



# Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2020 – 05:06 am BST

PDB ID : 1F9N  
Title : CRYSTAL STRUCTURE OF AHRC, THE ARGININE REPRESSOR/ACTIVATOR PROTEIN FROM BACILLUS SUBTILIS  
Authors : Dennis, C.A.; Glykos, N.M.; Parsons, M.R.; Phillips, S.E.V.  
Deposited on : 2000-07-11  
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

<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  
Xtrriage (Phenix) : 1.13  
EDS : 2.11  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

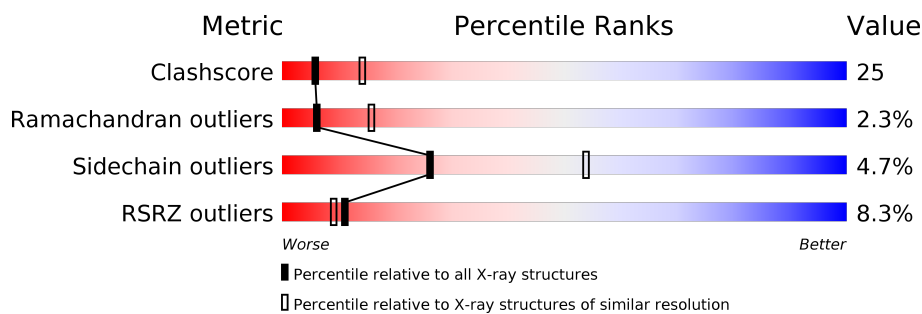
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.70 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))
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (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 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	149	<div style="display: flex; align-items: center;"> <div style="width: 8%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 49%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 49%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: grey; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">8%      49%      49%      ..</p>
1	B	149	<div style="display: flex; align-items: center;"> <div style="width: 3%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 54%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 42%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 5%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">3%      54%      42%      5%</p>
1	C	149	<div style="display: flex; align-items: center;"> <div style="width: 20%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 51%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 39%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 8%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: grey; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">20%      51%      39%      8%      ..</p>
1	D	149	<div style="display: flex; align-items: center;"> <div style="width: 0%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 58%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 36%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: grey; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">%      58%      36%      . .</p>
1	E	149	<div style="display: flex; align-items: center;"> <div style="width: 0%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 63%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 35%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: grey; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">%      63%      35%      ..</p>
1	F	149	<div style="display: flex; align-items: center;"> <div style="width: 16%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 54%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 42%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">16%      54%      42%      .</p>

## 2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 7108 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 ARGININE REPRESSOR/ACTIVATOR PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	147	Total 1143	C 715	N 195	O 224	S 9	0	0	0
1	B	149	Total 1157	C 725	N 196	O 226	S 10	0	0	0
1	C	148	Total 1132	C 708	N 193	O 222	S 9	0	0	0
1	D	146	Total 1124	C 703	N 189	O 223	S 9	0	0	0
1	E	148	Total 1159	C 725	N 199	O 226	S 9	0	0	0
1	F	149	Total 1153	C 724	N 195	O 224	S 10	0	0	0

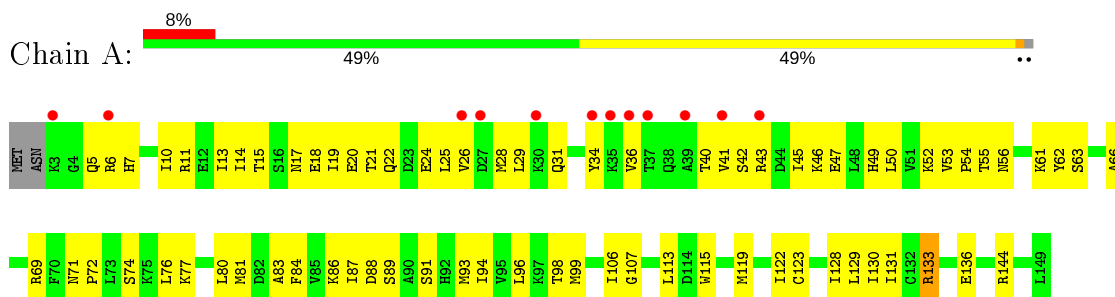
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	21	Total 21	O 21	0	0
2	B	56	Total 56	O 56	0	0
2	C	34	Total 34	O 34	0	0
2	D	36	Total 36	O 36	0	0
2	E	60	Total 60	O 60	0	0
2	F	33	Total 33	O 33	0	0

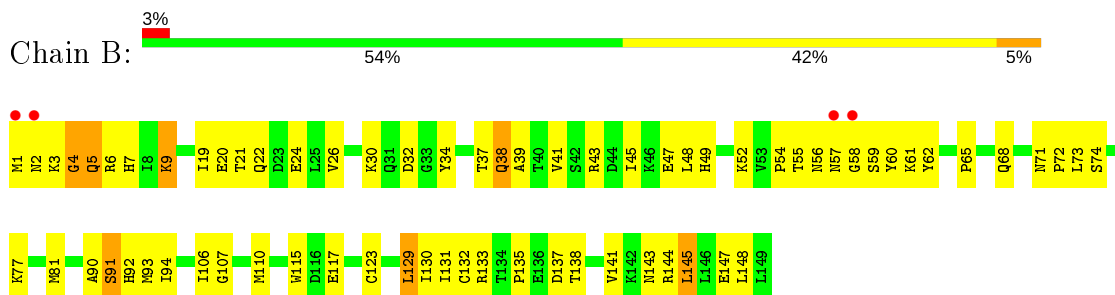
### 3 Residue-property plots [i](#)

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

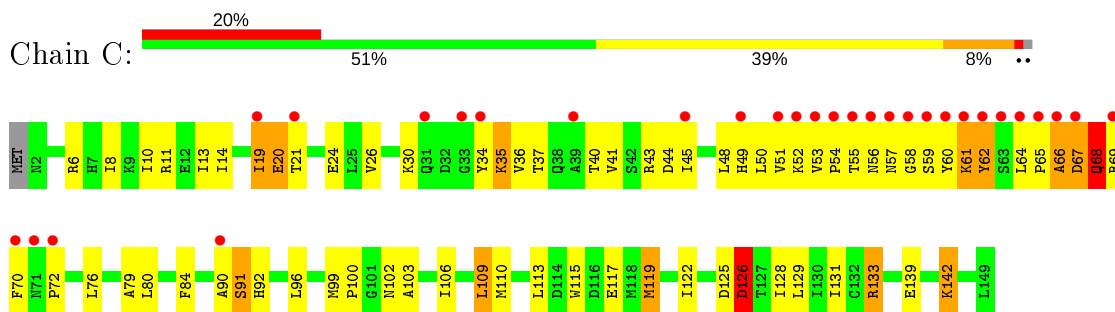
- Molecule 1: ARGININE REPRESSOR/ACTIVATOR PROTEIN



- Molecule 1: ARGININE REPRESSOR/ACTIVATOR PROTEIN

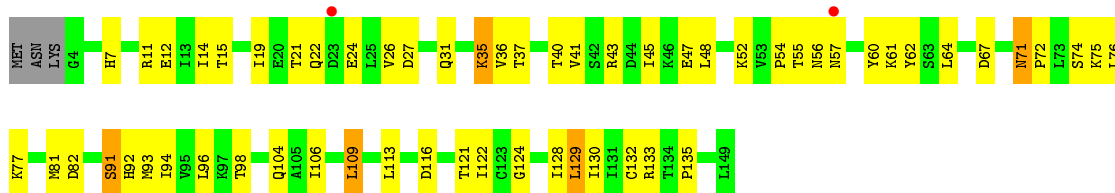


- Molecule 1: ARGININE REPRESSOR/ACTIVATOR PROTEIN

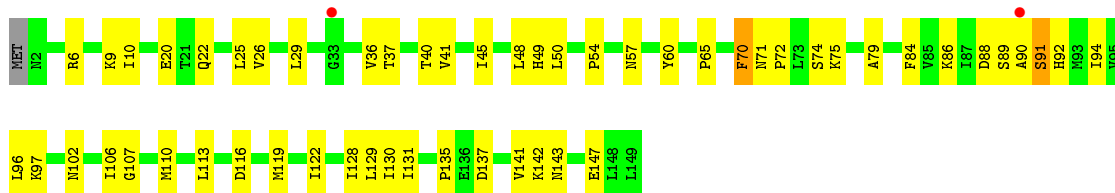


- Molecule 1: ARGININE REPRESSOR/ACTIVATOR PROTEIN

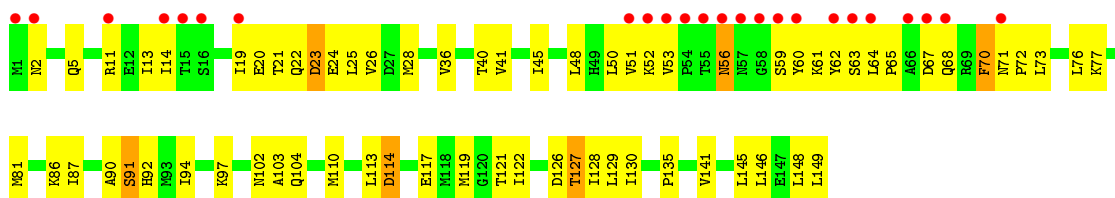




• Molecule 1: ARGININE REPRESSOR/ACTIVATOR PROTEIN



• Molecule 1: ARGININE REPRESSOR/ACTIVATOR PROTEIN



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	230.31Å 73.86Å 138.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.70 34.73 – 2.70	Depositor EDS
% Data completeness (in resolution range)	88.7 (30.00-2.70) 85.0 (34.73-2.70)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.36 (at 2.68Å)	Xtrriage
Refinement program	CNS 0.9	Depositor
R, $R_{free}$	0.220 , 0.270 0.215 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	56.4	Xtrriage
Anisotropy	0.044	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 87.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	7108	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	66.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.40% 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	0.26	0/1157	0.52	0/1563
1	B	0.24	0/1171	0.50	0/1580
1	C	0.29	0/1145	0.58	0/1548
1	D	0.24	0/1138	0.49	0/1541
1	E	0.26	0/1173	0.52	0/1582
1	F	0.37	0/1167	0.61	1/1574 (0.1%)
All	All	0.28	0/6951	0.54	1/9388 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	56	ASN	CA-CB-CG	-5.83	100.57	113.40

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	1143	0	1153	56	1
1	B	1157	0	1176	59	0
1	C	1132	0	1135	88	0
1	D	1124	0	1118	42	0
1	E	1159	0	1181	40	0
1	F	1153	0	1169	76	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	21	0	0	0	0
2	B	56	0	0	1	0
2	C	34	0	0	0	0
2	D	36	0	0	0	0
2	E	60	0	0	1	0
2	F	33	0	0	2	0
All	All	7108	0	6932	347	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 25.

All (347) 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:C:48:LEU:O	1:C:65:PRO:HG2	1.64	0.97
1:F:56:ASN:ND2	2:F:171:HOH:O	1.98	0.95
1:F:51:VAL:HG12	1:F:52:LYS:N	1.82	0.93
1:C:62:TYR:HD1	1:C:62:TYR:H	1.04	0.93
1:C:62:TYR:N	1:C:62:TYR:CD1	2.34	0.90
1:B:26:VAL:HG22	1:B:41:VAL:HG21	1.54	0.90
1:E:57:ASN:OD1	2:E:200:HOH:O	1.89	0.90
1:A:25:LEU:O	1:A:29:LEU:HG	1.70	0.90
1:F:11:ARG:O	1:F:14:ILE:HG22	1.73	0.87
1:A:77:LYS:HG2	1:A:81:MET:HE2	1.57	0.87
1:B:93:MET:HG2	1:B:129:LEU:HD21	1.57	0.85
1:F:14:ILE:HD11	1:F:62:TYR:HB3	1.58	0.85
1:F:20:GLU:HG3	1:F:60:TYR:HB2	1.57	0.84
1:C:51:VAL:HG12	1:C:52:LYS:H	1.41	0.84
1:F:51:VAL:CG1	1:F:52:LYS:N	2.40	0.82
1:F:21:THR:HG22	1:F:23:ASP:H	1.42	0.82
1:F:26:VAL:HG22	1:F:41:VAL:HG21	1.62	0.81
1:F:51:VAL:CG1	1:F:52:LYS:H	1.94	0.80
1:F:70:PHE:HD1	1:F:71:ASN:N	1.80	0.79
1:B:71:ASN:ND2	1:B:74:SER:HB3	1.98	0.78
1:A:7:HIS:O	1:A:11:ARG:HG3	1.84	0.77
1:B:55:THR:HG22	1:B:56:ASN:H	1.49	0.77
1:F:52:LYS:O	1:F:53:VAL:CG2	2.33	0.77
1:B:1:MET:HG3	1:B:3:LYS:HE2	1.65	0.76
1:C:51:VAL:HG12	1:C:52:LYS:N	2.01	0.75
1:C:53:VAL:HG13	1:C:54:PRO:HD2	1.68	0.75
1:F:86:LYS:HG2	1:F:87:ILE:N	1.99	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:70:PHE:CD1	1:C:70:PHE:O	2.40	0.75
1:C:53:VAL:CG1	1:C:54:PRO:HD2	2.17	0.75
1:A:94:ILE:CG2	1:A:130:ILE:HB	2.17	0.74
1:C:62:TYR:N	1:C:62:TYR:HD1	1.79	0.74
1:D:35:LYS:HA	1:D:35:LYS:HE3	1.70	0.74
1:C:142:LYS:NZ	1:C:142:LYS:HB2	2.02	0.74
1:A:18:GLU:H	1:D:57:ASN:HB2	1.53	0.73
1:C:53:VAL:HB	1:C:61:LYS:HD2	1.71	0.73
1:F:19:ILE:O	1:F:61:LYS:HA	1.89	0.72
1:A:94:ILE:HG23	1:A:130:ILE:HB	1.72	0.72
1:D:26:VAL:HG22	1:D:41:VAL:HG21	1.73	0.70
1:F:86:LYS:HG2	1:F:87:ILE:H	1.54	0.70
1:F:14:ILE:HD11	1:F:62:TYR:CB	2.22	0.69
1:D:93:MET:HE3	1:D:129:LEU:HD21	1.74	0.69
1:A:87:ILE:HD11	1:A:94:ILE:HD11	1.75	0.69
1:B:92:HIS:CE1	1:B:135:PRO:HD3	2.27	0.69
1:B:110:MET:HE1	1:B:145:LEU:HD21	1.75	0.68
1:C:76:LEU:HA	1:C:109:LEU:HD23	1.76	0.68
1:E:110:MET:HE1	1:E:128:ILE:HG21	1.75	0.68
1:C:49:HIS:O	1:C:65:PRO:HG3	1.93	0.68
1:F:52:LYS:C	1:F:53:VAL:HG23	2.14	0.68
1:E:88:ASP:OD2	1:F:119:MET:HG2	1.94	0.68
1:F:52:LYS:O	1:F:53:VAL:HG22	1.93	0.67
1:C:19:ILE:H	1:C:19:ILE:HD12	1.60	0.67
1:F:48:LEU:HB2	1:F:50:LEU:HG	1.77	0.66
1:F:64:LEU:HG	1:F:65:PRO:HD2	1.77	0.66
1:D:19:ILE:HG22	1:D:62:TYR:HD2	1.61	0.66
1:C:35:LYS:N	1:C:35:LYS:HD3	2.10	0.66
1:C:64:LEU:HG	1:C:65:PRO:HD2	1.76	0.66
1:C:66:ALA:O	1:C:68:GLN:N	2.27	0.66
1:B:57:ASN:ND2	1:B:58:GLY:H	1.93	0.66
1:F:21:THR:HB	1:F:24:GLU:HG3	1.78	0.66
1:D:11:ARG:NH1	1:D:64:LEU:HD21	2.12	0.65
1:B:43:ARG:O	1:B:47:GLU:HG3	1.96	0.65
1:C:45:ILE:HG23	1:C:50:LEU:HB2	1.77	0.65
1:B:71:ASN:N	1:B:72:PRO:HD3	2.11	0.65
1:F:51:VAL:HG12	1:F:52:LYS:O	1.97	0.65
1:A:36:VAL:HG13	1:A:40:THR:HB	1.78	0.65
1:F:14:ILE:HD11	1:F:62:TYR:C	2.17	0.64
1:C:19:ILE:HD12	1:C:19:ILE:N	2.11	0.64
1:F:76:LEU:HD11	1:F:110:MET:HG2	1.78	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:110:MET:HE1	1:E:128:ILE:HD13	1.80	0.64
1:C:35:LYS:HD3	1:C:35:LYS:H	1.63	0.64
1:B:4:GLY:HA2	2:B:154:HOH:O	1.97	0.64
1:B:71:ASN:HD21	1:B:74:SER:HB3	1.61	0.64
1:E:26:VAL:HG22	1:E:41:VAL:HG21	1.79	0.64
1:C:142:LYS:HZ2	1:C:142:LYS:HB2	1.63	0.64
1:C:65:PRO:C	1:C:67:ASP:H	2.00	0.63
1:F:13:ILE:HG21	1:F:28:MET:HG2	1.79	0.63
1:B:77:LYS:O	1:B:81:MET:HG3	1.98	0.63
1:F:76:LEU:CD1	1:F:113:LEU:HD11	2.28	0.63
1:E:71:ASN:HB3	1:E:74:SER:OG	1.99	0.63
1:F:51:VAL:HG12	1:F:52:LYS:H	1.58	0.62
1:F:13:ILE:HD13	1:F:28:MET:HG3	1.82	0.62
1:B:55:THR:HG22	1:B:56:ASN:N	2.14	0.62
1:F:67:ASP:HB3	1:F:70:PHE:CD2	2.35	0.62
1:F:70:PHE:HD1	1:F:71:ASN:H	1.45	0.62
1:B:94:ILE:HB	1:B:130:ILE:HB	1.82	0.62
1:A:119:MET:HE3	1:A:131:ILE:HG22	1.81	0.61
1:B:20:GLU:OE1	1:B:59:SER:HB2	2.01	0.61
1:F:13:ILE:CG2	1:F:28:MET:HG2	2.31	0.61
1:D:21:THR:OG1	1:D:24:GLU:HG3	1.99	0.61
1:C:53:VAL:CG1	1:C:54:PRO:CD	2.79	0.61
1:C:102:ASN:O	1:C:106:ILE:HG12	2.01	0.60
1:C:52:LYS:HD3	1:C:60:TYR:CB	2.31	0.60
1:F:103:ALA:HB3	2:F:154:HOH:O	2.01	0.60
1:F:52:LYS:C	1:F:53:VAL:CG2	2.69	0.60
1:D:54:PRO:HG3	1:D:60:TYR:CE2	2.36	0.59
1:F:67:ASP:HB3	1:F:70:PHE:CE2	2.37	0.59
1:B:133:ARG:HH11	1:B:133:ARG:HG2	1.67	0.59
1:B:30:LYS:HA	1:B:34:TYR:O	2.03	0.59
1:C:19:ILE:HG23	1:C:24:GLU:HB2	1.84	0.59
1:A:5:GLN:HA	1:A:5:GLN:OE1	2.03	0.58
1:C:14:ILE:HD13	1:C:19:ILE:HD13	1.85	0.58
1:C:51:VAL:CG1	1:C:52:LYS:H	2.12	0.58
1:C:113:LEU:HD13	1:C:115:TRP:CZ2	2.39	0.58
1:C:125:ASP:O	1:C:126:ASP:CB	2.51	0.58
1:D:7:HIS:HB3	1:D:48:LEU:HD21	1.84	0.58
1:C:65:PRO:O	1:C:67:ASP:N	2.36	0.58
1:B:49:HIS:O	1:B:65:PRO:HG3	2.04	0.57
1:D:43:ARG:O	1:D:47:GLU:HG3	2.03	0.57
1:E:106:ILE:O	1:E:110:MET:HG3	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:70:PHE:O	1:E:72:PRO:HD3	2.05	0.57
1:A:6:ARG:HB3	1:A:34:TYR:CE2	2.39	0.57
1:F:22:GLN:O	1:F:26:VAL:HG23	2.04	0.57
1:C:19:ILE:HG22	1:C:19:ILE:O	2.05	0.57
1:F:41:VAL:O	1:F:45:ILE:HD13	2.03	0.57
1:B:115:TRP:CH2	1:B:144:ARG:HD2	2.39	0.57
1:B:20:GLU:CD	1:B:59:SER:HB2	2.25	0.57
1:A:6:ARG:O	1:A:10:ILE:HG13	2.04	0.57
1:D:27:ASP:O	1:D:31:GLN:HG3	2.04	0.57
1:F:14:ILE:HD12	1:F:50:LEU:HD13	1.87	0.57
1:F:92:HIS:CE1	1:F:135:PRO:HD3	2.40	0.57
1:E:70:PHE:C	1:E:72:PRO:HD3	2.25	0.56
1:E:89:SER:OG	1:E:142:LYS:HD2	2.05	0.56
1:F:77:LYS:HG3	1:F:148:LEU:O	2.05	0.56
1:C:76:LEU:HD22	1:C:113:LEU:HD11	1.87	0.56
1:C:6:ARG:HH12	1:C:44:ASP:CG	2.09	0.56
1:E:107:GLY:HA2	1:E:110:MET:HE2	1.88	0.56
1:F:70:PHE:CD1	1:F:71:ASN:N	2.69	0.56
1:B:20:GLU:OE2	1:B:59:SER:HB2	2.06	0.55
1:B:3:LYS:O	1:B:5:GLN:N	2.38	0.55
1:B:4:GLY:C	1:B:6:ARG:H	2.10	0.55
1:C:21:THR:OG1	1:C:24:GLU:HG3	2.05	0.55
1:F:91:SER:O	1:F:135:PRO:HA	2.05	0.55
1:D:55:THR:HG22	1:D:56:ASN:N	2.22	0.55
1:A:22:GLN:O	1:A:26:VAL:HG23	2.06	0.55
1:E:90:ALA:O	1:E:91:SER:C	2.45	0.55
1:E:94:ILE:HB	1:E:130:ILE:HB	1.89	0.55
1:E:110:MET:CE	1:E:128:ILE:HG21	2.37	0.55
1:C:53:VAL:HG13	1:C:54:PRO:CD	2.36	0.55
1:A:80:LEU:O	1:A:84:PHE:HB2	2.07	0.55
1:C:122:ILE:HG23	1:C:129:LEU:HB3	1.90	0.54
1:C:68:GLN:HE22	1:D:82:ASP:HB3	1.73	0.54
1:C:125:ASP:O	1:C:126:ASP:HB3	2.08	0.54
1:A:88:ASP:OD1	1:A:89:SER:N	2.40	0.54
1:F:14:ILE:HG12	1:F:62:TYR:O	2.06	0.54
1:B:7:HIS:HB3	1:B:48:LEU:HD21	1.88	0.54
1:C:122:ILE:CG2	1:C:129:LEU:HB3	2.37	0.54
1:C:110:MET:CE	1:C:128:ILE:HG21	2.37	0.54
1:C:51:VAL:CG1	1:C:52:LYS:N	2.70	0.54
1:F:94:ILE:HB	1:F:130:ILE:HB	1.90	0.54
1:C:36:VAL:HG22	1:C:37:THR:N	2.21	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:141:VAL:O	1:F:145:LEU:HG	2.08	0.54
1:F:51:VAL:HG13	1:F:52:LYS:H	1.73	0.53
1:B:143:ASN:O	1:B:147:GLU:HG3	2.08	0.53
1:C:61:LYS:HB2	1:C:61:LYS:NZ	2.23	0.53
1:D:72:PRO:HB2	1:D:113:LEU:HD23	1.90	0.53
1:E:48:LEU:HB2	1:E:50:LEU:HD13	1.88	0.53
1:F:52:LYS:O	1:F:53:VAL:HG23	2.05	0.53
1:C:48:LEU:O	1:C:65:PRO:CG	2.50	0.53
1:E:20:GLU:HG2	1:E:60:TYR:O	2.09	0.53
1:F:36:VAL:CG2	1:F:40:THR:HB	2.39	0.53
1:A:18:GLU:H	1:D:57:ASN:CB	2.19	0.53
1:A:54:PRO:HG2	1:A:69:ARG:HH21	1.74	0.53
1:B:90:ALA:N	1:C:119:MET:HE2	2.24	0.53
1:C:41:VAL:O	1:C:45:ILE:HG13	2.08	0.53
1:C:40:THR:O	1:C:43:ARG:HB3	2.09	0.53
1:D:36:VAL:HG22	1:D:37:THR:N	2.24	0.53
1:B:19:ILE:HG22	1:B:62:TYR:HD2	1.74	0.53
1:A:14:ILE:HG21	1:A:63:SER:HA	1.92	0.53
1:E:6:ARG:O	1:E:10:ILE:HG13	2.08	0.53
1:B:137:ASP:O	1:B:141:VAL:HG23	2.09	0.52
1:A:41:VAL:HG13	1:A:42:SER:N	2.23	0.52
1:C:11:ARG:HG2	1:C:64:LEU:HD21	1.90	0.52
1:E:36:VAL:HG22	1:E:37:THR:N	2.23	0.52
1:A:21:THR:HG23	1:A:24:GLU:H	1.74	0.52
1:C:57:ASN:C	1:C:59:SER:H	2.13	0.52
1:C:84:PHE:CE1	1:C:96:LEU:HD11	2.44	0.52
1:E:49:HIS:O	1:E:65:PRO:HG3	2.09	0.52
1:D:122:ILE:HG22	1:E:122:ILE:HD13	1.92	0.52
1:D:93:MET:CE	1:D:129:LEU:HD21	2.40	0.52
1:A:19:ILE:O	1:A:61:LYS:HA	2.10	0.51
1:B:20:GLU:HG3	1:B:60:TYR:O	2.10	0.51
1:A:45:ILE:HG13	1:A:50:LEU:HD12	1.91	0.51
1:A:113:LEU:HD13	1:A:115:TRP:CZ2	2.46	0.51
1:E:102:ASN:O	1:E:106:ILE:HG12	2.10	0.51
1:E:25:LEU:O	1:E:29:LEU:HB2	2.10	0.51
1:F:76:LEU:O	1:F:76:LEU:HD23	2.11	0.51
1:C:61:LYS:HD3	1:C:61:LYS:O	2.11	0.51
1:D:22:GLN:HG3	1:D:45:ILE:HD12	1.93	0.51
1:D:92:HIS:HB3	1:D:132:CYS:O	2.11	0.50
1:E:91:SER:O	1:E:135:PRO:HA	2.11	0.50
1:A:13:ILE:O	1:A:17:ASN:HB2	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:41:VAL:O	1:A:45:ILE:HD13	2.11	0.50
1:C:65:PRO:C	1:C:67:ASP:N	2.65	0.50
1:B:1:MET:O	1:B:3:LYS:N	2.44	0.50
1:D:72:PRO:HB2	1:D:113:LEU:CD2	2.41	0.50
1:C:26:VAL:HG22	1:C:41:VAL:HG21	1.93	0.50
1:B:68:GLN:O	1:B:72:PRO:HG3	2.12	0.50
1:C:53:VAL:HG12	1:C:54:PRO:HD2	1.91	0.50
1:F:77:LYS:O	1:F:81:MET:HB2	2.11	0.50
1:F:110:MET:HE1	1:F:145:LEU:HD11	1.93	0.50
1:A:43:ARG:O	1:A:47:GLU:HG3	2.11	0.49
1:D:11:ARG:HH12	1:D:64:LEU:HD21	1.75	0.49
1:D:98:THR:HG21	1:D:106:ILE:HD12	1.93	0.49
1:E:86:LYS:HG2	1:E:97:LYS:HB2	1.94	0.49
1:F:87:ILE:HG21	1:F:146:LEU:HD21	1.94	0.49
1:B:19:ILE:N	1:B:19:ILE:HD12	2.27	0.49
1:A:53:VAL:HG13	1:A:69:ARG:HG3	1.96	0.48
1:F:19:ILE:CD1	1:F:25:LEU:HD13	2.43	0.48
1:D:71:ASN:HB3	1:D:74:SER:HB3	1.96	0.48
1:B:91:SER:HB3	1:C:92:HIS:CD2	2.48	0.48
1:C:30:LYS:HA	1:C:34:TYR:O	2.14	0.48
1:C:55:THR:OG1	1:C:61:LYS:HB3	2.14	0.48
1:D:91:SER:O	1:D:135:PRO:HA	2.12	0.48
1:D:52:LYS:HA	1:D:61:LYS:O	2.14	0.47
1:A:14:ILE:CG2	1:A:63:SER:HA	2.43	0.47
1:E:79:ALA:HB1	1:E:106:ILE:CD1	2.43	0.47
1:C:106:ILE:HB	1:C:128:ILE:HD11	1.97	0.47
1:A:55:THR:HG22	1:A:56:ASN:N	2.30	0.47
1:E:22:GLN:O	1:E:26:VAL:HG23	2.15	0.47
1:A:96:LEU:HD23	1:A:128:ILE:HD12	1.95	0.47
1:C:50:LEU:HD23	1:C:65:PRO:HD3	1.97	0.47
1:A:107:GLY:HA3	1:A:123:CYS:SG	2.55	0.47
1:B:92:HIS:HA	1:B:138:THR:OG1	2.15	0.47
1:E:143:ASN:O	1:E:147:GLU:HG3	2.14	0.47
1:E:79:ALA:HB1	1:E:106:ILE:HD11	1.96	0.47
1:F:48:LEU:O	1:F:65:PRO:HG3	2.15	0.47
1:D:12:GLU:O	1:D:15:THR:HG22	2.14	0.47
1:A:66:ALA:HB1	1:F:102:ASN:ND2	2.29	0.47
1:D:71:ASN:N	1:D:72:PRO:HD3	2.31	0.46
1:E:36:VAL:CG2	1:E:41:VAL:HG23	2.45	0.46
1:F:13:ILE:HG22	1:F:19:ILE:HD11	1.98	0.46
1:A:83:ALA:HA	1:A:99:MET:HG3	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:93:MET:HG2	1:B:129:LEU:CD2	2.38	0.46
1:A:24:GLU:O	1:A:28:MET:HG2	2.16	0.46
1:B:90:ALA:HB2	1:C:119:MET:HG2	1.97	0.46
1:D:60:TYR:CD1	1:D:60:TYR:N	2.83	0.46
1:F:149:LEU:C	1:F:149:LEU:HD12	2.36	0.46
1:F:97:LYS:HD3	1:F:127:THR:HB	1.96	0.46
1:B:71:ASN:HD21	1:B:74:SER:CB	2.26	0.46
1:C:117:GLU:O	1:C:133:ARG:HD3	2.15	0.46
1:B:115:TRP:HB3	1:B:117:GLU:OE1	2.16	0.46
1:B:71:ASN:CG	1:B:74:SER:HB3	2.36	0.46
1:B:129:LEU:HD12	1:C:122:ILE:HD12	1.98	0.46
1:C:20:GLU:HA	1:C:62:TYR:CZ	2.51	0.46
1:F:76:LEU:HD12	1:F:113:LEU:HD11	1.97	0.46
1:C:10:ILE:O	1:C:14:ILE:HG12	2.15	0.45
1:D:36:VAL:CG2	1:D:40:THR:HB	2.46	0.45
1:A:52:LYS:HA	1:A:61:LYS:O	2.15	0.45
1:B:106:ILE:O	1:B:110:MET:HG3	2.17	0.45
1:B:132:CYS:SG	1:B:138:THR:HA	2.57	0.45
1:C:36:VAL:CG2	1:C:40:THR:HB	2.47	0.45
1:E:119:MET:HE2	1:E:131:ILE:HG22	1.98	0.45
1:F:21:THR:HG22	1:F:23:ASP:N	2.22	0.45
1:A:133:ARG:HE	1:A:133:ARG:HB2	1.56	0.45
1:D:22:GLN:HG3	1:D:45:ILE:CD1	2.47	0.45
1:F:121:THR:O	1:F:122:ILE:HD13	2.16	0.45
1:A:86:LYS:HG2	1:A:87:ILE:N	2.32	0.45
1:F:117:GLU:H	1:F:117:GLU:CD	2.20	0.45
1:B:3:LYS:HD3	1:B:3:LYS:HA	1.73	0.45
1:B:4:GLY:C	1:B:6:ARG:N	2.70	0.45
1:C:52:LYS:HA	1:C:61:LYS:O	2.16	0.45
1:C:8:ILE:O	1:C:11:ARG:HB2	2.17	0.45
1:F:103:ALA:HB1	1:F:128:ILE:CG1	2.46	0.45
1:B:52:LYS:HA	1:B:61:LYS:O	2.17	0.44
1:C:131:ILE:N	1:C:131:ILE:HD12	2.32	0.44
1:C:139:GLU:O	1:C:142:LYS:HG3	2.16	0.44
1:C:19:ILE:CD1	1:C:19:ILE:H	2.29	0.44
1:C:13:ILE:HG22	1:C:19:ILE:HD11	1.97	0.44
1:F:73:LEU:HD11	1:F:148:LEU:HD21	2.00	0.44
1:F:36:VAL:HG23	1:F:40:THR:HB	1.99	0.44
1:A:98:THR:HG21	1:A:106:ILE:HD12	1.99	0.44
1:B:91:SER:O	1:B:135:PRO:HA	2.17	0.44
1:A:6:ARG:O	1:A:10:ILE:N	2.41	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:71:ASN:N	1:A:72:PRO:HD3	2.33	0.44
1:C:57:ASN:O	1:C:59:SER:N	2.51	0.44
1:C:67:ASP:O	1:C:69:ARG:N	2.38	0.44
1:A:81:MET:HA	1:F:11:ARG:HH22	1.83	0.44
1:E:22:GLN:HG2	1:E:45:ILE:CD1	2.47	0.43
1:F:103:ALA:HB1	1:F:128:ILE:HG13	2.00	0.43
1:B:73:LEU:HD13	1:B:73:LEU:C	2.38	0.43
1:C:79:ALA:HB1	1:C:106:ILE:CD1	2.48	0.43
1:D:19:ILE:HD12	1:D:19:ILE:N	2.33	0.43
1:E:71:ASN:O	1:E:75:LYS:HG3	2.18	0.43
1:A:15:THR:O	1:A:15:THR:HG22	2.18	0.43
1:A:18:GLU:OE1	1:A:61:LYS:HD3	2.18	0.43
1:E:41:VAL:O	1:E:45:ILE:HG13	2.19	0.43
1:F:73:LEU:CD1	1:F:148:LEU:HD21	2.48	0.43
1:D:121:THR:HA	1:D:129:LEU:O	2.18	0.43
1:A:6:ARG:HH11	1:A:6:ARG:HG2	1.84	0.43
1:B:22:GLN:HG3	1:B:45:ILE:CD1	2.49	0.43
1:B:55:THR:CG2	1:B:56:ASN:H	2.27	0.43
1:C:90:ALA:O	1:C:91:SER:C	2.57	0.43
1:E:84:PHE:CE2	1:E:96:LEU:HD11	2.54	0.43
1:F:68:GLN:HG3	1:F:68:GLN:H	1.59	0.43
1:B:92:HIS:HE1	1:B:135:PRO:HD3	1.78	0.43
1:C:57:ASN:C	1:C:59:SER:N	2.71	0.43
1:D:76:LEU:HA	1:D:109:LEU:HD23	2.01	0.43
1:D:77:LYS:O	1:D:81:MET:HG3	2.19	0.43
1:E:137:ASP:O	1:E:141:VAL:HG23	2.19	0.43
1:B:107:GLY:HA3	1:B:123:CYS:SG	2.59	0.42
1:F:48:LEU:CB	1:F:50:LEU:HG	2.48	0.42
1:A:21:THR:HG22	1:A:24:GLU:CG	2.50	0.42
1:A:93:MET:HE1	1:B:131:ILE:HG23	2.00	0.42
1:B:21:THR:OG1	1:B:24:GLU:HG3	2.19	0.42
1:D:104:GLN:NE2	1:D:124:GLY:O	2.52	0.42
1:D:94:ILE:HB	1:D:130:ILE:HB	2.01	0.42
1:A:71:ASN:ND2	1:A:74:SER:HB2	2.34	0.42
1:B:37:THR:HG22	1:B:39:ALA:H	1.84	0.42
1:D:96:LEU:HD23	1:D:128:ILE:HD12	2.00	0.42
1:E:48:LEU:CB	1:E:50:LEU:HD13	2.50	0.42
1:C:48:LEU:O	1:C:49:HIS:HB2	2.19	0.42
1:B:54:PRO:HG3	1:B:60:TYR:CE2	2.55	0.42
1:A:21:THR:HG22	1:A:24:GLU:HG3	2.01	0.42
1:F:14:ILE:CD1	1:F:62:TYR:HB3	2.40	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6:ARG:HH21	1:A:40:THR:HG22	1.84	0.42
1:C:53:VAL:HG12	1:C:54:PRO:CD	2.49	0.42
1:C:36:VAL:CG2	1:C:37:THR:N	2.83	0.42
1:F:126:ASP:OD1	1:F:127:THR:HG22	2.18	0.42
1:F:90:ALA:O	1:F:91:SER:C	2.59	0.41
1:A:10:ILE:O	1:A:14:ILE:HG13	2.20	0.41
1:B:9:LYS:HE3	1:B:9:LYS:HA	2.02	0.41
1:C:99:MET:HG3	1:C:100:PRO:HD2	2.02	0.41
1:D:116:ASP:O	1:D:133:ARG:NH2	2.53	0.41
1:C:103:ALA:CB	1:C:126:ASP:H	2.32	0.41
1:F:113:LEU:O	1:F:114:ASP:OD1	2.39	0.41
1:C:125:ASP:O	1:C:126:ASP:OD1	2.38	0.41
1:C:50:LEU:HD21	1:C:64:LEU:HD12	2.01	0.41
1:B:37:THR:HG22	1:B:38:GLN:N	2.35	0.41
1:A:136:GLU:N	1:A:136:GLU:OE1	2.42	0.41
1:A:14:ILE:HG23	1:A:62:TYR:O	2.21	0.41
1:A:46:LYS:O	1:A:49:HIS:HD2	2.04	0.41
1:A:77:LYS:HE2	1:A:81:MET:HE2	2.03	0.41
1:B:90:ALA:O	1:B:91:SER:C	2.59	0.41
1:E:92:HIS:CE1	1:E:135:PRO:HD3	2.55	0.41
1:C:142:LYS:HD3	1:C:142:LYS:C	2.41	0.41
1:E:113:LEU:HA	1:E:113:LEU:HD12	1.90	0.41
1:F:2:ASN:OD1	1:F:5:GLN:HG3	2.20	0.41
1:F:64:LEU:HA	1:F:65:PRO:HD3	1.86	0.41
1:A:119:MET:CE	1:C:90:ALA:HB2	2.52	0.40
1:A:122:ILE:HD11	1:C:129:LEU:HB2	2.03	0.40
1:E:36:VAL:HG23	1:E:40:THR:OG1	2.21	0.40
1:E:54:PRO:HG3	1:E:60:TYR:CZ	2.56	0.40
1:D:14:ILE:HD11	1:D:64:LEU:CD1	2.51	0.40
1:F:77:LYS:HE3	1:F:81:MET:HE3	2.03	0.40
1:C:6:ARG:HH11	1:C:6:ARG:HG2	1.86	0.40
1:D:35:LYS:CA	1:D:35:LYS:HE3	2.47	0.40
1:D:67:ASP:O	1:D:75:LYS:HE2	2.21	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:144:ARG:NH2	1:A:144:ARG:NH2[3_555]	1.76	0.44



## 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	145/149 (97%)	137 (94%)	7 (5%)	1 (1%)	22	46
1	B	147/149 (99%)	136 (92%)	7 (5%)	4 (3%)	5	12
1	C	146/149 (98%)	125 (86%)	11 (8%)	10 (7%)	1	1
1	D	144/149 (97%)	135 (94%)	8 (6%)	1 (1%)	22	46
1	E	146/149 (98%)	141 (97%)	3 (2%)	2 (1%)	11	28
1	F	147/149 (99%)	132 (90%)	13 (9%)	2 (1%)	11	28
All	All	875/894 (98%)	806 (92%)	49 (6%)	20 (2%)	6	16

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	91	SER
1	B	2	ASN
1	F	72	PRO
1	B	91	SER
1	C	20	GLU
1	C	56	ASN
1	C	66	ALA
1	C	67	ASP
1	C	91	SER
1	C	126	ASP
1	D	91	SER
1	E	70	PHE
1	E	91	SER
1	F	91	SER
1	B	4	GLY
1	C	68	GLN
1	B	5	GLN
1	C	19	ILE
1	C	72	PRO
1	C	58	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	128/133 (96%)	123 (96%)	5 (4%)	32	61
1	B	130/133 (98%)	124 (95%)	6 (5%)	27	54
1	C	125/133 (94%)	115 (92%)	10 (8%)	12	27
1	D	125/133 (94%)	121 (97%)	4 (3%)	39	68
1	E	131/133 (98%)	128 (98%)	3 (2%)	50	78
1	F	128/133 (96%)	120 (94%)	8 (6%)	18	40
All	All	767/798 (96%)	731 (95%)	36 (5%)	26	54

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	GLU
1	A	31	GLN
1	A	76	LEU
1	A	129	LEU
1	A	133	ARG
1	B	9	LYS
1	B	32	ASP
1	B	38	GLN
1	B	129	LEU
1	B	145	LEU
1	B	148	LEU
1	C	35	LYS
1	C	61	LYS
1	C	62	TYR
1	C	68	GLN
1	C	80	LEU
1	C	109	LEU
1	C	119	MET
1	C	126	ASP
1	C	133	ARG
1	C	142	LYS
1	D	35	LYS

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	D	71	ASN
1	D	109	LEU
1	D	129	LEU
1	E	9	LYS
1	E	116	ASP
1	E	129	LEU
1	F	23	ASP
1	F	59	SER
1	F	63	SER
1	F	70	PHE
1	F	104	GLN
1	F	114	ASP
1	F	127	THR
1	F	129	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	A	49	HIS
1	A	102	ASN
1	A	104	GLN
1	A	112	ASN
1	B	38	GLN
1	B	57	ASN
1	B	71	ASN
1	B	92	HIS
1	B	102	ASN
1	B	143	ASN
1	C	68	GLN
1	C	143	ASN
1	D	17	ASN
1	D	22	GLN
1	D	71	ASN
1	D	143	ASN
1	F	38	GLN
1	F	71	ASN
1	F	92	HIS
1	F	102	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 carbohydrates in this entry.

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

There are no ligands in this entry.

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

There are no such residues in this entry.

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

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

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	147/149 (98%)	0.38	12 (8%) <b>11</b> <b>9</b>	35, 67, 133, 146	0
1	B	149/149 (100%)	0.21	4 (2%) 54 55	21, 47, 91, 124	0
1	C	148/149 (99%)	1.06	30 (20%) <b>1</b> <b>0</b>	33, 66, 150, 164	0
1	D	146/149 (97%)	0.12	2 (1%) 75 77	31, 59, 109, 127	0
1	E	148/149 (99%)	0.08	2 (1%) 75 77	24, 43, 74, 83	0
1	F	149/149 (100%)	0.88	24 (16%) <b>1</b> <b>1</b>	39, 76, 151, 165	0
All	All	887/894 (99%)	0.46	74 (8%) <b>11</b> <b>9</b>	21, 59, 130, 165	0

All (74) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	59	SER	7.8
1	A	39	ALA	7.7
1	C	53	VAL	7.5
1	C	54	PRO	7.4
1	C	70	PHE	7.2
1	C	55	THR	7.1
1	C	67	ASP	6.8
1	C	57	ASN	6.2
1	C	60	TYR	6.1
1	F	56	ASN	5.7
1	C	65	PRO	5.7
1	F	60	TYR	5.5
1	F	53	VAL	5.1
1	F	54	PRO	4.9
1	C	69	ARG	4.9
1	F	57	ASN	4.9
1	F	52	LYS	4.6
1	C	66	ALA	4.4
1	C	71	ASN	4.2

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	C	58	GLY	4.0
1	F	16	SER	3.9
1	C	56	ASN	3.9
1	A	34	TYR	3.8
1	C	19	ILE	3.7
1	C	59	SER	3.7
1	F	66	ALA	3.5
1	A	41	VAL	3.4
1	F	55	THR	3.2
1	F	68	GLN	3.1
1	A	27	ASP	3.1
1	F	67	ASP	3.1
1	C	61	LYS	3.0
1	F	63	SER	3.0
1	C	72	PRO	3.0
1	F	71	ASN	3.0
1	F	62	TYR	3.0
1	F	58	GLY	3.0
1	B	2	ASN	2.9
1	F	15	THR	2.8
1	C	51	VAL	2.8
1	C	63	SER	2.7
1	F	51	VAL	2.7
1	F	64	LEU	2.7
1	C	90	ALA	2.7
1	C	52	LYS	2.6
1	B	57	ASN	2.5
1	A	3	LYS	2.5
1	B	58	GLY	2.4
1	A	35	LYS	2.4
1	C	49	HIS	2.4
1	C	45	ILE	2.4
1	D	57	ASN	2.3
1	F	19	ILE	2.3
1	E	33	GLY	2.3
1	C	21	THR	2.3
1	C	31	GLN	2.3
1	C	33	GLY	2.3
1	A	37	THR	2.3
1	A	26	VAL	2.3
1	A	30	LYS	2.2
1	A	43	ARG	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	39	ALA	2.2
1	D	23	ASP	2.2
1	F	2	ASN	2.2
1	F	11	ARG	2.2
1	A	36	VAL	2.2
1	C	62	TYR	2.1
1	B	1	MET	2.1
1	F	1	MET	2.1
1	C	64	LEU	2.1
1	F	14	ILE	2.1
1	C	34	TYR	2.1
1	E	90	ALA	2.0
1	A	6	ARG	2.0

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

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

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.