



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:50 PM UTC

PDB ID : 3LMU / pdb_00003lmu
Title : Crystal structure of DTD from Plasmodium falciparum
Authors : Manickam, Y.; Bhatt, T.K.; Khan, S.; Sharma, A.
Deposited on : 2010-02-01
Resolution : 3.30 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

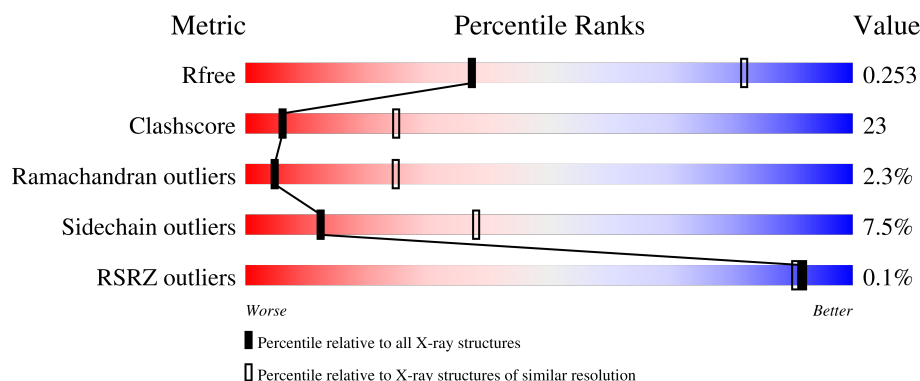
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1169 (3.32-3.28)
Clashscore	190562	1209 (3.32-3.28)
Ramachandran outliers	187476	1188 (3.32-3.28)
Sidechain outliers	187428	1187 (3.32-3.28)
RSRZ outliers	180081	1169 (3.32-3.28)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 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	164	
1	B	164	
1	C	164	
1	D	164	
1	E	164	

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Mol	Chain	Length	Quality of chain
1	F	164	
1	G	164	
1	H	164	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	IOD	D	167	-	-	X	-
2	IOD	G	165	-	-	X	-

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9279 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 D-tyrosyl-tRNA(Tyr) deacylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	155	Total	C	N	O	S	0	0	0
			1204	778	201	222	3			
1	B	148	Total	C	N	O	S	2	0	0
			1174	756	198	217	3			
1	C	148	Total	C	N	O	S	0	0	0
			1176	760	196	217	3			
1	D	148	Total	C	N	O	S	2	0	0
			1172	757	194	218	3			
1	E	145	Total	C	N	O	S	0	0	0
			1134	735	185	211	3			
1	F	144	Total	C	N	O	S	0	0	0
			1126	730	184	209	3			
1	G	148	Total	C	N	O	S	0	0	0
			1120	720	186	211	3			
1	H	148	Total	C	N	O	S	0	0	0
			1147	743	190	211	3			

- Molecule 2 is IODIDE ION (CCD ID: IOD) (formula: I).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	I	0	0
			1	1		
2	B	5	Total	I	0	0
			5	5		
2	C	4	Total	I	0	0
			4	4		
2	D	6	Total	I	0	0
			6	6		
2	E	1	Total	I	0	0
			1	1		
2	F	4	Total	I	0	0
			4	4		

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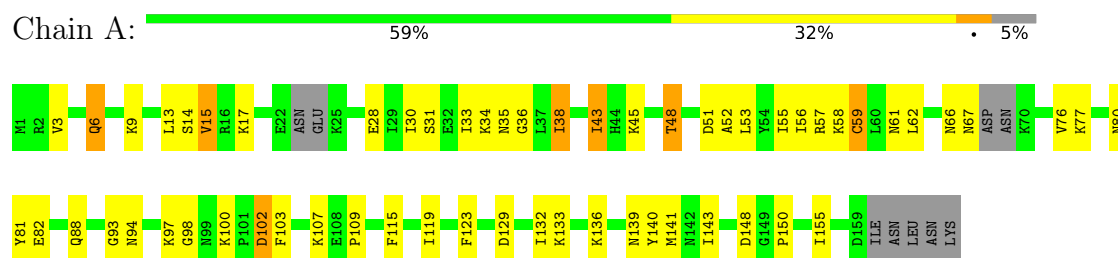
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	G	3	Total	I	0	0
			3	3		
2	H	2	Total	I	0	0
			2	2		

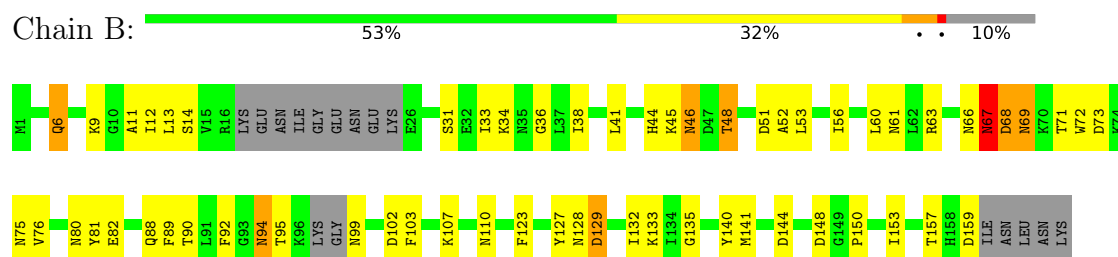
3 Residue-property plots [i](#)

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.

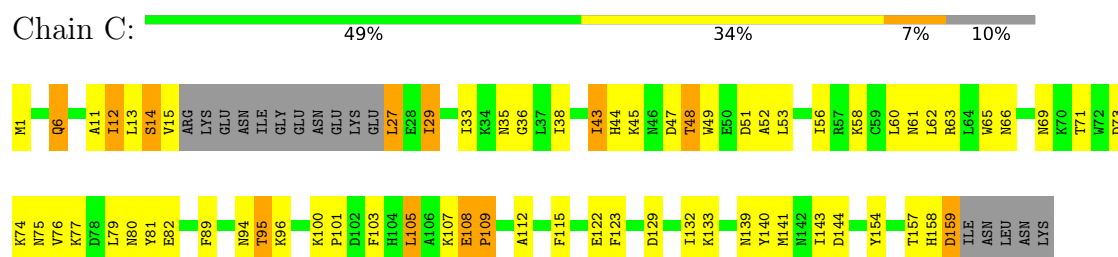
• Molecule 1: D-tyrosyl-tRNA(Tyr) deacylase



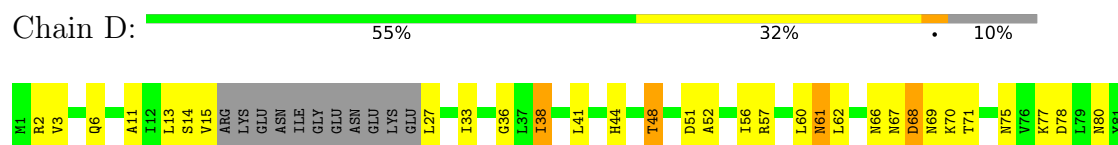
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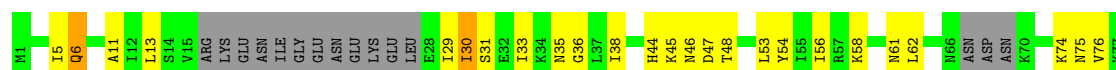




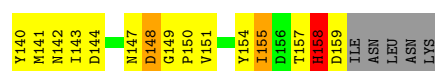
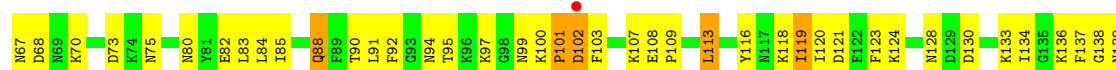
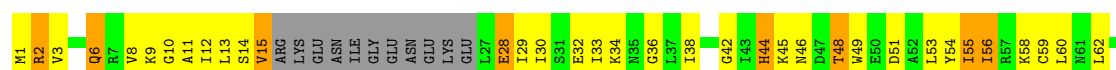
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4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	77.18Å 77.18Å 214.52Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	33.96 – 3.30 33.96 – 3.30	Depositor EDS
% Data completeness (in resolution range)	99.9 (33.96-3.30) 99.8 (33.96-3.30)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.63 (at 3.33Å)	Xtriage
Refinement program	PHENIX 1.5_2	Depositor
R, R_{free}	0.185 , 0.256 0.184 , 0.253	Depositor DCC
R_{free} test set	971 reflections (5.15%)	wwPDB-VP
Wilson B-factor (Å ²)	56.9	Xtriage
Anisotropy	0.295	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 42.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.045 for h,-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	9279	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.11% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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 ⓘ

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

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	0.74	5/1225 (0.4%)	0.92	1/1661 (0.1%)
1	B	0.79	5/1195 (0.4%)	0.89	0/1621
1	C	0.72	5/1199 (0.4%)	0.90	3/1628 (0.2%)
1	D	0.73	2/1195 (0.2%)	0.94	3/1623 (0.2%)
1	E	0.76	7/1154 (0.6%)	0.89	2/1567 (0.1%)
1	F	0.68	4/1148 (0.3%)	0.91	3/1560 (0.2%)
1	G	0.79	7/1141 (0.6%)	0.99	8/1560 (0.5%)
1	H	0.75	5/1169 (0.4%)	0.97	0/1591
All	All	0.75	40/9426 (0.4%)	0.93	20/12811 (0.2%)

All (40) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	35	ASN	CG-OD1	5.86	1.34	1.23
1	C	69	ASN	CG-OD1	5.85	1.34	1.23
1	B	94	ASN	CG-OD1	5.78	1.34	1.23
1	E	88	GLN	CD-OE1	5.78	1.34	1.23
1	E	6	GLN	CD-OE1	5.77	1.34	1.23
1	E	67	ASN	CG-OD1	5.68	1.34	1.23
1	F	6	GLN	CD-OE1	5.65	1.34	1.23
1	A	94	ASN	CG-OD1	5.57	1.34	1.23
1	G	158	HIS	ND1-CE1	5.52	1.38	1.32
1	C	6	GLN	CD-OE1	5.47	1.33	1.23
1	G	6	GLN	CD-OE1	5.47	1.33	1.23
1	H	6	GLN	CD-OE1	5.45	1.33	1.23
1	G	94	ASN	CG-OD1	5.45	1.33	1.23
1	B	6	GLN	CD-OE1	5.43	1.33	1.23
1	G	88	GLN	CD-OE1	5.41	1.33	1.23
1	C	139	ASN	CG-OD1	5.40	1.33	1.23
1	H	88	GLN	CD-OE1	5.36	1.33	1.23
1	A	35	ASN	CG-ND2	-5.36	1.22	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	67	ASN	CG-OD1	5.35	1.33	1.23
1	F	158	HIS	ND1-CE1	5.32	1.37	1.32
1	H	80	ASN	CG-OD1	5.28	1.33	1.23
1	C	35	ASN	CG-OD1	5.27	1.33	1.23
1	H	99	ASN	CG-OD1	5.27	1.33	1.23
1	G	147	ASN	CG-OD1	5.26	1.33	1.23
1	H	147	ASN	CG-OD1	5.25	1.33	1.23
1	G	44	HIS	ND1-CE1	5.24	1.37	1.32
1	B	46	ASN	CG-OD1	5.17	1.33	1.23
1	A	139	ASN	CG-OD1	5.16	1.33	1.23
1	D	61	ASN	CG-OD1	5.16	1.33	1.23
1	A	6	GLN	CD-OE1	5.15	1.33	1.23
1	B	110	ASN	CG-OD1	5.14	1.33	1.23
1	E	139	ASN	CG-OD1	5.13	1.33	1.23
1	D	139	ASN	CG-OD1	5.11	1.33	1.23
1	F	139	ASN	CG-OD1	5.10	1.33	1.23
1	A	6	GLN	CD-NE2	-5.08	1.22	1.33
1	E	35	ASN	CG-OD1	5.08	1.33	1.23
1	E	35	ASN	CG-ND2	-5.07	1.22	1.33
1	G	67	ASN	CG-OD1	5.05	1.33	1.23
1	E	94	ASN	CG-OD1	5.04	1.33	1.23
1	C	35	ASN	CG-ND2	-5.04	1.22	1.33

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	100	LYS	CA-C-N	7.85	129.66	119.84
1	G	100	LYS	C-N-CA	7.85	129.66	119.84
1	D	100	LYS	CA-C-N	7.16	128.79	119.84
1	D	100	LYS	C-N-CA	7.16	128.79	119.84
1	G	158	HIS	CB-CG-CD2	-6.55	122.68	131.20
1	G	44	HIS	CB-CG-CD2	-6.52	122.72	131.20
1	E	100	LYS	CA-C-N	6.45	127.91	119.84
1	E	100	LYS	C-N-CA	6.45	127.91	119.84
1	C	108	GLU	CA-C-N	6.44	127.89	119.84
1	C	108	GLU	C-N-CA	6.44	127.89	119.84
1	D	68	ASP	N-CA-C	-6.31	102.19	110.33
1	F	158	HIS	CB-CG-CD2	-6.28	123.03	131.20
1	G	28	GLU	N-CA-C	5.82	117.88	108.34
1	G	158	HIS	CB-CG-ND1	5.80	131.40	122.70
1	G	44	HIS	CB-CG-ND1	5.69	131.23	122.70
1	A	143	ILE	N-CA-C	5.61	115.93	107.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	143	ILE	N-CA-C	5.59	115.90	107.80
1	F	158	HIS	CB-CG-ND1	5.55	131.03	122.70
1	G	9	LYS	N-CA-C	-5.33	106.28	112.89
1	C	143	ILE	N-CA-C	5.06	115.13	107.80

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	1204	0	1147	35	0
1	B	1174	0	1126	40	0
1	C	1176	0	1143	55	0
1	D	1172	0	1127	47	0
1	E	1134	0	1080	50	0
1	F	1126	0	1065	53	0
1	G	1120	0	1029	80	0
1	H	1147	0	1093	76	0
2	A	1	0	0	0	0
2	B	5	0	0	0	0
2	C	4	0	0	1	0
2	D	6	0	0	4	0
2	E	1	0	0	0	0
2	F	4	0	0	0	0
2	G	3	0	0	4	0
2	H	2	0	0	2	0
All	All	9279	0	8810	409	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 23.

All (409) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:99:ASN:OD1	2:D:167:IOD:I	2.19	1.30
1:D:99:ASN:CG	2:D:167:IOD:I	2.54	1.19
1:D:99:ASN:ND2	2:D:167:IOD:I	2.49	1.15
1:E:61:ASN:HA	1:E:75:ASN:HD22	1.22	1.03
1:H:9:LYS:HG2	1:H:148:ASP:HB2	1.39	1.03
1:G:6:GLN:HB2	1:G:38:ILE:HG22	1.48	0.95
1:G:33:ILE:HG22	1:G:133:LYS:HG3	1.49	0.94
1:G:56:ILE:HD11	1:G:119:ILE:HG22	1.51	0.93
1:H:90:THR:HA	1:H:103:PHE:HE1	1.33	0.91
1:C:105:LEU:HD12	1:C:105:LEU:H	1.35	0.91
1:B:94:ASN:HB3	1:B:102:ASP:HB3	1.53	0.90
1:D:68:ASP:O	1:D:70:LYS:N	2.07	0.87
1:H:90:THR:HA	1:H:103:PHE:CE1	2.13	0.83
1:E:105:LEU:HD12	1:E:105:LEU:H	1.47	0.79
1:F:107:LYS:HG2	1:F:112:ALA:HB2	1.65	0.79
1:G:141:MET:HE1	2:G:165:IOD:I	2.52	0.79
1:G:99:ASN:OD1	1:H:58:LYS:NZ	2.18	0.77
1:G:82:GLU:C	1:G:83:LEU:HD12	2.11	0.75
1:D:68:ASP:C	1:D:70:LYS:H	1.95	0.75
1:E:61:ASN:HA	1:E:75:ASN:ND2	2.01	0.75
1:G:6:GLN:HB2	1:G:38:ILE:CG2	2.16	0.74
1:C:94:ASN:HD21	1:C:96:LYS:HE3	1.54	0.73
1:F:76:VAL:HG23	1:F:81:TYR:HB2	1.69	0.72
1:A:43:ILE:HD12	1:A:43:ILE:H	1.53	0.71
1:H:5:ILE:HB	1:H:153:ILE:HB	1.72	0.71
1:C:12:ILE:HG22	1:C:29:ILE:HG12	1.72	0.70
1:G:56:ILE:HD11	1:G:119:ILE:CG2	2.21	0.70
1:C:43:ILE:H	1:C:43:ILE:HD12	1.55	0.70
1:D:93:GLY:HA2	1:D:102:ASP:O	1.92	0.70
1:H:56:ILE:HD11	1:H:119:ILE:HG23	1.74	0.69
1:C:105:LEU:HD12	1:C:105:LEU:N	2.08	0.69
1:G:116:TYR:O	1:G:120:ILE:HG12	1.93	0.69
1:A:76:VAL:HG23	1:A:81:TYR:HB2	1.75	0.68
1:C:14:SER:HB3	1:C:27:LEU:HD22	1.73	0.68
1:A:15:VAL:HG22	1:A:30:ILE:HD13	1.74	0.68
1:F:44:HIS:HE2	1:F:105:LEU:HD23	1.59	0.68
1:G:157:THR:C	1:G:159:ASP:H	2.02	0.68
1:G:2:ARG:HG3	1:G:92:PHE:CZ	2.30	0.67
1:H:9:LYS:CG	1:H:148:ASP:HB2	2.22	0.67
1:C:49:TRP:HB2	2:C:168:IOD:I	2.65	0.67
1:G:2:ARG:HG3	1:G:92:PHE:CE2	2.29	0.67
1:G:137:PHE:CD2	1:G:138:GLY:N	2.62	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:137:PHE:HD2	1:G:138:GLY:N	1.94	0.66
1:A:15:VAL:HG11	1:A:136:LYS:HB3	1.77	0.66
1:A:52:ALA:O	1:A:56:ILE:HG13	1.96	0.66
1:H:44:HIS:HE1	1:H:46:ASN:HB2	1.62	0.65
1:G:44:HIS:HD2	1:G:46:ASN:H	1.44	0.65
1:H:7:ARG:HB2	1:H:151:VAL:HB	1.79	0.65
1:G:84:LEU:HD12	1:G:133:LYS:HB2	1.79	0.64
1:E:99:ASN:O	1:F:62:LEU:HD23	1.96	0.64
1:G:136:LYS:O	1:G:139:ASN:HB2	1.98	0.64
1:C:105:LEU:H	1:C:105:LEU:CD1	2.09	0.62
1:H:71:THR:HG1	1:H:72:TRP:CD1	2.16	0.62
1:B:61:ASN:HA	1:B:75:ASN:HD22	1.64	0.62
1:H:69:ASN:OD1	2:H:166:IOD:I	2.87	0.62
1:G:33:ILE:CG2	1:G:133:LYS:HG3	2.27	0.62
1:F:90:THR:HA	1:F:103:PHE:CE1	2.34	0.62
1:H:76:VAL:HG23	1:H:81:TYR:HB2	1.82	0.62
1:A:45:LYS:HA	1:A:107:LYS:HB2	1.82	0.61
1:E:58:LYS:HD2	1:E:62:LEU:HD21	1.82	0.61
1:E:94:ASN:OD1	1:E:94:ASN:C	2.43	0.61
1:E:150:PRO:HB2	1:F:88:GLN:OE1	1.99	0.61
1:H:44:HIS:CE1	1:H:46:ASN:H	2.18	0.61
1:H:123:PHE:HB3	1:H:132:ILE:HD12	1.82	0.61
1:B:52:ALA:O	1:B:56:ILE:HG13	2.00	0.61
1:H:15:VAL:HG12	1:H:140:TYR:O	2.00	0.61
1:H:44:HIS:CE1	1:H:46:ASN:HB2	2.35	0.61
1:F:45:LYS:HA	1:F:107:LYS:HE3	1.81	0.61
1:H:136:LYS:HG2	1:H:139:ASN:HB2	1.83	0.61
1:A:6:GLN:HB2	1:A:38:ILE:HG22	1.83	0.60
1:D:13:LEU:C	1:D:13:LEU:HD23	2.27	0.60
1:E:95:THR:HG21	1:F:62:LEU:HD11	1.83	0.60
1:F:45:LYS:HG3	1:F:46:ASN:OD1	2.02	0.60
1:E:108:GLU:HG2	1:E:109:PRO:HD2	1.83	0.60
1:F:54:TYR:O	1:F:58:LYS:HG2	2.00	0.60
1:H:11:ALA:HA	1:H:144:ASP:O	2.01	0.60
1:F:74:LYS:HB2	1:F:79:LEU:HD11	1.83	0.60
1:D:94:ASN:HB3	1:D:102:ASP:HB3	1.84	0.60
1:C:6:GLN:HB2	1:C:38:ILE:HG22	1.83	0.59
1:H:41:LEU:HD12	1:H:42:GLY:N	2.17	0.59
1:H:6:GLN:HB2	1:H:38:ILE:HG22	1.83	0.59
1:F:6:GLN:HB2	1:F:38:ILE:HG22	1.84	0.59
1:C:76:VAL:HG23	1:C:81:TYR:HB2	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:99:ASN:HD22	1:F:61:ASN:HB3	1.68	0.59
1:C:43:ILE:H	1:C:43:ILE:CD1	2.13	0.59
1:A:15:VAL:CG2	1:A:30:ILE:HG21	2.34	0.58
1:G:44:HIS:CD2	1:G:46:ASN:H	2.20	0.58
1:G:116:TYR:CZ	2:G:165:IOD:I	3.26	0.58
1:G:120:ILE:O	1:G:124:LYS:HG3	2.03	0.58
1:H:107:LYS:HG2	1:H:112:ALA:HB2	1.85	0.58
1:C:94:ASN:ND2	1:C:96:LYS:HE3	2.18	0.58
1:A:93:GLY:HA2	1:A:103:PHE:CD1	2.38	0.58
1:C:52:ALA:O	1:C:56:ILE:HG13	2.04	0.58
1:D:108:GLU:OE1	1:D:110:ASN:HB2	2.04	0.57
1:G:154:TYR:O	1:G:155:ILE:HG12	2.04	0.57
1:F:61:ASN:HA	1:F:75:ASN:ND2	2.20	0.57
1:C:103:PHE:CD2	1:C:103:PHE:N	2.73	0.57
1:H:54:TYR:CD2	1:H:157:THR:HG22	2.39	0.57
1:C:6:GLN:HB2	1:C:38:ILE:CG2	2.33	0.57
1:H:6:GLN:HB2	1:H:38:ILE:CG2	2.34	0.57
1:E:42:GLY:C	1:E:92:PHE:HE2	2.13	0.57
1:D:6:GLN:HB2	1:D:38:ILE:HG22	1.87	0.56
1:G:3:VAL:HG12	1:G:155:ILE:HB	1.88	0.56
1:D:68:ASP:C	1:D:70:LYS:N	2.62	0.56
1:A:15:VAL:HG13	1:A:141:MET:HG2	1.87	0.56
1:B:13:LEU:HD22	1:B:135:GLY:HA2	1.88	0.56
1:D:14:SER:HB3	1:D:27:LEU:HD22	1.88	0.56
1:F:6:GLN:HB2	1:F:38:ILE:CG2	2.35	0.55
1:F:108:GLU:HG2	1:F:109:PRO:HD2	1.87	0.55
1:G:140:TYR:C	1:G:140:TYR:CD2	2.84	0.55
1:G:154:TYR:CZ	1:H:154:TYR:CE2	2.95	0.55
1:A:6:GLN:HB2	1:A:38:ILE:CG2	2.36	0.55
1:B:6:GLN:HB2	1:B:38:ILE:HG22	1.87	0.55
1:E:123:PHE:HB3	1:E:132:ILE:HG13	1.89	0.55
1:G:157:THR:O	1:G:159:ASP:N	2.34	0.55
1:H:42:GLY:C	1:H:92:PHE:HE2	2.15	0.55
1:C:14:SER:HB3	1:C:27:LEU:CD2	2.37	0.55
1:G:88:GLN:HA	1:G:88:GLN:OE1	2.06	0.55
1:A:15:VAL:CG2	1:A:30:ILE:HD13	2.36	0.55
1:E:158:HIS:O	1:E:159:ASP:OD1	2.25	0.55
1:G:13:LEU:HD11	1:G:141:MET:SD	2.47	0.55
1:D:56:ILE:HG22	1:D:60:LEU:HD12	1.89	0.54
1:E:6:GLN:HB2	1:E:38:ILE:HG22	1.89	0.54
1:E:105:LEU:HD12	1:E:105:LEU:N	2.20	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:57:ARG:O	1:D:61:ASN:HB2	2.08	0.54
1:D:6:GLN:HB2	1:D:38:ILE:CG2	2.37	0.54
1:G:44:HIS:HD1	1:G:92:PHE:HE1	1.52	0.54
1:H:100:LYS:HG3	1:H:101:PRO:CD	2.38	0.54
1:C:140:TYR:C	1:C:140:TYR:CD2	2.84	0.54
1:G:113:LEU:O	1:G:113:LEU:HD12	2.07	0.53
1:H:87:SER:HB3	2:H:165:IOD:I	2.78	0.53
1:G:85:ILE:N	1:G:85:ILE:HD13	2.23	0.53
1:D:75:ASN:C	1:D:75:ASN:OD1	2.52	0.53
1:E:88:GLN:NE2	1:F:150:PRO:HB2	2.23	0.53
1:A:123:PHE:HB3	1:A:132:ILE:HG13	1.91	0.53
1:F:105:LEU:N	1:F:105:LEU:HD12	2.24	0.53
1:H:149:GLY:HA3	1:H:150:PRO:C	2.33	0.53
1:E:89:PHE:CD1	1:H:69:ASN:ND2	2.76	0.53
1:F:140:TYR:C	1:F:140:TYR:CD2	2.86	0.53
1:G:116:TYR:OH	2:G:165:IOD:I	2.90	0.53
1:E:62:LEU:HD11	1:F:95:THR:HG21	1.91	0.53
1:B:6:GLN:HB2	1:B:38:ILE:CG2	2.39	0.53
1:F:90:THR:HA	1:F:103:PHE:HE1	1.73	0.53
1:G:12:ILE:HG21	1:G:29:ILE:HD13	1.89	0.53
1:B:123:PHE:HB3	1:B:132:ILE:HG13	1.90	0.53
1:G:13:LEU:HD23	1:G:14:SER:N	2.23	0.53
1:B:44:HIS:HB2	1:B:92:PHE:CZ	2.44	0.52
1:D:105:LEU:HD13	2:D:170:IOD:I	2.79	0.52
1:F:76:VAL:HG23	1:F:81:TYR:CB	2.36	0.52
1:A:13:LEU:HD21	1:A:141:MET:SD	2.49	0.52
1:G:13:LEU:HD23	1:G:13:LEU:C	2.33	0.52
1:E:6:GLN:HB2	1:E:38:ILE:CG2	2.39	0.52
1:F:44:HIS:HB2	1:F:92:PHE:CZ	2.44	0.52
1:F:108:GLU:OE2	1:F:110:ASN:HB2	2.10	0.52
1:E:64:LEU:HD23	1:E:72:TRP:HZ3	1.74	0.52
1:G:11:ALA:HA	1:G:144:ASP:O	2.10	0.52
1:A:88:GLN:OE1	1:B:150:PRO:HB2	2.10	0.52
1:F:123:PHE:HB3	1:F:132:ILE:HG13	1.92	0.52
1:H:100:LYS:HG3	1:H:101:PRO:HD2	1.91	0.52
1:H:157:THR:C	1:H:159:ASP:H	2.17	0.52
1:C:63:ARG:HH21	1:C:73:ASP:HA	1.75	0.51
1:D:60:LEU:O	1:D:77:LYS:HE2	2.10	0.51
1:B:63:ARG:HG2	1:B:73:ASP:O	2.11	0.51
1:E:44:HIS:HB2	1:E:92:PHE:CZ	2.46	0.51
1:H:7:ARG:CB	1:H:151:VAL:HB	2.40	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:55:ILE:O	1:A:59:CYS:HB2	2.10	0.51
1:E:156:ASP:O	1:E:158:HIS:N	2.44	0.51
1:A:53:LEU:HA	1:A:56:ILE:HD12	1.93	0.51
1:A:66:ASN:O	1:A:67:ASN:OD1	2.29	0.51
1:B:11:ALA:HA	1:B:144:ASP:O	2.11	0.51
1:F:107:LYS:O	1:F:108:GLU:C	2.54	0.51
1:B:44:HIS:HE1	1:B:46:ASN:OD1	1.94	0.51
1:E:140:TYR:CD2	1:E:140:TYR:C	2.88	0.51
1:B:67:ASN:O	1:B:68:ASP:O	2.28	0.50
1:E:52:ALA:O	1:E:56:ILE:HG13	2.12	0.50
1:H:148:ASP:OD1	1:H:148:ASP:C	2.55	0.50
1:G:157:THR:C	1:G:159:ASP:N	2.69	0.50
1:G:48:THR:O	1:G:51:ASP:HB2	2.11	0.50
1:H:44:HIS:HB2	1:H:92:PHE:CZ	2.47	0.50
1:G:116:TYR:CE2	2:G:165:IOD:I	3.35	0.50
1:H:49:TRP:NE1	1:H:53:LEU:HD11	2.27	0.50
1:E:63:ARG:HG3	1:F:99:ASN:O	2.12	0.49
1:F:58:LYS:O	1:F:62:LEU:HB2	2.11	0.49
1:F:107:LYS:HG2	1:F:112:ALA:CB	2.40	0.49
1:G:136:LYS:HG2	1:G:139:ASN:OD1	2.12	0.49
1:C:61:ASN:HB3	1:D:99:ASN:HD22	1.76	0.49
1:B:13:LEU:HB3	1:B:31:SER:HB3	1.94	0.49
1:F:13:LEU:HB3	1:F:31:SER:HB3	1.94	0.49
1:F:56:ILE:CD1	1:F:122:GLU:HB3	2.43	0.49
1:B:33:ILE:CG2	1:B:133:LYS:HG3	2.42	0.49
1:B:140:TYR:C	1:B:140:TYR:CD2	2.91	0.49
1:D:123:PHE:HB3	1:D:132:ILE:HG13	1.93	0.49
1:H:127:TYR:CD1	1:H:127:TYR:C	2.87	0.49
1:B:13:LEU:HD21	1:B:141:MET:SD	2.52	0.48
1:C:157:THR:C	1:C:159:ASP:H	2.21	0.48
1:G:2:ARG:HA	1:G:155:ILE:O	2.13	0.48
1:C:45:LYS:HA	1:C:107:LYS:HE3	1.96	0.48
1:C:61:ASN:HB3	1:D:99:ASN:ND2	2.28	0.48
1:E:36:GLY:HA3	1:E:82:GLU:O	2.13	0.48
1:C:13:LEU:HD23	1:C:13:LEU:C	2.38	0.48
1:B:103:PHE:N	1:B:103:PHE:CD2	2.81	0.48
1:F:29:ILE:HG22	1:F:30:ILE:N	2.29	0.48
1:G:12:ILE:CG2	1:G:29:ILE:HD13	2.44	0.48
1:B:60:LEU:HD11	1:B:127:TYR:HB2	1.94	0.48
1:H:38:ILE:HG23	1:H:38:ILE:O	2.12	0.48
1:H:75:ASN:O	1:H:78:ASP:HB2	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:127:TYR:CZ	1:H:131:LYS:HG3	2.49	0.48
1:C:61:ASN:CB	1:D:99:ASN:ND2	2.77	0.48
1:E:42:GLY:O	1:E:92:PHE:HE2	1.97	0.48
1:E:156:ASP:O	1:E:157:THR:C	2.57	0.48
1:F:107:LYS:HE3	1:F:107:LYS:HB2	1.56	0.48
1:B:89:PHE:CE1	1:B:103:PHE:CD1	3.01	0.47
1:C:75:ASN:OD1	1:C:75:ASN:C	2.56	0.47
1:H:94:ASN:OD1	1:H:96:LYS:HE3	2.14	0.47
1:D:140:TYR:C	1:D:140:TYR:CD2	2.91	0.47
1:A:58:LYS:O	1:A:62:LEU:HG	2.14	0.47
1:E:94:ASN:OD1	1:E:95:THR:N	2.47	0.47
1:G:10:GLY:HA2	1:G:34:LYS:O	2.15	0.47
1:A:76:VAL:HG23	1:A:81:TYR:CB	2.43	0.47
1:G:149:GLY:HA3	1:G:150:PRO:C	2.39	0.47
1:G:84:LEU:CD1	1:G:133:LYS:HB2	2.43	0.47
1:A:13:LEU:HB3	1:A:31:SER:HB3	1.96	0.47
1:C:1:MET:H3	1:C:157:THR:HG1	1.59	0.47
1:C:36:GLY:HA3	1:C:82:GLU:O	2.14	0.47
1:C:158:HIS:C	1:C:159:ASP:OD1	2.58	0.47
1:G:150:PRO:O	1:G:151:VAL:HG23	2.15	0.47
1:G:150:PRO:O	1:G:151:VAL:CG2	2.63	0.47
1:H:113:LEU:O	1:H:117:ASN:HB2	2.15	0.47
1:D:56:ILE:CD1	1:D:122:GLU:HB3	2.45	0.47
1:B:71:THR:O	1:B:72:TRP:HB2	2.15	0.47
1:C:56:ILE:HG22	1:C:60:LEU:HD12	1.96	0.47
1:C:123:PHE:HB3	1:C:132:ILE:HG13	1.96	0.47
1:F:33:ILE:CG2	1:F:133:LYS:HG3	2.45	0.47
1:F:56:ILE:HD13	1:F:122:GLU:HB3	1.96	0.47
1:A:36:GLY:HA3	1:A:82:GLU:O	2.15	0.47
1:C:154:TYR:CE2	1:D:154:TYR:CZ	3.03	0.47
1:D:11:ALA:HA	1:D:144:ASP:O	2.15	0.47
1:A:140:TYR:C	1:A:140:TYR:CD2	2.92	0.46
1:B:157:THR:C	1:B:159:ASP:H	2.23	0.46
1:D:2:ARG:NH1	1:D:92:PHE:CD1	2.82	0.46
1:G:33:ILE:HB	1:G:82:GLU:HB2	1.97	0.46
1:C:33:ILE:CG2	1:C:133:LYS:HG3	2.45	0.46
1:G:15:VAL:HG22	1:G:28:GLU:O	2.16	0.46
1:H:123:PHE:HB3	1:H:132:ILE:CD1	2.46	0.46
1:F:122:GLU:OE1	1:F:122:GLU:HA	2.16	0.46
1:B:76:VAL:HG23	1:B:81:TYR:HB2	1.98	0.46
1:H:42:GLY:C	1:H:92:PHE:CE2	2.94	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:154:TYR:CZ	1:D:154:TYR:CE2	3.03	0.46
1:G:148:ASP:C	1:G:148:ASP:OD1	2.58	0.46
1:G:62:LEU:HD23	1:G:62:LEU:HA	1.75	0.46
1:H:94:ASN:HD21	1:H:96:LYS:HB2	1.81	0.46
1:B:90:THR:HG22	1:B:103:PHE:HE1	1.81	0.46
1:G:75:ASN:OD1	1:G:75:ASN:C	2.58	0.46
1:A:98:GLY:C	1:A:100:LYS:H	2.24	0.45
1:E:13:LEU:HD21	1:E:141:MET:SD	2.56	0.45
1:G:107:LYS:O	1:G:108:GLU:C	2.59	0.45
1:B:67:ASN:O	1:B:68:ASP:C	2.58	0.45
1:D:52:ALA:O	1:D:56:ILE:HG13	2.15	0.45
1:G:12:ILE:HG13	1:G:32:GLU:HG3	1.98	0.45
1:G:116:TYR:CE1	1:G:120:ILE:HD11	2.51	0.45
1:G:128:ASN:HD21	1:G:130:ASP:HB2	1.81	0.45
1:C:107:LYS:HG2	1:C:112:ALA:HB2	1.97	0.45
1:D:94:ASN:O	1:D:101:PRO:HA	2.16	0.45
1:G:54:TYR:O	1:G:55:ILE:C	2.59	0.45
1:G:8:VAL:HG22	1:G:36:GLY:O	2.16	0.45
1:D:75:ASN:O	1:D:78:ASP:HB2	2.16	0.45
1:F:103:PHE:O	1:F:106:ALA:HB3	2.16	0.45
1:G:151:VAL:HG13	1:H:90:THR:HG21	1.97	0.45
1:D:36:GLY:HA3	1:D:82:GLU:O	2.16	0.45
1:A:150:PRO:HB2	1:B:88:GLN:OE1	2.17	0.45
1:C:58:LYS:NZ	1:D:98:GLY:O	2.50	0.45
1:E:44:HIS:O	1:E:47:ASP:HB2	2.17	0.45
1:E:128:ASN:OD1	1:E:129:ASP:N	2.50	0.45
1:G:158:HIS:N	1:G:158:HIS:CD2	2.83	0.45
1:H:52:ALA:HB1	1:H:119:ILE:HG12	1.97	0.45
1:H:58:LYS:HA	1:H:58:LYS:HD3	1.63	0.45
1:B:53:LEU:HA	1:B:56:ILE:HD12	1.99	0.45
1:F:75:ASN:O	1:F:78:ASP:HB2	2.17	0.44
1:F:128:ASN:OD1	1:F:129:ASP:N	2.50	0.44
1:G:88:GLN:HE21	1:H:152:THR:CB	2.30	0.44
1:H:79:LEU:HD13	1:H:81:TYR:CE2	2.52	0.44
1:C:43:ILE:HG21	1:C:115:PHE:HD1	1.82	0.44
1:G:83:LEU:HD12	1:G:83:LEU:N	2.32	0.44
1:H:44:HIS:O	1:H:45:LYS:C	2.60	0.44
1:D:14:SER:CB	1:D:27:LEU:HD22	2.46	0.44
1:F:36:GLY:HA3	1:F:82:GLU:O	2.18	0.44
1:G:49:TRP:NE1	1:G:53:LEU:HD21	2.32	0.44
1:D:13:LEU:HD23	1:D:14:SER:N	2.33	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:49:TRP:HE1	1:H:53:LEU:HD11	1.82	0.44
1:A:57:ARG:O	1:A:61:ASN:HB2	2.17	0.44
1:D:93:GLY:HA2	1:D:103:PHE:CD1	2.52	0.44
1:F:13:LEU:C	1:F:13:LEU:HD23	2.43	0.44
1:C:66:ASN:OD1	1:C:71:THR:HA	2.18	0.44
1:D:27:LEU:HD23	1:D:27:LEU:HA	1.71	0.44
1:G:99:ASN:HD22	1:H:61:ASN:CB	2.30	0.44
1:E:109:PRO:HG2	1:H:70:LYS:HA	2.00	0.44
1:G:13:LEU:HD23	1:G:14:SER:O	2.17	0.44
1:G:99:ASN:HD22	1:H:61:ASN:HB2	1.83	0.44
1:H:88:GLN:HG3	1:H:91:LEU:HG	1.99	0.44
1:D:44:HIS:HB2	1:D:92:PHE:CZ	2.53	0.43
1:E:64:LEU:HD22	1:E:151:VAL:HG11	2.00	0.43
1:F:94:ASN:HB3	1:F:102:ASP:HB3	2.00	0.43
1:G:55:ILE:O	1:G:56:ILE:C	2.60	0.43
1:H:136:LYS:HG2	1:H:139:ASN:CB	2.48	0.43
1:C:89:PHE:CE1	1:C:103:PHE:CD1	3.06	0.43
1:D:33:ILE:CG2	1:D:133:LYS:HG3	2.48	0.43
1:E:62:LEU:CD1	1:F:95:THR:HG21	2.49	0.43
1:G:154:TYR:CZ	1:H:154:TYR:CD2	3.06	0.43
1:H:56:ILE:HG23	1:H:123:PHE:CE1	2.53	0.43
1:E:105:LEU:H	1:E:105:LEU:CD1	2.25	0.43
1:C:53:LEU:HD23	1:C:53:LEU:HA	1.85	0.43
1:E:100:LYS:HA	1:E:101:PRO:HD3	1.80	0.43
1:H:123:PHE:C	1:H:125:LYS:N	2.77	0.43
1:B:128:ASN:OD1	1:B:129:ASP:N	2.51	0.43
1:H:156:ASP:O	1:H:158:HIS:N	2.52	0.43
1:A:33:ILE:CG2	1:A:133:LYS:HG3	2.49	0.43
1:F:100:LYS:HA	1:F:101:PRO:HD3	1.83	0.43
1:C:108:GLU:HA	1:C:109:PRO:HD2	1.76	0.43
1:E:64:LEU:HD23	1:E:72:TRP:CZ3	2.53	0.43
1:A:43:ILE:HG21	1:A:115:PHE:HD1	1.84	0.43
1:C:48:THR:HG23	1:C:51:ASP:OD2	2.19	0.43
1:E:9:LYS:HG2	1:E:148:ASP:HB2	2.00	0.43
1:B:9:LYS:HG2	1:B:148:ASP:HB2	2.00	0.42
1:B:61:ASN:HA	1:B:75:ASN:ND2	2.32	0.42
1:D:66:ASN:OD1	1:D:71:THR:HA	2.19	0.42
1:E:53:LEU:HD23	1:E:53:LEU:HA	1.85	0.42
1:F:11:ALA:HA	1:F:144:ASP:O	2.19	0.42
1:G:1:MET:HA	1:G:42:GLY:O	2.19	0.42
1:C:122:GLU:OE1	1:C:122:GLU:HA	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:101:PRO:O	1:G:102:ASP:HB2	2.19	0.42
1:H:94:ASN:ND2	1:H:96:LYS:H	2.17	0.42
1:H:153:ILE:HA	1:H:153:ILE:HD13	1.79	0.42
1:B:45:LYS:HA	1:B:107:LYS:HB2	2.02	0.42
1:D:118:LYS:HE3	1:D:118:LYS:HB2	1.87	0.42
1:B:33:ILE:HG22	1:B:133:LYS:HG3	2.01	0.42
1:C:62:LEU:HA	1:C:62:LEU:HD23	1.82	0.42
1:G:56:ILE:HG23	1:G:123:PHE:CD1	2.54	0.42
1:G:70:LYS:O	1:G:73:ASP:HB2	2.19	0.42
1:B:9:LYS:HE3	1:B:148:ASP:CG	2.44	0.42
1:B:53:LEU:HA	1:B:53:LEU:HD23	1.81	0.42
1:F:53:LEU:HD23	1:F:53:LEU:HA	1.86	0.42
1:H:14:SER:HB2	1:H:27:LEU:CD2	2.49	0.42
1:D:66:ASN:OD1	1:D:71:THR:HB	2.19	0.42
1:E:56:ILE:HD11	1:E:119:ILE:HG23	2.01	0.42
1:A:14:SER:HA	1:A:28:GLU:O	2.20	0.42
1:C:11:ALA:HA	1:C:144:ASP:O	2.20	0.42
1:E:33:ILE:CG2	1:E:133:LYS:HG3	2.49	0.42
1:E:158:HIS:C	1:E:159:ASP:OD1	2.63	0.42
1:G:90:THR:HG22	1:G:103:PHE:CE1	2.55	0.42
1:G:92:PHE:O	1:G:103:PHE:HA	2.20	0.42
1:H:88:GLN:OE1	1:H:88:GLN:HA	2.19	0.42
1:D:48:THR:HG23	1:D:51:ASP:OD2	2.20	0.42
1:C:100:LYS:HA	1:C:101:PRO:HD3	1.82	0.41
1:E:63:ARG:HB3	1:E:73:ASP:O	2.19	0.41
1:H:54:TYR:HD2	1:H:157:THR:HG22	1.82	0.41
1:H:56:ILE:CD1	1:H:119:ILE:HG23	2.48	0.41
1:H:123:PHE:O	1:H:124:LYS:C	2.62	0.41
1:B:48:THR:HG23	1:B:51:ASP:OD2	2.20	0.41
1:E:58:LYS:O	1:E:62:LEU:HG	2.19	0.41
1:E:157:THR:C	1:E:159:ASP:H	2.28	0.41
1:F:44:HIS:O	1:F:47:ASP:HB2	2.19	0.41
1:B:71:THR:O	1:B:72:TRP:CB	2.67	0.41
1:C:65:TRP:CG	1:C:79:LEU:HD11	2.55	0.41
1:G:49:TRP:O	1:G:53:LEU:HG	2.20	0.41
1:H:128:ASN:HB3	1:H:131:LYS:HG2	2.03	0.41
1:B:41:LEU:HD12	1:B:41:LEU:HA	1.78	0.41
1:C:107:LYS:HE3	1:C:107:LYS:HB2	1.86	0.41
1:G:45:LYS:HA	1:G:107:LYS:HA	2.03	0.41
1:H:9:LYS:HE3	1:H:148:ASP:CG	2.46	0.41
1:H:13:LEU:HD21	1:H:141:MET:SD	2.61	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:153:ILE:HD13	1:B:153:ILE:HA	1.91	0.41
1:D:41:LEU:HD12	1:D:41:LEU:HA	1.77	0.41
1:E:5:ILE:HB	1:E:153:ILE:HB	2.01	0.41
1:F:76:VAL:CG2	1:F:81:TYR:HB2	2.44	0.41
1:F:91:LEU:HD23	1:F:91:LEU:HA	1.73	0.41
1:G:55:ILE:O	1:G:58:LYS:N	2.54	0.41
1:A:3:VAL:HG12	1:A:155:ILE:HB	2.03	0.41
1:E:63:ARG:HD3	1:F:100:LYS:CB	2.50	0.41
1:B:36:GLY:HA3	1:B:82:GLU:O	2.21	0.41
1:B:66:ASN:OD1	1:B:71:THR:HA	2.21	0.41
1:C:44:HIS:O	1:C:47:ASP:HB2	2.20	0.41
1:F:61:ASN:HA	1:F:75:ASN:HD22	1.84	0.41
1:G:118:LYS:O	1:G:121:ASP:HB2	2.21	0.41
1:A:9:LYS:HG2	1:A:148:ASP:HB2	2.02	0.41
1:A:48:THR:HG23	1:A:51:ASP:OD2	2.21	0.41
1:A:52:ALA:HB1	1:A:119:ILE:HG12	2.03	0.41
1:C:43:ILE:CG2	1:C:115:PHE:HD1	2.33	0.41
1:D:3:VAL:HG12	1:D:155:ILE:HB	2.03	0.41
1:E:9:LYS:HE3	1:E:148:ASP:CG	2.45	0.41
1:F:5:ILE:HB	1:F:153:ILE:HB	2.03	0.41
1:H:88:GLN:CD	1:H:90:THR:HG1	2.27	0.41
1:A:102:ASP:OD1	1:A:102:ASP:C	2.63	0.41
1:C:13:LEU:HD11	1:C:141:MET:SD	2.61	0.41
1:H:116:TYR:CE1	1:H:120:ILE:HD11	2.56	0.41
1:C:43:ILE:CG2	1:C:115:PHE:CD1	3.05	0.40
1:C:60:LEU:O	1:C:77:LYS:HE2	2.21	0.40
1:C:79:LEU:HD13	1:C:81:TYR:CE2	2.56	0.40
1:F:157:THR:C	1:F:159:ASP:H	2.29	0.40
1:H:56:ILE:HG23	1:H:123:PHE:CD1	2.56	0.40
1:H:71:THR:O	1:H:72:TRP:C	2.64	0.40
1:H:127:TYR:O	1:H:128:ASN:HB2	2.20	0.40
1:A:43:ILE:CG2	1:A:115:PHE:HD1	2.34	0.40
1:E:107:LYS:HG2	1:E:108:GLU:N	2.36	0.40
1:H:13:LEU:HD21	1:H:141:MET:HG2	2.03	0.40
1:G:60:LEU:HD11	1:G:123:PHE:HD1	1.86	0.40
1:C:107:LYS:HG2	1:C:112:ALA:N	2.36	0.40
1:C:95:THR:HG21	1:D:62:LEU:CD2	2.51	0.40
1:D:153:ILE:HD13	1:D:153:ILE:HA	1.94	0.40
1:G:108:GLU:O	1:G:109:PRO:C	2.65	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	149/164 (91%)	128 (86%)	18 (12%)	3 (2%)	6	27
1	B	142/164 (87%)	126 (89%)	14 (10%)	2 (1%)	9	33
1	C	144/164 (88%)	128 (89%)	15 (10%)	1 (1%)	18	49
1	D	144/164 (88%)	124 (86%)	16 (11%)	4 (3%)	4	21
1	E	137/164 (84%)	120 (88%)	14 (10%)	3 (2%)	5	26
1	F	138/164 (84%)	114 (83%)	23 (17%)	1 (1%)	18	49
1	G	144/164 (88%)	118 (82%)	18 (12%)	8 (6%)	1	9
1	H	144/164 (88%)	120 (83%)	20 (14%)	4 (3%)	4	21
All	All	1142/1312 (87%)	978 (86%)	138 (12%)	26 (2%)	5	25

All (26) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	97	LYS
1	B	68	ASP
1	D	69	ASN
1	G	97	LYS
1	G	101	PRO
1	G	102	ASP
1	B	69	ASN
1	E	157	THR
1	G	55	ILE
1	G	158	HIS
1	E	158	HIS
1	G	68	ASP
1	H	158	HIS
1	D	67	ASN
1	D	97	LYS
1	D	101	PRO
1	G	56	ILE

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Mol	Chain	Res	Type
1	G	59	CYS
1	H	157	THR
1	A	109	PRO
1	H	124	LYS
1	A	102	ASP
1	C	109	PRO
1	F	30	ILE
1	E	101	PRO
1	H	149	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	121/151 (80%)	111 (92%)	10 (8%)	10	34
1	B	122/151 (81%)	112 (92%)	10 (8%)	10	35
1	C	125/151 (83%)	112 (90%)	13 (10%)	7	25
1	D	123/151 (82%)	117 (95%)	6 (5%)	22	51
1	E	117/151 (78%)	113 (97%)	4 (3%)	32	59
1	F	115/151 (76%)	110 (96%)	5 (4%)	26	54
1	G	111/151 (74%)	97 (87%)	14 (13%)	4	18
1	H	117/151 (78%)	108 (92%)	9 (8%)	12	37
All	All	951/1208 (79%)	880 (92%)	71 (8%)	12	38

All (71) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	VAL
1	A	17	LYS
1	A	34	LYS
1	A	38	ILE
1	A	43	ILE
1	A	48	THR

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Mol	Chain	Res	Type
1	A	59	CYS
1	A	77	LYS
1	A	80	ASN
1	A	129	ASP
1	B	12	ILE
1	B	14	SER
1	B	34	LYS
1	B	48	THR
1	B	67	ASN
1	B	69	ASN
1	B	80	ASN
1	B	95	THR
1	B	99	ASN
1	B	129	ASP
1	C	12	ILE
1	C	14	SER
1	C	15	VAL
1	C	27	LEU
1	C	29	ILE
1	C	43	ILE
1	C	48	THR
1	C	74	LYS
1	C	80	ASN
1	C	95	THR
1	C	105	LEU
1	C	129	ASP
1	C	159	ASP
1	D	15	VAL
1	D	38	ILE
1	D	48	THR
1	D	80	ASN
1	D	95	THR
1	D	129	ASP
1	E	27	LEU
1	E	48	THR
1	E	105	LEU
1	E	129	ASP
1	F	48	THR
1	F	80	ASN
1	F	95	THR
1	F	104	HIS
1	F	107	LYS

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Mol	Chain	Res	Type
1	G	2	ARG
1	G	15	VAL
1	G	30	ILE
1	G	48	THR
1	G	80	ASN
1	G	91	LEU
1	G	95	THR
1	G	113	LEU
1	G	119	ILE
1	G	134	ILE
1	G	142	ASN
1	G	143	ILE
1	G	148	ASP
1	G	155	ILE
1	H	8	VAL
1	H	14	SER
1	H	27	LEU
1	H	31	SER
1	H	41	LEU
1	H	48	THR
1	H	69	ASN
1	H	145	VAL
1	H	153	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	67	ASN
1	A	80	ASN
1	B	147	ASN
1	D	99	ASN
1	E	75	ASN
1	F	80	ASN
1	F	94	ASN
1	G	80	ASN
1	H	67	ASN
1	H	69	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 26 ligands modelled in this entry, 26 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th 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 ²)	Q<0.9
1	A	155/164 (94%)	-0.30	0	100	100	27, 42, 57, 68	0
1	B	148/164 (90%)	-0.43	0	100	100	31, 42, 66, 83	1 (0%)
1	C	148/164 (90%)	-0.46	0	100	100	29, 43, 61, 73	0
1	D	148/164 (90%)	-0.50	0	100	100	32, 41, 60, 66	2 (1%)
1	E	145/164 (88%)	-0.42	0	100	100	31, 44, 64, 73	0
1	F	144/164 (87%)	-0.23	0	100	100	33, 47, 65, 83	0
1	G	148/164 (90%)	-0.35	1 (0%)	84	70	34, 63, 87, 100	0
1	H	148/164 (90%)	-0.47	0	100	100	37, 55, 74, 88	0
All	All	1184/1312 (90%)	-0.39	1 (0%)	92	90	27, 46, 72, 100	3 (0%)

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	102	ASP	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 oligosaccharides 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, 95th 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 ²)	Q<0.9
2	IOD	B	167	1/1	0.54	0.20	85,85,85,85	1
2	IOD	B	169	1/1	0.85	0.09	76,76,76,76	1
2	IOD	C	168	1/1	0.90	0.07	55,55,55,55	1
2	IOD	C	167	1/1	0.92	0.06	74,74,74,74	1
2	IOD	G	167	1/1	0.94	0.07	97,97,97,97	1
2	IOD	F	168	1/1	0.95	0.06	44,44,44,44	1
2	IOD	D	168	1/1	0.95	0.06	54,54,54,54	1
2	IOD	B	168	1/1	0.96	0.05	74,74,74,74	1
2	IOD	F	167	1/1	0.96	0.06	77,77,77,77	0
2	IOD	D	169	1/1	0.97	0.05	56,56,56,56	1
2	IOD	D	170	1/1	0.97	0.06	63,63,63,63	1
2	IOD	D	166	1/1	0.97	0.04	72,72,72,72	0
2	IOD	D	167	1/1	0.97	0.04	46,46,46,46	1
2	IOD	C	166	1/1	0.97	0.04	62,62,62,62	1
2	IOD	G	166	1/1	0.98	0.03	72,72,72,72	0
2	IOD	B	165	1/1	0.98	0.03	73,73,73,73	0
2	IOD	H	166	1/1	0.98	0.03	61,61,61,61	1
2	IOD	F	165	1/1	0.99	0.03	90,90,90,90	0
2	IOD	F	166	1/1	0.99	0.09	79,79,79,79	1
2	IOD	A	165	1/1	0.99	0.04	48,48,48,48	0
2	IOD	C	165	1/1	0.99	0.03	65,65,65,65	0
2	IOD	G	165	1/1	0.99	0.02	75,75,75,75	0
2	IOD	D	165	1/1	0.99	0.02	54,54,54,54	0
2	IOD	B	166	1/1	0.99	0.03	80,80,80,80	0
2	IOD	H	165	1/1	0.99	0.02	63,63,63,63	1
2	IOD	E	165	1/1	0.99	0.02	54,54,54,54	0

6.5 Other polymers [i](#)

There are no such residues in this entry.