105:VAL:HG12	1:A:106:ASN:O	2.13	0.48
1:A:89:THR:OG1	2:C:30:GLU:HG3	2.14	0.48
1:A:175:GLU:O	1:A:179:LEU:HB2	2.14	0.48
2:C:34:VAL:HA	2:C:35:PRO:HD3	1.73	0.48
2:E:6:VAL:CG2	2:E:62:LEU:HD22	2.43	0.48
2:D:6:VAL:HG21	2:D:62:LEU:HB2	1.96	0.48
1:A:187:LEU:HD13	1:A:210:TYR:HD1	1.79	0.48
1:A:14:SER:HB3	1:A:287:GLY:N	2.29	0.47
1:B:229:PHE:O	1:B:233:VAL:HG23	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:274:ARG:N	1:B:274:ARG:HE	2.12	0.47
2:C:56:ALA:HB3	2:C:62:LEU:HB2	1.96	0.47
1:B:103:ILE:HD11	1:B:294:ALA:HB1	1.95	0.47
1:B:246:ASP:HB2	1:B:273:ILE:HG22	1.96	0.47
1:B:249:LEU:HD22	1:B:278:THR:CG2	2.44	0.47
1:B:7:GLU:HB3	1:B:287:GLY:CA	2.45	0.47
1:B:216:ALA:O	1:B:220:ILE:HG13	2.15	0.47
1:B:116:ARG:HH12	1:B:325:ILE:HG12	1.80	0.47
1:B:203:ILE:H	1:B:203:ILE:CD1	2.26	0.47
2:D:45:LEU:HD23	2:D:46:GLU:N	2.30	0.47
1:A:26:MET:SD	1:A:286:ASN:HB2	2.55	0.47
1:A:70:ASP:O	1:A:100:LYS:HB3	2.15	0.47
1:A:324:TRP:H	1:A:325:ILE:HD12	1.80	0.47
2:E:13:GLU:HB2	2:E:49:LYS:HE3	1.96	0.47
1:A:41:LEU:HD12	1:A:41:LEU:N	2.29	0.47
1:A:185:LYS:HG2	1:A:186:LEU:H	1.80	0.47
1:B:7:GLU:O	1:B:13:ILE:HD12	2.14	0.47
2:D:69:TYR:O	2:D:73:ILE:HG13	2.15	0.47
2:F:6:VAL:CG2	2:F:62:LEU:HD22	2.44	0.47
2:F:23:VAL:HG11	2:F:73:ILE:HG23	1.97	0.47
1:A:2:ILE:O	1:A:70:ASP:HB2	2.15	0.47
1:A:9:THR:HA	1:A:84:LEU:HD23	1.97	0.47
1:B:175:GLU:O	1:B:179:LEU:HB2	2.14	0.47
1:A:77:GLY:HA2	1:A:315:PHE:HD1	1.80	0.46
1:A:196:MET:CG	1:A:238:ARG:HG2	2.41	0.46
2:F:49:LYS:HB2	2:F:49:LYS:HZ2	1.77	0.46
1:A:325:ILE:HG22	1:A:325:ILE:O	2.16	0.46
1:B:78:PRO:HD3	1:B:315:PHE:HB3	1.98	0.46
1:B:208:LEU:CD1	2:D:35:PRO:HB3	2.41	0.46
1:A:78:PRO:HG3	1:A:315:PHE:O	2.14	0.46
1:A:189:LEU:HB2	1:A:227:THR:HG21	1.97	0.46
2:E:14:PHE:CD1	2:E:77:ILE:HD11	2.51	0.46
1:A:126:MET:HG2	1:A:137:ILE:HD11	1.98	0.46
1:A:152:ASP:OD2	1:A:199:ALA:HB1	2.15	0.46
1:B:70:ASP:O	1:B:100:LYS:HB3	2.15	0.46
1:B:176:ILE:HG23	1:B:221:SER:HB2	1.97	0.46
2:C:45:LEU:HD23	2:C:46:GLU:H	1.81	0.46
1:B:148:GLY:O	1:B:149:GLU:HB3	2.15	0.46
2:C:24:TYR:O	2:C:28:LEU:HB2	2.16	0.46
2:F:56:ALA:HB3	2:F:62:LEU:HB2	1.98	0.46
2:E:25:GLU:O	2:E:29:TYR:HD1	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:9:LYS:HA	2:F:52:LEU:O	2.15	0.46
1:A:212:LYS:C	1:A:214:GLY:N	2.69	0.46
2:E:10:ILE:N	2:E:10:ILE:HD12	2.30	0.46
1:B:240:LEU:C	1:B:242:VAL:H	2.19	0.46
2:C:14:PHE:CE1	2:C:77:ILE:HD11	2.51	0.46
2:E:42:ASP:HB3	2:E:53:ASP:HB2	1.97	0.46
1:A:18:ILE:HG22	1:A:23:ILE:HD12	1.98	0.45
1:B:124:PRO:HA	1:B:247:GLU:HG3	1.97	0.45
2:E:24:TYR:CE2	2:E:28:LEU:HG	2.51	0.45
1:B:19:ASP:OD1	1:B:19:ASP:C	2.54	0.45
1:B:61:LYS:HB2	1:B:61:LYS:NZ	2.31	0.45
2:C:5:ARG:N	2:C:5:ARG:CD	2.79	0.45
1:A:206:ALA:O	1:A:210:TYR:CD2	2.69	0.45
1:B:213:THR:HG22	1:B:213:THR:O	2.17	0.45
2:E:7:GLN:HE22	2:E:55:LYS:CE	2.26	0.45
1:A:307:GLU:HA	2:C:29:TYR:HE2	1.80	0.45
2:E:18:ASP:O	2:E:21:LYS:HB2	2.17	0.45
1:B:106:ASN:HB2	1:B:315:PHE:CE1	2.52	0.45
1:B:170:PHE:N	1:B:171:PRO:CD	2.80	0.45
1:A:143:ARG:CG	1:A:321:ASP:HA	2.43	0.45
1:B:152:ASP:O	2:D:29:TYR:CE2	2.70	0.45
1:B:189:LEU:HA	1:B:190:PRO:HD3	1.85	0.45
1:A:23:ILE:HG13	1:A:288:ILE:HD13	1.98	0.45
1:B:23:ILE:HG13	1:B:288:ILE:CD1	2.47	0.45
2:F:34:VAL:HG21	2:F:37:ARG:NH2	2.32	0.45
1:A:170:PHE:N	1:A:171:PRO:CD	2.80	0.45
1:A:129:VAL:HB	1:A:254:VAL:HB	1.99	0.44
1:A:158:MET:HG2	1:A:204:LEU:HB2	1.99	0.44
1:B:236:LEU:HD23	1:B:236:LEU:HA	1.79	0.44
1:B:246:ASP:C	1:B:273:ILE:HB	2.38	0.44
1:B:100:LYS:HB3	1:B:101:PRO:HD2	1.98	0.44
1:A:55:ILE:O	1:A:59:LEU:HG	2.17	0.44
1:A:105:VAL:HG12	1:A:106:ASN:N	2.32	0.44
2:D:5:ARG:HD2	2:D:5:ARG:N	2.33	0.44
1:A:61:LYS:HB2	1:A:61:LYS:NZ	2.32	0.44
1:A:76:MET:HE2	1:A:310:ALA:O	2.17	0.44
1:A:116:ARG:HH12	1:A:325:ILE:HG12	1.83	0.44
1:B:169:PRO:C	1:B:171:PRO:HD2	2.38	0.44
1:B:307:GLU:HG2	1:B:308:GLU:N	2.32	0.44
2:D:49:LYS:HB2	2:D:49:LYS:HZ2	1.82	0.44
2:D:66:VAL:O	2:D:70:LEU:HB2	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:136:VAL:HG13	1:A:136:VAL:O	2.16	0.44
1:A:317:ILE:N	1:A:317:ILE:CD1	2.81	0.44
1:B:41:LEU:N	1:B:41:LEU:HD12	2.32	0.44
2:D:6:VAL:CG2	2:D:62:LEU:HD22	2.46	0.44
1:A:5:GLY:O	1:A:15:CYS:HA	2.17	0.44
1:A:265:THR:O	1:A:268:ALA:HB3	2.17	0.44
2:C:34:VAL:HG21	2:C:37:ARG:NH2	2.33	0.44
1:A:186:LEU:HD23	1:A:223:SER:OG	2.16	0.44
1:A:193:VAL:HG23	1:A:231:MET:HG2	2.00	0.44
1:B:247:GLU:O	1:B:248:ILE:HD12	2.18	0.44
2:F:74:LYS:HE2	2:F:78:ASP:OD2	2.18	0.44
1:B:72:ILE:O	1:B:102:ILE:HG13	2.17	0.43
1:B:114:ILE:HG23	1:B:115:GLY:N	2.32	0.43
1:B:206:ALA:O	1:B:210:TYR:CD2	2.71	0.43
2:F:46:GLU:O	2:F:49:LYS:N	2.51	0.43
2:C:49:LYS:HB2	2:C:49:LYS:HZ2	1.83	0.43
2:E:54:ILE:N	2:E:54:ILE:HD12	2.33	0.43
1:A:3:VAL:HG22	1:A:71:LEU:HB3	1.99	0.43
1:B:14:SER:HB3	1:B:287:GLY:N	2.33	0.43
1:B:41:LEU:O	1:B:45:VAL:HG23	2.18	0.43
2:E:6:VAL:HG21	2:E:62:LEU:HB2	2.00	0.43
2:E:74:LYS:HD2	2:F:60:SER:HB2	2.00	0.43
1:A:59:LEU:HG	1:A:59:LEU:H	1.64	0.43
1:A:186:LEU:HD23	1:A:223:SER:CB	2.49	0.43
1:B:200:PHE:HA	1:B:203:ILE:HD13	1.98	0.43
2:C:43:PHE:HE1	2:C:50:ILE:CG2	2.31	0.43
2:D:16:SER:OG	2:D:19:VAL:HG23	2.18	0.43
1:A:54:VAL:HG23	1:A:55:ILE:N	2.34	0.43
1:A:273:ILE:HA	1:A:274:ARG:HH21	1.83	0.43
1:B:315:PHE:CZ	1:B:320:VAL:HG23	2.54	0.43
2:C:67:ASN:O	2:C:71:ARG:HB2	2.18	0.43
1:A:14:SER:HB2	1:A:28:SER:OG	2.19	0.43
1:A:307:GLU:HG2	1:A:308:GLU:N	2.33	0.43
1:B:12:THR:HG23	1:B:28:SER:HB3	2.01	0.43
1:B:46:HIS:O	1:B:50:VAL:HG22	2.18	0.43
2:E:34:VAL:HA	2:E:35:PRO:HD3	1.85	0.43
1:A:182:LYS:O	1:A:182:LYS:CG	2.66	0.43
1:A:236:LEU:O	1:A:237:GLU:C	2.56	0.43
1:A:249:LEU:HD22	1:A:278:THR:CG2	2.48	0.43
1:A:176:ILE:CG2	1:A:221:SER:HB2	2.49	0.43
2:C:36:TYR:CE1	2:E:19:VAL:HG22	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:10:ILE:HD12	2:F:10:ILE:N	2.34	0.43
1:A:169:PRO:C	1:A:171:PRO:HD2	2.39	0.43
1:A:236:LEU:HD23	1:A:236:LEU:HA	1.75	0.43
1:B:31:TYR:HB2	1:B:46:HIS:CG	2.54	0.43
1:B:177:GLU:C	1:B:179:LEU:H	2.22	0.43
2:C:41:ILE:N	2:C:41:ILE:HD12	2.33	0.43
2:F:41:ILE:HD12	2:F:41:ILE:N	2.34	0.43
1:A:123:ASP:O	1:A:245:LYS:HD3	2.19	0.43
1:A:203:ILE:HD12	1:A:203:ILE:H	1.84	0.43
1:B:240:LEU:HD21	1:B:248:ILE:HD11	1.98	0.43
2:D:10:ILE:N	2:D:10:ILE:HD12	2.34	0.43
2:E:19:VAL:O	2:E:23:VAL:HG23	2.19	0.43
2:E:27:VAL:HB	2:E:43:PHE:HZ	1.84	0.43
1:B:286:ASN:O	1:B:289:MET:HG2	2.19	0.42
1:A:89:THR:O	1:A:93:THR:CG2	2.68	0.42
1:A:286:ASN:HD22	1:A:289:MET:H	1.64	0.42
1:B:152:ASP:OD2	1:B:199:ALA:HB1	2.19	0.42
2:C:10:ILE:HD12	2:C:10:ILE:N	2.34	0.42
2:D:14:PHE:CD1	2:D:77:ILE:HD11	2.55	0.42
1:B:93:THR:HG21	2:F:76:ALA:HA	2.00	0.42
1:B:128:TYR:CD2	1:B:135:GLN:HG3	2.54	0.42
1:B:128:TYR:HD2	1:B:135:GLN:HG3	1.84	0.42
1:B:182:LYS:O	1:B:182:LYS:CG	2.67	0.42
1:A:187:LEU:HB3	1:A:210:TYR:CE1	2.54	0.42
1:B:9:THR:HA	1:B:84:LEU:HD23	2.01	0.42
1:B:187:LEU:HB3	1:B:210:TYR:CE1	2.53	0.42
2:C:38:ARG:HB3	2:E:81:GLU:HB3	2.02	0.42
1:B:36:GLY:HA2	1:B:170:PHE:CE2	2.55	0.42
1:B:286:ASN:ND2	1:B:289:MET:N	2.59	0.42
2:C:13:GLU:HB2	2:C:49:LYS:HE3	2.00	0.42
2:C:23:VAL:HG11	2:C:73:ILE:HG23	2.01	0.42
2:E:76:ALA:O	2:E:77:ILE:C	2.58	0.42
1:A:3:VAL:CG1	1:A:4:LEU:N	2.83	0.42
1:A:54:VAL:HG23	1:A:55:ILE:H	1.85	0.42
2:F:31:HIS:HD2	2:F:43:PHE:HB2	1.84	0.42
1:A:180:ALA:O	1:A:222:TYR:HD1	2.02	0.42
2:C:72:TRP:O	2:C:75:ALA:HB3	2.19	0.42
1:A:1:MET:SD	1:A:1:MET:C	2.98	0.42
1:A:246:ASP:HB2	1:A:273:ILE:HG22	2.02	0.42
1:A:226:GLU:O	1:A:230:ALA:HB2	2.20	0.42
2:E:14:PHE:HD1	2:E:77:ILE:HD11	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:24:LEU:H	1:A:24:LEU:HD12	1.84	0.41
1:A:47:HIS:O	1:A:49:GLU:N	2.53	0.41
2:D:44:LYS:HE3	2:D:44:LYS:HB2	1.67	0.41
1:A:324:TRP:O	1:A:324:TRP:CD1	2.73	0.41
1:B:27:GLU:OE2	1:B:57:ARG:HB3	2.21	0.41
1:B:273:ILE:HA	1:B:274:ARG:HH21	1.84	0.41
1:A:92:ARG:NH2	2:C:30:GLU:OE1	2.52	0.41
1:A:189:LEU:HA	1:A:190:PRO:HD3	1.86	0.41
1:A:238:ARG:O	1:A:242:VAL:HG23	2.20	0.41
2:E:34:VAL:HG11	2:E:37:ARG:HH21	1.84	0.41
1:B:47:HIS:O	1:B:51:ILE:HB	2.20	0.41
1:B:110:GLY:O	1:B:289:MET:HB2	2.20	0.41
2:C:70:LEU:HD23	2:C:70:LEU:HA	1.77	0.41
1:A:7:GLU:HB3	1:A:287:GLY:CA	2.48	0.41
1:A:177:GLU:C	1:A:179:LEU:H	2.24	0.41
1:A:193:VAL:HG22	1:A:198:THR:HG22	2.03	0.41
2:E:67:ASN:HB3	2:E:71:ARG:NH2	2.35	0.41
1:B:23:ILE:HG13	1:B:288:ILE:HD13	2.01	0.41
1:B:193:VAL:HG22	1:B:198:THR:HG22	2.03	0.41
1:A:128:TYR:CD2	1:A:135:GLN:HG3	2.55	0.41
1:A:180:ALA:CB	1:A:222:TYR:HA	2.48	0.41
1:B:140:VAL:HG12	1:B:141:ASN:N	2.36	0.41
1:A:18:ILE:CG2	1:A:23:ILE:HD12	2.50	0.41
1:A:31:TYR:CE2	1:A:33:PRO:HG3	2.56	0.41
1:A:129:VAL:O	1:A:254:VAL:HB	2.21	0.41
1:B:63:LYS:HA	1:B:63:LYS:HD3	1.88	0.41
2:C:11:GLU:HG2	2:C:51:ILE:HD12	2.01	0.41
1:A:15:CYS:SG	1:A:55:ILE:HA	2.60	0.40
1:B:139:HIS:HB2	1:B:144:TYR:CE1	2.57	0.40
1:B:180:ALA:CB	1:B:222:TYR:HA	2.49	0.40
2:D:54:ILE:HD12	2:D:54:ILE:N	2.36	0.40
2:F:14:PHE:HD1	2:F:77:ILE:HD11	1.81	0.40
1:A:216:ALA:O	1:A:220:ILE:HG13	2.21	0.40
1:B:170:PHE:C	1:B:172:GLY:N	2.74	0.40
1:A:3:VAL:HG12	1:A:4:LEU:N	2.34	0.40
1:A:7:GLU:O	1:A:13:ILE:HD12	2.21	0.40
2:D:28:LEU:HD22	2:D:28:LEU:HA	1.88	0.40
1:A:19:ASP:OD1	1:A:19:ASP:C	2.60	0.40
1:A:116:ARG:HH12	1:A:325:ILE:HG13	1.86	0.40
1:B:155:ILE:HG23	1:B:254:VAL:HG21	2.03	0.40
2:C:31:HIS:CE1	2:C:37:ARG:HH12	2.40	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:14:PHE:CD1	2:F:3:ALA:N	2.89	0.40
1:A:72:ILE:HD12	1:A:95:SER:CA	2.48	0.40
1:B:105:VAL:CG1	1:B:106:ASN:N	2.84	0.40
1:B:240:LEU:HD21	1:B:248:ILE:HD12	2.01	0.40
2:D:18:ASP:O	2:D:21:LYS:HB2	2.21	0.40
2:D:66:VAL:HG13	2:D:70:LEU:HD12	2.03	0.40
2:F:4:LYS:HB3	2:F:4:LYS:NZ	2.36	0.40
2:F:62:LEU:O	2:F:66:VAL:HG23	2.21	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	327/334 (98%)	271 (83%)	49 (15%)	7 (2%)	7	31
1	B	326/334 (98%)	272 (83%)	44 (14%)	10 (3%)	4	21
2	C	75/87 (86%)	68 (91%)	5 (7%)	2 (3%)	5	25
2	D	74/87 (85%)	63 (85%)	8 (11%)	3 (4%)	3	15
2	E	75/87 (86%)	64 (85%)	8 (11%)	3 (4%)	3	16
2	F	77/87 (88%)	62 (80%)	11 (14%)	4 (5%)	2	11
All	All	954/1016 (94%)	800 (84%)	125 (13%)	29 (3%)	4	22

All (29) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	35	THR
1	A	48	SER
2	D	47	GLY
2	F	68	SER

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Mol	Chain	Res	Type
1	A	173	GLY
2	C	47	GLY
2	D	76	ALA
2	E	47	GLY
2	E	76	ALA
2	F	35	PRO
1	A	26	MET
1	B	61	LYS
1	B	167	GLY
1	B	244	GLY
2	C	35	PRO
2	D	77	ILE
2	F	39	SER
1	A	167	GLY
1	B	182	LYS
1	B	240	LEU
2	E	77	ILE
1	A	77	GLY
1	B	77	GLY
1	A	190	PRO
1	B	173	GLY
2	F	47	GLY
1	B	102	ILE
1	B	190	PRO
1	B	8	GLY

### 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	265/268 (99%)	225 (85%)	40 (15%)	3	13
1	B	264/268 (98%)	233 (88%)	31 (12%)	5	21
2	C	68/75 (91%)	61 (90%)	7 (10%)	7	26
2	D	67/75 (89%)	58 (87%)	9 (13%)	4	16
2	E	68/75 (91%)	58 (85%)	10 (15%)	3	14

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	F	69/75 (92%)	59 (86%)	10 (14%)	3	14
All	All	801/836 (96%)	694 (87%)	107 (13%)	4	16

All (107) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-2	MET
1	A	-1	ASP
1	A	1	MET
1	A	21	SER
1	A	22	ARG
1	A	35	THR
1	A	39	ARG
1	A	41	LEU
1	A	49	GLU
1	A	51	ILE
1	A	53	THR
1	A	55	ILE
1	A	61	LYS
1	A	63	LYS
1	A	71	LEU
1	A	87	THR
1	A	93	THR
1	A	109	LEU
1	A	126	MET
1	A	130	SER
1	A	135	GLN
1	A	137	ILE
1	A	152	ASP
1	A	175	GLU
1	A	179	LEU
1	A	186	LEU
1	A	197	ASP
1	A	211	LEU
1	A	212	LYS
1	A	217	ILE
1	A	219	ASP
1	A	226	GLU
1	A	236	LEU
1	A	238	ARG
1	A	247	GLU
1	A	269	ARG

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Mol	Chain	Res	Type
1	A	270	GLU
1	A	274	ARG
1	A	305	SER
1	A	321	ASP
1	B	-1	ASP
1	B	22	ARG
1	B	39	ARG
1	B	41	LEU
1	B	49	GLU
1	B	51	ILE
1	B	53	THR
1	B	55	ILE
1	B	61	LYS
1	B	63	LYS
1	B	71	LEU
1	B	93	THR
1	B	109	LEU
1	B	130	SER
1	B	135	GLN
1	B	137	ILE
1	B	175	GLU
1	B	179	LEU
1	B	182	LYS
1	B	186	LEU
1	B	197	ASP
1	B	211	LEU
1	B	212	LYS
1	B	217	ILE
1	B	219	ASP
1	B	226	GLU
1	B	238	ARG
1	B	269	ARG
1	B	270	GLU
1	B	274	ARG
1	B	321	ASP
2	C	5	ARG
2	C	12	MET
2	C	34	VAL
2	C	37	ARG
2	C	38	ARG
2	C	40	GLU
2	C	43	PHE

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Mol	Chain	Res	Type
2	D	5	ARG
2	D	6	VAL
2	D	12	MET
2	D	18	ASP
2	D	28	LEU
2	D	38	ARG
2	D	40	GLU
2	D	43	PHE
2	D	46	GLU
2	E	5	ARG
2	E	12	MET
2	E	18	ASP
2	E	28	LEU
2	E	38	ARG
2	E	40	GLU
2	E	41	ILE
2	E	43	PHE
2	E	46	GLU
2	E	60	SER
2	F	5	ARG
2	F	6	VAL
2	F	12	MET
2	F	34	VAL
2	F	37	ARG
2	F	38	ARG
2	F	40	GLU
2	F	43	PHE
2	F	46	GLU
2	F	81	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	HIS
1	A	135	GLN
1	A	139	HIS
1	A	225	GLN
1	A	286	ASN
1	B	11	HIS
1	B	107	HIS
1	B	135	GLN
1	B	139	HIS

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Mol	Chain	Res	Type
1	B	225	GLN
1	B	286	ASN
2	C	7	GLN
2	C	31	HIS
2	D	7	GLN
2	E	7	GLN
2	E	31	HIS
2	F	7	GLN
2	F	31	HIS

### 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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 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

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	329/334 (98%)	-0.09	0 <span>100</span> <span>100</span>	56, 79, 108, 121	0
1	B	328/334 (98%)	-0.09	2 (0%) <span>89</span> <span>72</span>	49, 78, 107, 120	0
2	C	77/87 (88%)	-0.01	0 <span>100</span> <span>100</span>	67, 86, 102, 117	0
2	D	76/87 (87%)	0.11	1 (1%) <span>77</span> <span>51</span>	65, 90, 111, 133	0
2	E	77/87 (88%)	-0.05	0 <span>100</span> <span>100</span>	69, 89, 110, 130	0
2	F	79/87 (90%)	-0.03	0 <span>100</span> <span>100</span>	65, 88, 101, 114	0
All	All	966/1016 (95%)	-0.06	3 (0%) <span>94</span> <span>83</span>	49, 82, 108, 133	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	34	LYS	3.2
2	D	7	GLN	3.0
1	B	72	ILE	2.0

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

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

### 6.3 Carbohydrates [i](#)

There are no monosaccharides 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. 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	B-factors(Å <sup>2</sup> )	Q<0.9
3	MG	A	600	1/1	0.90	0.29	47,47,47,47	0
3	MG	B	600	1/1	0.96	0.15	73,73,73,73	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.