



# Full wwPDB X-ray Structure Validation Report ⓘ

May 21, 2020 – 11:40 pm BST

PDB ID : 1LQM  
Title : ESCHERICHIA COLI URACIL-DNA GLYCOSYLASE COMPLEX WITH  
URACIL-DNA GLYCOSYLASE INHIBITOR PROTEIN  
Authors : Saikrishnan, K.; Sagar, M.B.; Ravishankar, R.; Roy, S.; Purnapatre, K.; Varsh-  
ney, U.; Vijayan, M.  
Deposited on : 2002-05-10  
Resolution : 3.20 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	:	2.11
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

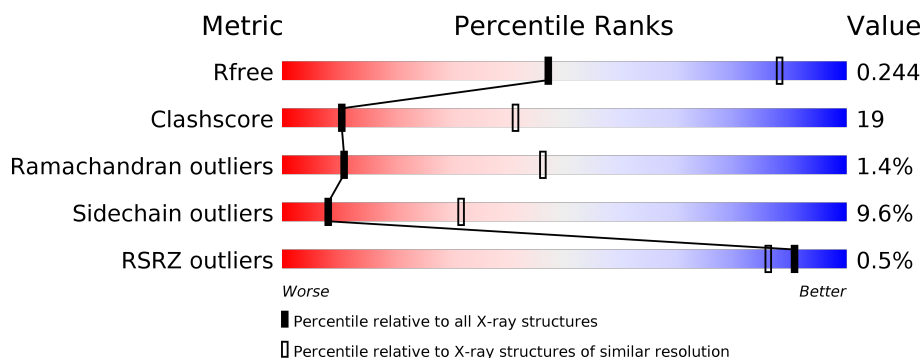
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.20 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	1133 (3.20-3.20)
Clashscore	141614	1253 (3.20-3.20)
Ramachandran outliers	138981	1234 (3.20-3.20)
Sidechain outliers	138945	1233 (3.20-3.20)
RSRZ outliers	127900	1095 (3.20-3.20)

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	229	<div> <div>58%</div> <div>33%</div> <div>5%</div> <div>• •</div> </div>
1	C	229	<div> <div>59%</div> <div>36%</div> <div>• •</div> </div>
1	E	229	<div> <div>58%</div> <div>35%</div> <div>5%</div> <div>•</div> </div>
1	G	229	<div> <div>57%</div> <div>37%</div> <div>• • •</div> </div>
2	B	84	<div> <div>61%</div> <div>32%</div> <div>• • •</div> </div>
2	D	84	<div> <div>61%</div> <div>33%</div> <div>• •</div> </div>

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Mol	Chain	Length	Quality of chain
2	F	84	<div><div></div><div>64%</div><div>31%</div><div></div><div>• •</div></div>
2	H	84	<div><div>2%</div><div></div><div>52%</div><div>39%</div><div>6%</div><div></div><div>•</div></div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9744 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 URACIL-DNA GLYCOSYLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	223	Total	C	N	O	S	0	0	0
			1764	1138	316	307	3			
1	C	224	Total	C	N	O	S	0	0	0
			1773	1143	317	310	3			
1	E	224	Total	C	N	O	S	0	0	0
			1773	1141	318	311	3			
1	G	225	Total	C	N	O	S	0	0	0
			1777	1145	319	310	3			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	CLONING ARTIFACT	UNP P12295
C	1	MET	-	CLONING ARTIFACT	UNP P12295
E	1	MET	-	CLONING ARTIFACT	UNP P12295
G	1	MET	-	CLONING ARTIFACT	UNP P12295

- Molecule 2 is a protein called URACIL-DNA GLYCOSYLASE INHIBITOR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	82	Total	C	N	O	S	0	0	0
			643	403	98	139	3			
2	D	83	Total	C	N	O	S	0	0	0
			652	409	100	140	3			
2	F	82	Total	C	N	O	S	0	0	0
			647	406	99	139	3			
2	H	84	Total	C	N	O	S	0	0	0
			657	412	101	141	3			

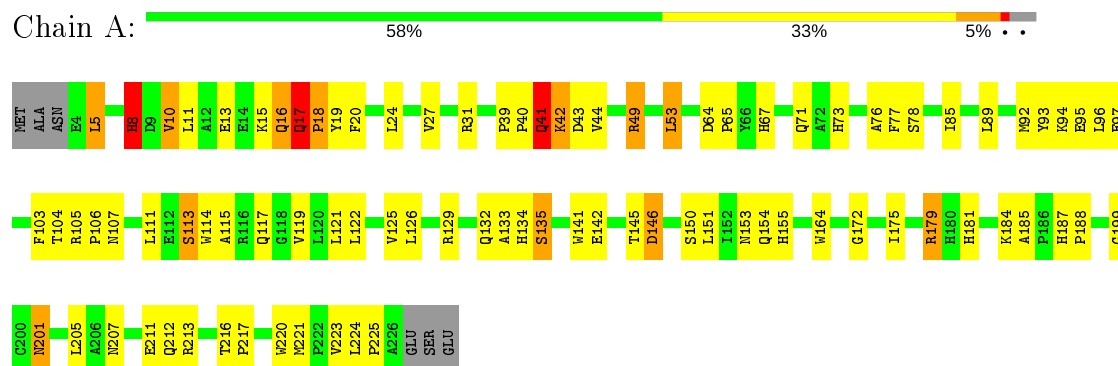
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	19	Total 19	O 19	0	0
3	B	5	Total 5	O 5	0	0
3	C	9	Total 9	O 9	0	0
3	D	3	Total 3	O 3	0	0
3	E	16	Total 16	O 16	0	0
3	F	1	Total 1	O 1	0	0
3	G	4	Total 4	O 4	0	0
3	H	1	Total 1	O 1	0	0

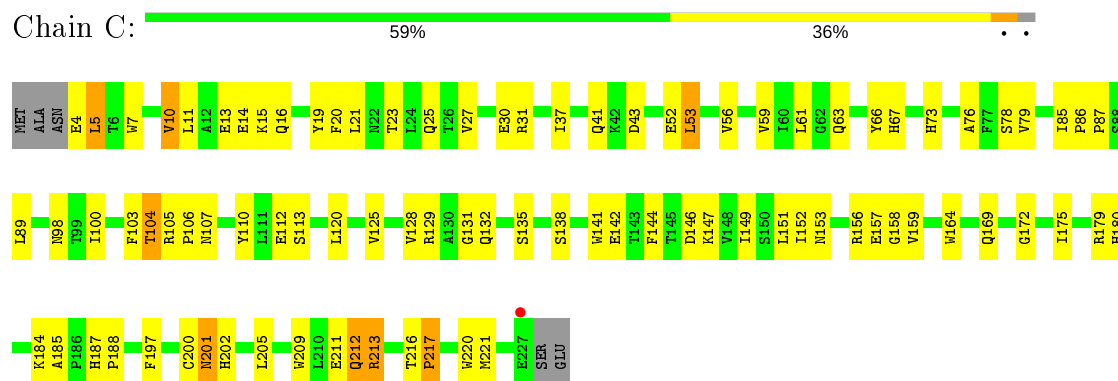
### 3 Residue-property plots [i](#)

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.

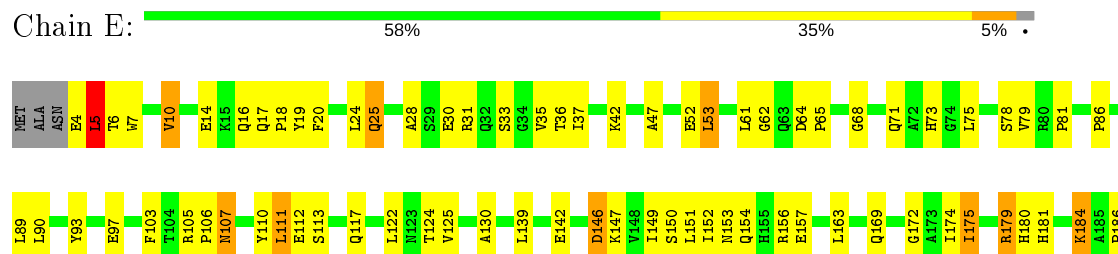
#### • Molecule 1: URACIL-DNA GLYCOSYLASE



#### • Molecule 1: URACIL-DNA GLYCOSYLASE

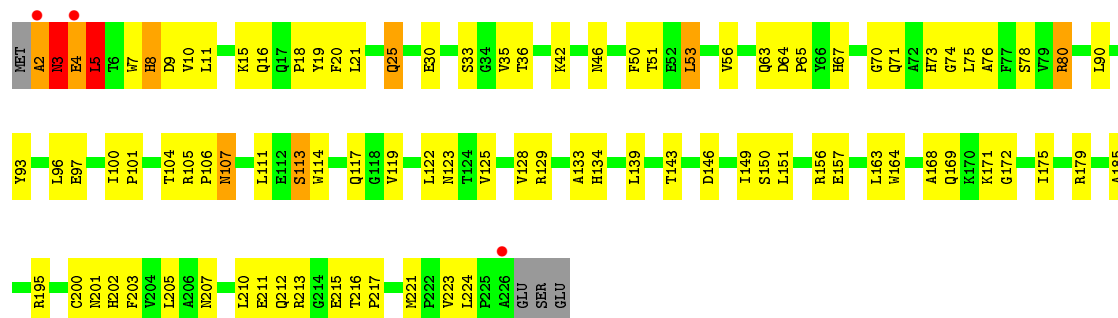


#### • Molecule 1: URACIL-DNA GLYCOSYLASE





• Molecule 1: URACIL-DNA GLYCOSYLASE



• Molecule 2: URACIL-DNA GLYCOSYLASE INHIBITOR



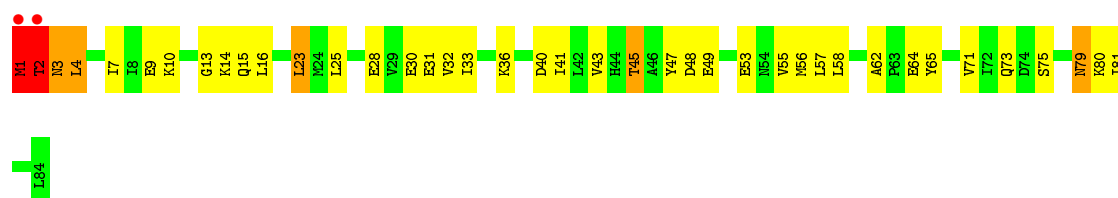
• Molecule 2: URACIL-DNA GLYCOSYLASE INHIBITOR



• Molecule 2: URACIL-DNA GLYCOSYLASE INHIBITOR



• Molecule 2: URACIL-DNA GLYCOSYLASE INHIBITOR



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	98.76 Å   158.88 Å   91.22 Å 90.00°   90.00°   90.00°	Depositor
Resolution (Å)	15.00 – 3.20 15.00 – 3.20	Depositor EDS
% Data completeness (in resolution range)	78.3 (15.00-3.20) 78.3 (15.00-3.20)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.17	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.57 (at 3.19 Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, $R_{free}$	0.188   ,   0.260 0.182   ,   0.244	Depositor DCC
$R_{free}$ test set	916 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	33.9	Xtriage
Anisotropy	0.377	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 63.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	9744	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.25% 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

### 5.1 Standard geometry

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	0.64	7/1820 (0.4%)	0.74	5/2483 (0.2%)
1	C	0.39	0/1829	0.64	0/2495
1	E	0.47	1/1829 (0.1%)	0.66	1/2495 (0.0%)
1	G	0.47	0/1833	0.84	8/2501 (0.3%)
2	B	0.67	3/651 (0.5%)	0.87	3/883 (0.3%)
2	D	0.63	2/660 (0.3%)	0.75	3/894 (0.3%)
2	F	0.37	0/655	0.58	0/887
2	H	0.64	2/665 (0.3%)	1.06	3/901 (0.3%)
All	All	0.53	15/9942 (0.2%)	0.75	23/13539 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	E	0	1
1	G	0	4
2	H	0	2
All	All	0	8

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	5	LEU	CA-CB	9.24	1.75	1.53
2	B	3	ASN	C-O	8.30	1.39	1.23
1	A	17	GLN	CB-CG	8.01	1.74	1.52
2	D	3	ASN	CB-CG	7.56	1.68	1.51
1	A	49	ARG	CB-CG	7.37	1.72	1.52
1	A	17	GLN	CA-CB	7.34	1.70	1.53
2	D	2	THR	C-O	6.89	1.36	1.23
2	H	1	MET	N-CA	6.16	1.58	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	3	ASN	CB-CG	-6.13	1.36	1.51
1	A	8	HIS	CA-CB	5.96	1.67	1.53
1	A	49	ARG	CZ-NH1	-5.82	1.25	1.33
2	B	4	LEU	C-O	5.68	1.34	1.23
1	A	17	GLN	CG-CD	5.45	1.63	1.51
1	A	49	ARG	CG-CD	5.22	1.65	1.51
2	H	2	THR	CA-C	5.12	1.66	1.52

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	1	MET	CA-C-N	-14.43	85.44	117.20
1	G	4	GLU	CB-CA-C	14.26	138.93	110.40
2	H	2	THR	N-CA-CB	-12.65	86.26	110.30
2	H	1	MET	O-C-N	12.26	142.31	122.70
1	G	3	ASN	N-CA-C	11.77	142.77	111.00
1	G	3	ASN	N-CA-CB	-9.98	92.64	110.60
1	E	4	GLU	C-N-CA	-9.78	97.26	121.70
2	B	4	LEU	CA-CB-CG	9.21	136.48	115.30
2	B	3	ASN	N-CA-C	8.16	133.02	111.00
2	B	3	ASN	C-N-CA	-7.80	102.20	121.70
1	G	2	ALA	CA-C-N	-7.58	100.52	117.20
1	A	17	GLN	CB-CA-C	7.32	125.03	110.40
1	A	49	ARG	CD-NE-CZ	6.94	133.32	123.60
1	G	4	GLU	CA-C-N	-6.83	102.19	117.20
1	A	41	GLN	CA-CB-CG	6.43	127.56	113.40
2	D	2	THR	CA-C-N	6.30	131.06	117.20
2	D	2	THR	C-N-CA	-5.98	106.74	121.70
1	G	2	ALA	CB-CA-C	5.88	118.93	110.10
1	G	4	GLU	CA-C-O	5.79	132.26	120.10
1	A	41	GLN	CB-CG-CD	5.75	126.54	111.60
2	D	3	ASN	N-CA-CB	5.71	120.88	110.60
1	A	8	HIS	CA-CB-CG	5.52	122.98	113.60
1	G	5	LEU	N-CA-C	5.34	125.42	111.00

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	8	HIS	Sidechain
1	E	5	LEU	Mainchain
1	G	2	ALA	Mainchain,Peptide

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Mol	Chain	Res	Type	Group
1	G	3	ASN	Mainchain
1	G	8	HIS	Sidechain
2	H	1	MET	Mainchain,Peptide

## 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	1764	0	1730	71	0
1	C	1773	0	1736	62	0
1	E	1773	0	1731	63	0
1	G	1777	0	1741	63	0
2	B	643	0	629	24	0
2	D	652	0	642	21	0
2	F	647	0	640	23	0
2	H	657	0	647	33	0
3	A	19	0	0	4	0
3	B	5	0	0	0	0
3	C	9	0	0	1	0
3	D	3	0	0	0	0
3	E	16	0	0	0	0
3	F	1	0	0	0	0
3	G	4	0	0	0	0
3	H	1	0	0	0	0
All	All	9744	0	9496	355	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 19.

All (355) 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:E:5:LEU:CB	1:E:5:LEU:CA	1.75	1.54
1:A:41:GLN:NE2	1:A:44:VAL:HG21	1.33	1.37
1:A:41:GLN:HE21	1:A:44:VAL:CG2	1.56	1.18
2:D:45:THR:HG23	2:D:56:MET:HG2	1.20	1.13

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:213:ARG:HB3	1:C:213:ARG:HH11	1.27	0.99
1:G:4:GLU:C	1:G:5:LEU:HD23	1.84	0.95
1:C:10:VAL:HB	1:C:151:LEU:HD13	1.49	0.94
2:F:30:GLU:HG3	2:F:36:LYS:HB2	1.50	0.93
2:H:4:LEU:HA	2:H:7:ILE:HD12	1.53	0.91
1:A:41:GLN:HA	1:A:44:VAL:HG23	1.54	0.87
2:D:30:GLU:HG3	2:D:36:LYS:HB2	1.57	0.86
2:D:45:THR:HG23	2:D:56:MET:CG	2.05	0.84
1:E:172:GLY:HA2	1:E:175:ILE:HD12	1.59	0.84
2:D:14:LYS:HD3	2:D:48:ASP:OD2	1.79	0.81
1:C:67:HIS:HB2	1:C:131:GLY:H	1.44	0.81
2:B:3:ASN:ND2	2:B:3:ASN:O	2.14	0.80
1:E:78:SER:HB2	1:E:111:LEU:HG	1.63	0.78
2:H:30:GLU:HG3	2:H:36:LYS:HB2	1.64	0.78
1:E:172:GLY:HA2	1:E:175:ILE:CD1	2.15	0.76
2:H:1:MET:N	2:H:2:THR:CB	2.49	0.76
2:F:45:THR:HG23	2:F:56:MET:HG2	1.69	0.75
1:G:172:GLY:HA2	1:G:175:ILE:HD12	1.68	0.75
2:H:45:THR:HG23	2:H:56:MET:HG2	1.70	0.74
1:A:73:HIS:HE2	1:A:78:SER:HG	1.34	0.73
1:C:213:ARG:HB3	1:C:213:ARG:NH1	2.02	0.73
1:A:67:HIS:CD2	1:A:133:ALA:HB2	2.24	0.72
1:G:63:GLN:O	1:G:123:ASN:HB3	1.90	0.72
1:G:4:GLU:O	1:G:5:LEU:HD23	1.88	0.71
2:F:41:ILE:HD13	2:F:58:LEU:HD13	1.73	0.71
1:A:10:VAL:HB	1:A:151:LEU:HD13	1.71	0.71
1:E:93:TYR:CZ	1:E:106:PRO:HG2	2.26	0.70
2:B:41:ILE:HD13	2:B:58:LEU:HD13	1.73	0.70
1:C:201:ASN:HB3	1:C:205:LEU:HD13	1.72	0.70
1:A:27:VAL:O	1:A:31:ARG:HG3	1.93	0.69
2:H:43:VAL:HG13	2:H:58:LEU:HD21	1.73	0.69
1:G:73:HIS:HE1	1:G:75:LEU:HD12	1.58	0.68
1:G:201:ASN:HB3	1:G:205:LEU:HD13	1.73	0.68
2:H:43:VAL:HG13	2:H:58:LEU:CD2	2.23	0.68
2:F:29:VAL:HG21	2:F:41:ILE:HG12	1.76	0.68
1:G:73:HIS:CE1	1:G:75:LEU:HD12	2.28	0.68
1:A:172:GLY:HA2	1:A:175:ILE:HD12	1.74	0.67
1:C:23:THR:O	1:C:27:VAL:HG23	1.95	0.67
2:H:1:MET:H3	2:H:2:THR:CB	2.09	0.67
1:C:59:VAL:HG22	1:C:120:LEU:HB3	1.76	0.66
2:H:23:LEU:HD11	2:H:40:ASP:HB3	1.77	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:201:ASN:O	1:C:205:LEU:HD13	1.96	0.66
1:A:41:GLN:NE2	1:A:44:VAL:CG2	2.28	0.65
2:B:3:ASN:CG	2:B:3:ASN:O	2.34	0.65
1:A:113:SER:O	1:A:117:GLN:HG3	1.96	0.64
1:C:63:GLN:OE1	2:D:23:LEU:HB3	1.97	0.64
1:C:43:ASP:HB3	1:C:73:HIS:O	1.97	0.64
1:G:10:VAL:HB	1:G:151:LEU:HD13	1.79	0.64
1:A:122:LEU:HD11	1:A:145:THR:HG22	1.80	0.64
2:D:47:TYR:CE2	2:D:49:GLU:HA	2.32	0.64
1:G:213:ARG:HB3	1:G:213:ARG:NH1	2.13	0.64
2:H:1:MET:C	2:H:3:ASN:H	2.01	0.64
2:F:71:VAL:CG1	2:F:79:ASN:HB2	2.29	0.63
1:A:211:GLU:HG2	1:A:217:PRO:HG3	1.81	0.63
1:E:90:LEU:HD12	1:E:90:LEU:O	1.98	0.63
1:A:129:ARG:HG3	1:A:135:SER:HB2	1.80	0.62
1:A:164:TRP:HA	1:A:185:ALA:O	2.00	0.62
1:E:163:LEU:HD13	1:E:169:GLN:HA	1.81	0.62
1:A:201:ASN:HB3	1:A:205:LEU:HD13	1.82	0.62
1:E:172:GLY:CA	1:E:175:ILE:HD12	2.30	0.62
1:E:201:ASN:HB3	1:E:205:LEU:HD13	1.81	0.61
1:G:53:LEU:O	1:G:156:ARG:NH1	2.33	0.61
1:G:70:GLY:O	1:G:80:ARG:HG3	2.00	0.61
1:A:201:ASN:O	1:A:205:LEU:HD13	2.00	0.61
1:C:4:GLU:N	3:C:234:HOH:O	2.34	0.61
1:E:78:SER:HB2	1:E:111:LEU:CG	2.30	0.61
1:C:211:GLU:C	1:C:213:ARG:H	2.04	0.61
2:B:41:ILE:HD13	2:B:58:LEU:HB3	1.83	0.60
1:G:113:SER:O	1:G:117:GLN:HG3	2.01	0.60
2:B:41:ILE:CD1	2:B:58:LEU:HD13	2.31	0.60
1:C:128:VAL:HG22	1:C:129:ARG:N	2.16	0.60
1:E:10:VAL:HB	1:E:151:LEU:HD13	1.82	0.59
1:E:61:LEU:HD21	1:E:149:ILE:HD11	1.83	0.59
2:F:30:GLU:HG3	2:F:36:LYS:CB	2.30	0.59
1:E:86:PRO:HD2	1:E:89:LEU:HD23	1.85	0.59
1:C:103:PHE:HE2	1:C:221:MET:HA	1.67	0.59
2:D:41:ILE:HD13	2:D:58:LEU:HD13	1.84	0.59
1:A:41:GLN:HA	1:A:44:VAL:CG2	2.30	0.59
2:H:73:GLN:HB2	2:H:79:ASN:HB3	1.84	0.58
2:F:41:ILE:CD1	2:F:58:LEU:HD13	2.33	0.58
2:F:58:LEU:N	2:F:58:LEU:HD23	2.18	0.58
1:E:5:LEU:CB	1:E:5:LEU:N	2.60	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:41:GLN:O	1:A:43:ASP:N	2.37	0.57
2:H:4:LEU:HB3	2:H:57:LEU:HD23	1.87	0.57
1:A:39:PRO:HB2	1:A:40:PRO:CD	2.35	0.57
2:B:73:GLN:HB2	2:B:79:ASN:HB3	1.86	0.57
2:D:71:VAL:CG1	2:D:79:ASN:HB2	2.34	0.57
1:G:128:VAL:HG22	1:G:129:ARG:N	2.19	0.57
1:E:110:TYR:CE2	1:E:112:GLU:HB2	2.40	0.57
2:F:30:GLU:OE1	2:F:36:LYS:HD2	2.04	0.57
1:G:90:LEU:O	1:G:93:TYR:HB2	2.05	0.57
2:H:23:LEU:CD1	2:H:40:ASP:HB3	2.35	0.57
1:G:213:ARG:HH11	1:G:213:ARG:HB3	1.70	0.56
1:A:41:GLN:HE21	1:A:44:VAL:HG21	0.74	0.56
2:H:1:MET:H2	2:H:2:THR:CB	2.18	0.56
1:C:212:GLN:HG2	1:C:212:GLN:O	2.05	0.56
1:G:172:GLY:HA2	1:G:175:ILE:CD1	2.36	0.56
1:A:142:GLU:O	1:A:146:ASP:HB2	2.05	0.56
2:H:43:VAL:HG22	2:H:58:LEU:HD22	1.88	0.56
2:D:29:VAL:HG21	2:D:41:ILE:HG12	1.88	0.56
1:C:66:TYR:OH	1:C:86:PRO:HG2	2.06	0.56
1:A:187:HIS:CG	1:A:188:PRO:HD2	2.40	0.56
1:C:201:ASN:HB3	1:C:205:LEU:CD1	2.36	0.56
2:H:1:MET:C	2:H:3:ASN:N	2.58	0.56
1:A:67:HIS:HD2	1:A:133:ALA:HB2	1.69	0.55
1:E:68:GLY:HA3	1:E:71:GLN:OE1	2.06	0.55
1:C:172:GLY:HA2	1:C:175:ILE:HD12	1.87	0.55
1:E:153:ASN:O	1:E:179:ARG:NH1	2.40	0.55
1:C:132:GLN:HB3	1:C:135:SER:HB3	1.89	0.55
1:G:106:PRO:O	1:G:107:ASN:ND2	2.40	0.55
1:G:212:GLN:HG2	1:G:212:GLN:O	2.06	0.55
2:H:33:ILE:HD11	2:H:71:VAL:HG21	1.88	0.55
2:H:45:THR:HG23	2:H:56:MET:CG	2.37	0.55
2:H:41:ILE:HD13	2:H:58:LEU:HD13	1.87	0.55
1:G:4:GLU:O	1:G:9:ASP:OD2	2.25	0.55
1:C:211:GLU:CG	1:C:217:PRO:HG3	2.38	0.54
2:F:30:GLU:CG	2:F:36:LYS:HB2	2.32	0.54
1:A:104:THR:O	1:A:221:MET:CE	2.55	0.54
1:G:172:GLY:CA	1:G:175:ILE:HD12	2.36	0.54
1:E:213:ARG:NH1	1:E:213:ARG:HB3	2.22	0.54
1:E:147:LYS:O	1:E:151:LEU:HG	2.08	0.54
1:A:213:ARG:NH1	1:A:213:ARG:HB3	2.23	0.53
1:C:125:VAL:HA	1:C:141:TRP:HB3	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:189:SER:O	1:E:193:ALA:HB2	2.08	0.53
2:F:33:ILE:HD11	2:F:71:VAL:HG21	1.90	0.53
1:C:104:THR:O	1:C:221:MET:HE1	2.08	0.53
1:E:113:SER:O	1:E:117:GLN:HG3	2.08	0.53
1:C:52:GLU:O	1:C:56:VAL:HG23	2.08	0.53
2:D:6:ASP:O	2:D:10:LYS:HB3	2.08	0.53
1:G:4:GLU:C	1:G:5:LEU:CD2	2.70	0.53
1:A:201:ASN:HB3	1:A:205:LEU:CD1	2.38	0.52
2:B:14:LYS:HD3	2:B:48:ASP:OD2	2.09	0.52
1:G:50:PHE:HD2	1:G:75:LEU:HD21	1.73	0.52
1:C:187:HIS:CG	1:C:188:PRO:HD2	2.44	0.52
1:E:64:ASP:HB2	1:E:65:PRO:CD	2.40	0.52
1:E:6:THR:HA	1:E:52:GLU:OE2	2.10	0.52
2:F:9:GLU:O	2:F:13:GLY:N	2.39	0.52
2:H:71:VAL:CG1	2:H:79:ASN:HB2	2.40	0.51
1:E:154:GLN:HA	1:E:179:ARG:HH12	1.74	0.51
1:E:195:ARG:HG3	2:F:32:VAL:HG22	1.92	0.51
1:A:95:GLU:CD	1:A:199:GLY:H	2.14	0.51
1:A:78:SER:HB2	1:A:111:LEU:HB2	1.92	0.51
1:A:85:ILE:HG23	1:A:89:LEU:HD23	1.91	0.51
1:C:164:TRP:HA	1:C:185:ALA:O	2.11	0.51
1:C:67:HIS:HB2	1:C:131:GLY:N	2.20	0.51
1:G:223:VAL:HG12	1:G:224:LEU:N	2.26	0.51
1:G:5:LEU:HD23	1:G:5:LEU:N	2.21	0.51
1:G:146:ASP:O	1:G:149:ILE:HB	2.12	0.50
1:C:106:PRO:O	1:C:107:ASN:CB	2.58	0.50
1:E:36:THR:HG22	1:E:130:ALA:HB2	1.92	0.50
1:A:115:ALA:HA	1:A:119:VAL:O	2.11	0.50
1:A:96:LEU:HD22	1:A:220:TRP:HB3	1.94	0.50
1:C:53:LEU:O	1:C:53:LEU:HD22	2.11	0.50
1:E:211:GLU:HG2	1:E:217:PRO:HG3	1.93	0.50
2:B:57:LEU:C	2:B:58:LEU:HD23	2.32	0.50
1:E:30:GLU:O	1:E:35:VAL:HB	2.12	0.50
2:D:58:LEU:N	2:D:58:LEU:HD23	2.26	0.50
1:A:93:TYR:CZ	1:A:106:PRO:HG2	2.46	0.50
1:A:114:TRP:O	1:A:119:VAL:HB	2.11	0.50
2:B:58:LEU:O	2:B:68:TRP:HB3	2.11	0.50
1:G:104:THR:O	1:G:221:MET:HE1	2.12	0.50
1:G:4:GLU:O	1:G:5:LEU:CD2	2.57	0.50
1:A:76:ALA:HB3	1:A:121:LEU:HB3	1.93	0.49
1:A:106:PRO:O	1:A:107:ASN:CB	2.60	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:30:GLU:HG3	2:B:36:LYS:HB2	1.94	0.49
1:G:210:LEU:O	1:G:215:GLU:HB2	2.12	0.49
1:C:153:ASN:O	1:C:179:ARG:NH1	2.46	0.49
2:H:14:LYS:HD3	2:H:48:ASP:OD2	2.13	0.48
1:C:7:TRP:CE3	1:C:7:TRP:HA	2.48	0.48
2:F:49:GLU:HG3	2:F:50:SER:N	2.29	0.48
1:E:14:GLU:HA	1:E:17:GLN:HG2	1.95	0.48
1:A:172:GLY:HA2	1:A:175:ILE:CD1	2.42	0.48
1:A:132:GLN:O	1:A:135:SER:OG	2.24	0.48
1:C:128:VAL:CG2	1:C:129:ARG:N	2.76	0.48
1:G:19:TYR:CG	1:G:20:PHE:N	2.82	0.48
1:G:50:PHE:CD2	1:G:75:LEU:HD21	2.49	0.48
1:C:110:TYR:CE2	1:C:112:GLU:HB2	2.49	0.48
1:G:63:GLN:OE1	2:H:23:LEU:HB3	2.13	0.48
2:B:73:GLN:HA	2:B:79:ASN:HA	1.95	0.47
1:E:163:LEU:O	1:E:184:LYS:HA	2.13	0.47
1:E:211:GLU:C	1:E:213:ARG:H	2.18	0.47
1:G:211:GLU:CG	1:G:217:PRO:HG3	2.44	0.47
1:A:126:LEU:HD12	1:A:141:TRP:CZ3	2.49	0.47
2:H:9:GLU:O	2:H:13:GLY:N	2.46	0.47
2:B:45:THR:HG23	2:B:56:MET:HG2	1.95	0.47
1:E:64:ASP:HB3	1:E:125:VAL:HB	1.95	0.47
2:B:23:LEU:HD11	2:B:40:ASP:HB3	1.97	0.47
1:G:96:LEU:HD21	1:G:203:PHE:CD1	2.50	0.47
1:G:93:TYR:CZ	1:G:106:PRO:HG2	2.50	0.47
1:G:46:ASN:ND2	1:G:73:HIS:O	2.47	0.47
1:E:105:ARG:HA	1:E:106:PRO:HD3	1.69	0.47
1:E:142:GLU:O	1:E:146:ASP:HB2	2.14	0.47
2:F:43:VAL:HG13	2:F:58:LEU:CD2	2.44	0.47
1:G:30:GLU:O	1:G:35:VAL:HB	2.15	0.47
1:G:46:ASN:HB2	1:G:74:GLY:HA3	1.97	0.47
2:H:3:ASN:O	2:H:4:LEU:C	2.53	0.47
1:E:211:GLU:CG	1:E:217:PRO:HG3	2.45	0.47
1:C:37:ILE:HG23	1:C:128:VAL:O	2.15	0.46
1:A:184:LYS:O	1:A:185:ALA:HB2	2.15	0.46
1:A:94:LYS:O	1:A:97:GLU:HB2	2.15	0.46
2:F:12:THR:OG1	2:F:14:LYS:HG3	2.15	0.46
2:F:26:PRO:HA	2:F:37:PRO:HG2	1.97	0.46
1:A:65:PRO:HD3	1:A:125:VAL:O	2.15	0.46
1:A:213:ARG:HH11	1:A:213:ARG:HB3	1.80	0.46
1:A:41:GLN:O	1:A:42:LYS:C	2.52	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:105:ARG:HA	1:C:106:PRO:HD3	1.70	0.46
1:C:103:PHE:CE2	1:C:221:MET:HA	2.49	0.46
1:A:153:ASN:O	1:A:179:ARG:NH1	2.48	0.46
2:B:25:LEU:HD23	2:B:25:LEU:N	2.30	0.46
2:B:60:SER:OG	2:B:64:GLU:HB3	2.16	0.46
1:C:14:GLU:OE2	1:C:147:LYS:HG2	2.15	0.46
2:D:23:LEU:HD11	2:D:40:ASP:HB3	1.98	0.46
2:F:25:LEU:HD23	2:F:25:LEU:N	2.31	0.46
1:E:73:HIS:HE2	1:E:78:SER:HG	1.62	0.46
2:B:59:THR:HA	2:B:66:LYS:O	2.16	0.46
2:D:56:MET:O	2:D:70:LEU:HD12	2.16	0.46
1:C:159:VAL:O	1:C:180:HIS:HB3	2.16	0.46
1:C:30:GLU:HB2	1:C:37:ILE:HD11	1.98	0.46
1:E:10:VAL:HG21	1:E:53:LEU:HD11	1.98	0.46
1:G:106:PRO:O	1:G:107:ASN:CB	2.64	0.46
1:A:154:GLN:OE1	1:A:155:HIS:CE1	2.69	0.46
1:E:207:ASN:CG	1:E:217:PRO:HB3	2.37	0.46
1:A:41:GLN:C	1:A:43:ASP:N	2.69	0.45
2:D:25:LEU:O	2:D:29:VAL:HG23	2.16	0.45
1:G:19:TYR:CZ	1:G:143:THR:HG22	2.51	0.45
2:D:43:VAL:HG13	2:D:58:LEU:CD2	2.46	0.45
2:D:43:VAL:HG13	2:D:58:LEU:HD22	1.98	0.45
2:D:24:MET:HA	2:D:28:GLU:OE1	2.17	0.45
2:B:29:VAL:HG21	2:B:41:ILE:HG12	1.99	0.45
2:D:3:ASN:O	2:D:7:ILE:HG13	2.16	0.45
2:H:62:ALA:HB2	2:H:65:TYR:CE1	2.52	0.45
1:C:172:GLY:HA2	1:C:175:ILE:CD1	2.47	0.45
1:A:41:GLN:O	1:A:44:VAL:N	2.48	0.45
1:C:211:GLU:C	1:C:213:ARG:N	2.68	0.45
2:F:71:VAL:HG12	2:F:79:ASN:HB2	1.98	0.45
1:A:17:GLN:C	3:A:230:HOH:O	2.55	0.44
1:E:25:GLN:O	1:E:28:ALA:HB3	2.17	0.44
1:A:122:LEU:HD11	1:A:145:THR:CG2	2.45	0.44
1:G:211:GLU:C	1:G:213:ARG:H	2.20	0.44
1:G:64:ASP:HB2	1:G:65:PRO:CD	2.47	0.44
2:B:58:LEU:N	2:B:58:LEU:HD23	2.31	0.44
1:A:105:ARG:HA	1:A:106:PRO:HD3	1.70	0.44
1:A:154:GLN:HA	1:A:179:ARG:HH12	1.82	0.44
1:G:53:LEU:O	1:G:56:VAL:HG23	2.17	0.44
1:G:64:ASP:HB2	1:G:65:PRO:HD2	2.00	0.44
1:A:221:MET:HE2	3:A:237:HOH:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:100:ILE:HD11	1:C:220:TRP:CD1	2.53	0.44
1:E:210:LEU:O	1:E:215:GLU:HB2	2.17	0.44
1:A:19:TYR:CG	1:A:20:PHE:N	2.86	0.43
2:B:56:MET:HE3	2:B:71:VAL:HB	1.99	0.43
2:D:9:GLU:O	2:D:13:GLY:N	2.42	0.43
1:E:62:GLY:O	1:E:124:THR:HG23	2.18	0.43
1:G:128:VAL:HG22	1:G:129:ARG:H	1.80	0.43
2:H:47:TYR:CE2	2:H:49:GLU:HA	2.53	0.43
1:A:11:LEU:O	1:A:15:LYS:HG2	2.18	0.43
1:C:20:PHE:CE1	1:C:144:PHE:CD1	3.06	0.43
1:E:169:GLN:O	1:E:172:GLY:N	2.51	0.43
1:G:114:TRP:O	1:G:119:VAL:HG23	2.19	0.43
1:G:67:HIS:CD2	1:G:133:ALA:HB2	2.53	0.43
2:H:16:LEU:HD11	2:H:55:VAL:HG21	2.01	0.43
1:A:16:GLN:NE2	3:A:238:HOH:O	2.51	0.43
1:A:207:ASN:O	1:A:211:GLU:HG3	2.18	0.43
1:C:211:GLU:HG2	1:C:217:PRO:HG3	2.00	0.43
1:A:175:ILE:H	1:A:175:ILE:HG13	1.61	0.43
1:E:47:ALA:HA	1:E:75:LEU:CD2	2.49	0.43
2:F:47:TYR:CE2	2:F:49:GLU:HA	2.54	0.43
1:G:105:ARG:HA	1:G:106:PRO:HD3	1.68	0.43
2:H:25:LEU:N	2:H:25:LEU:HD23	2.33	0.43
1:A:97:GLU:HA	1:A:103:PHE:HD1	1.83	0.43
1:C:152:ILE:O	1:C:156:ARG:HB2	2.19	0.43
1:E:156:ARG:O	1:E:180:HIS:HE1	2.02	0.43
1:E:187:HIS:CG	1:E:188:PRO:HD2	2.54	0.43
2:B:4:LEU:HB3	2:B:7:ILE:HD12	2.01	0.43
1:C:85:ILE:HG23	1:C:89:LEU:HD23	2.00	0.43
2:F:56:MET:O	2:F:70:LEU:HD12	2.18	0.43
1:A:5:LEU:HA	1:A:5:LEU:HD23	1.87	0.43
1:E:97:GLU:HA	1:E:103:PHE:HD1	1.84	0.43
2:D:4:LEU:HA	2:D:7:ILE:HD12	2.01	0.43
1:G:168:ALA:O	1:G:171:LYS:HB2	2.19	0.43
1:G:21:LEU:O	1:G:25:GLN:HB2	2.19	0.43
2:H:80:LYS:HB2	2:H:80:LYS:HE2	1.89	0.43
1:A:39:PRO:HB2	1:A:40:PRO:HD2	2.00	0.42
1:C:200:CYS:HB2	1:C:202:HIS:CD2	2.54	0.42
1:G:163:LEU:HD13	1:G:169:GLN:HA	2.01	0.42
2:B:71:VAL:HG12	2:B:79:ASN:HB2	2.01	0.42
1:E:213:ARG:HH11	1:E:213:ARG:HB3	1.83	0.42
1:E:81:PRO:HD3	1:E:110:TYR:CD1	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:10:VAL:HG21	1:A:53:LEU:HD11	2.01	0.42
2:F:10:LYS:HA	2:F:10:LYS:HD2	1.82	0.42
1:C:197:PHE:O	1:C:200:CYS:SG	2.62	0.42
1:E:147:LYS:O	1:E:150:SER:HB3	2.19	0.42
1:E:64:ASP:HB2	1:E:65:PRO:HD2	2.01	0.42
1:G:3:ASN:HA	1:G:5:LEU:H	1.84	0.42
1:G:73:HIS:HE1	1:G:75:LEU:CD1	2.29	0.42
2:B:66:LYS:HA	2:B:67:PRO:HD3	1.94	0.42
1:E:106:PRO:O	1:E:107:ASN:CB	2.67	0.42
2:H:71:VAL:HG22	2:H:81:ILE:HG12	2.02	0.42
1:C:209:TRP:O	1:C:213:ARG:HB2	2.19	0.42
1:E:186:PRO:HG2	1:E:196:GLY:HA3	2.02	0.42
1:G:164:TRP:HA	1:G:185:ALA:O	2.20	0.42
1:A:64:ASP:HB2	1:A:65:PRO:HD2	2.02	0.41
1:C:104:THR:HG23	1:C:221:MET:HE1	2.01	0.41
1:E:7:TRP:HH2	1:E:152:ILE:HD11	1.84	0.41
1:A:211:GLU:C	1:A:213:ARG:H	2.24	0.41
1:E:174:ILE:H	1:E:174:ILE:HG13	1.60	0.41
1:E:90:LEU:HD12	1:E:90:LEU:C	2.40	0.41
1:C:142:GLU:O	1:C:146:ASP:HB2	2.20	0.41
1:C:86:PRO:HA	1:C:87:PRO:HD3	1.92	0.41
1:E:212:GLN:HG2	1:E:212:GLN:O	2.20	0.41
1:G:195:ARG:HG3	2:H:32:VAL:HG22	2.01	0.41
1:G:8:HIS:O	1:G:9:ASP:C	2.59	0.41
1:A:187:HIS:CD2	1:A:188:PRO:HD2	2.54	0.41
1:C:20:PHE:HE1	1:C:144:PHE:CD1	2.38	0.41
1:G:65:PRO:HD3	1:G:125:VAL:O	2.21	0.41
2:H:28:GLU:O	2:H:31:GLU:HB3	2.20	0.41
1:A:223:VAL:O	1:A:225:PRO:HD3	2.20	0.41
1:A:76:ALA:O	1:A:78:SER:N	2.45	0.41
2:B:30:GLU:O	2:B:34:GLY:HA2	2.21	0.41
1:E:19:TYR:CG	1:E:20:PHE:N	2.88	0.41
1:G:207:ASN:OD1	1:G:217:PRO:HB3	2.21	0.41
1:A:104:THR:O	1:A:221:MET:HE3	2.21	0.41
1:E:90:LEU:O	1:E:93:TYR:HB2	2.21	0.41
1:C:76:ALA:C	1:C:78:SER:H	2.23	0.41
1:G:100:ILE:HA	1:G:101:PRO:HD3	1.72	0.41
1:C:11:LEU:O	1:C:15:LYS:CG	2.69	0.41
1:C:21:LEU:HA	1:C:21:LEU:HD23	1.82	0.41
1:E:103:PHE:HE2	1:E:221:MET:HA	1.86	0.41
1:E:31:ARG:HG2	1:E:37:ILE:HD12	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:28:GLU:O	2:B:31:GLU:HB3	2.20	0.40
1:C:159:VAL:HB	1:C:180:HIS:ND1	2.36	0.40
1:C:19:TYR:CG	1:C:20:PHE:N	2.89	0.40
1:C:21:LEU:O	1:C:25:GLN:CG	2.69	0.40
1:C:61:LEU:HD21	1:C:149:ILE:HD11	2.03	0.40
1:G:201:ASN:O	1:G:202:HIS:C	2.59	0.40
1:G:76:ALA:O	1:G:78:SER:N	2.48	0.40
2:H:56:MET:HE3	2:H:71:VAL:HB	2.01	0.40
1:A:18:PRO:N	3:A:230:HOH:O	2.55	0.40
2:D:71:VAL:HG12	2:D:79:ASN:HB2	2.01	0.40
1:G:200:CYS:HB2	1:G:202:HIS:CD2	2.56	0.40
1:A:224:LEU:HD12	1:A:224:LEU:HA	1.93	0.40
1:A:41:GLN:CA	1:A:44:VAL:HG23	2.39	0.40
1:C:27:VAL:O	1:C:31:ARG:HG3	2.22	0.40
1:E:172:GLY:HA2	1:E:175:ILE:HD11	1.97	0.40
1:E:216:THR:HA	1:E:217:PRO:HD3	1.94	0.40
1:E:195:ARG:HG3	2:F:32:VAL:CG2	2.51	0.40
1:G:93:TYR:O	1:G:97:GLU:HG3	2.21	0.40
1:A:89:LEU:O	1:A:92:MET:HB2	2.21	0.40
1:C:5:LEU:HD23	1:C:5:LEU:HA	1.87	0.40
1:C:85:ILE:HA	1:C:86:PRO:HD3	1.90	0.40
1:G:7:TRP:O	1:G:11:LEU:HB2	2.21	0.40
2:H:16:LEU:CD1	2:H:55:VAL:HG21	2.52	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	221/229 (96%)	193 (87%)	23 (10%)	5 (2%)	6	34
1	C	222/229 (97%)	197 (89%)	20 (9%)	5 (2%)	6	34

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	222/229 (97%)	197 (89%)	23 (10%)	2 (1%)	17	56
1	G	223/229 (97%)	194 (87%)	26 (12%)	3 (1%)	12	47
2	B	80/84 (95%)	71 (89%)	9 (11%)	0	100	100
2	D	81/84 (96%)	76 (94%)	5 (6%)	0	100	100
2	F	80/84 (95%)	69 (86%)	11 (14%)	0	100	100
2	H	82/84 (98%)	72 (88%)	8 (10%)	2 (2%)	6	34
All	All	1211/1252 (97%)	1069 (88%)	125 (10%)	17 (1%)	11	46

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	71	GLN
2	H	2	THR
1	A	42	LYS
1	C	158	GLY
1	C	169	GLN
1	G	71	GLN
2	H	3	ASN
1	A	18	PRO
1	C	212	GLN
1	E	212	GLN
1	G	42	LYS
1	A	77	PHE
1	C	41	GLN
1	G	18	PRO
1	A	212	GLN
1	C	217	PRO
1	E	18	PRO

### 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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	186/194 (96%)	167 (90%)	19 (10%)	7 29
1	C	187/194 (96%)	172 (92%)	15 (8%)	12 42
1	E	187/194 (96%)	168 (90%)	19 (10%)	7 29
1	G	187/194 (96%)	168 (90%)	19 (10%)	7 29
2	B	75/78 (96%)	69 (92%)	6 (8%)	12 42
2	D	76/78 (97%)	68 (90%)	8 (10%)	7 28
2	F	76/78 (97%)	70 (92%)	6 (8%)	12 43
2	H	76/78 (97%)	67 (88%)	9 (12%)	5 23
All	All	1050/1088 (96%)	949 (90%)	101 (10%)	8 32

All (101) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	LEU
1	A	8	HIS
1	A	10	VAL
1	A	13	GLU
1	A	16	GLN
1	A	17	GLN
1	A	24	LEU
1	A	41	GLN
1	A	49	ARG
1	A	53	LEU
1	A	113	SER
1	A	134	HIS
1	A	135	SER
1	A	146	ASP
1	A	150	SER
1	A	179	ARG
1	A	181	HIS
1	A	201	ASN
1	A	216	THR
2	B	4	LEU
2	B	15	GLN
2	B	64	GLU
2	B	75	SER
2	B	79	ASN

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Mol	Chain	Res	Type
2	B	80	LYS
1	C	5	LEU
1	C	10	VAL
1	C	13	GLU
1	C	16	GLN
1	C	53	LEU
1	C	79	VAL
1	C	98	ASN
1	C	104	THR
1	C	113	SER
1	C	138	SER
1	C	157	GLU
1	C	184	LYS
1	C	201	ASN
1	C	213	ARG
1	C	216	THR
2	D	3	ASN
2	D	15	GLN
2	D	21	SER
2	D	45	THR
2	D	58	LEU
2	D	64	GLU
2	D	75	SER
2	D	79	ASN
1	E	10	VAL
1	E	16	GLN
1	E	24	LEU
1	E	25	GLN
1	E	33	SER
1	E	42	LYS
1	E	53	LEU
1	E	79	VAL
1	E	107	ASN
1	E	111	LEU
1	E	122	LEU
1	E	139	LEU
1	E	146	ASP
1	E	157	GLU
1	E	175	ILE
1	E	179	ARG
1	E	181	HIS
1	E	184	LYS

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Mol	Chain	Res	Type
1	E	216	THR
2	F	10	LYS
2	F	15	GLN
2	F	35	ASN
2	F	64	GLU
2	F	75	SER
2	F	79	ASN
1	G	5	LEU
1	G	15	LYS
1	G	16	GLN
1	G	25	GLN
1	G	33	SER
1	G	36	THR
1	G	51	THR
1	G	53	LEU
1	G	80	ARG
1	G	107	ASN
1	G	111	LEU
1	G	113	SER
1	G	122	LEU
1	G	134	HIS
1	G	139	LEU
1	G	150	SER
1	G	157	GLU
1	G	179	ARG
1	G	216	THR
2	H	4	LEU
2	H	10	LYS
2	H	15	GLN
2	H	23	LEU
2	H	45	THR
2	H	53	GLU
2	H	64	GLU
2	H	75	SER
2	H	79	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	GLN
1	A	22	ASN
1	A	41	GLN

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Mol	Chain	Res	Type
1	A	98	ASN
1	A	154	GLN
1	A	155	HIS
1	A	169	GLN
1	A	208	GLN
2	B	3	ASN
2	B	76	ASN
1	C	22	ASN
1	C	208	GLN
2	D	76	ASN
1	E	8	HIS
1	E	22	ASN
1	E	169	GLN
1	E	208	GLN
2	F	19	GLN
2	F	76	ASN
1	G	22	ASN
1	G	98	ASN
1	G	154	GLN
1	G	169	GLN
1	G	194	HIS
1	G	208	GLN
2	H	76	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

### 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

### 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

### 5.6 Ligand geometry ⓘ

There are no ligands in this entry.

## 5.7 Other polymers

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	223/229 (97%)	-0.79	0 100 100	2, 20, 54, 83	0
1	C	224/229 (97%)	-0.71	1 (0%) 92 89	2, 22, 57, 79	0
1	E	224/229 (97%)	-0.72	0 100 100	2, 21, 57, 74	0
1	G	225/229 (98%)	-0.75	3 (1%) 77 65	2, 21, 52, 71	0
2	B	82/84 (97%)	-0.61	0 100 100	2, 27, 64, 79	0
2	D	83/84 (98%)	-0.63	0 100 100	2, 30, 54, 74	0
2	F	82/84 (97%)	-0.58	0 100 100	2, 28, 59, 70	0
2	H	84/84 (100%)	-0.56	2 (2%) 59 44	2, 28, 60, 70	0
All	All	1227/1252 (98%)	-0.70	6 (0%) 91 86	2, 23, 58, 83	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	226	ALA	3.3
1	C	227	GLU	3.3
2	H	2	THR	2.8
2	H	1	MET	2.5
1	G	2	ALA	2.4
1	G	4	GLU	2.1

### 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.