



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 06:33 AM UTC

PDB ID : 1JOG / pdb_00001jog
Title : Structure of HI0074 from Heamophilus Influenzae reveals the fold of a substrate binding domain of a nucleotidyltransferase
Authors : Lehmann, C.; Lim, K.; Herzberg, O.; Structure 2 Function Project (S2F)
Deposited on : 2001-07-29
Resolution : 2.40 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

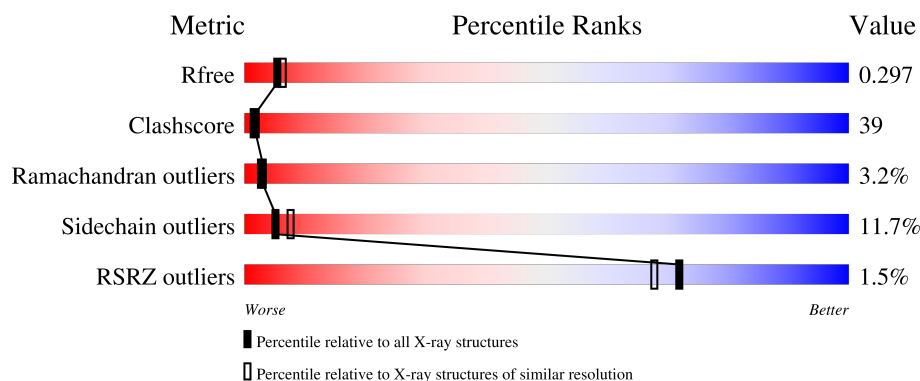
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	146	 3% 48% 32% 12% • 8%
1	B	146	 43% 38% 9% • 8%
1	C	146	 2% 30% 45% 16% • 8%
1	D	146	 41% 38% 13% • 7%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4575 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 HYPOTHETICAL PROTEIN HI0074.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	135	Total	C	N	O	Se	0	0	0
			1106	706	183	211	6			
1	B	135	Total	C	N	O	Se	0	0	0
			1106	706	183	211	6			
1	C	135	Total	C	N	O	Se	0	0	0
			1106	706	183	211	6			
1	D	136	Total	C	N	O	Se	0	0	0
			1117	712	187	212	6			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	modified residue	UNP P43934
A	2	MSE	MET	modified residue	UNP P43934
A	33	MSE	MET	modified residue	UNP P43934
A	60	MSE	MET	modified residue	UNP P43934
A	61	MSE	MET	modified residue	UNP P43934
A	96	MSE	MET	modified residue	UNP P43934
A	105	MSE	MET	modified residue	UNP P43934
A	119	MSE	MET	modified residue	UNP P43934
B	1	MSE	MET	modified residue	UNP P43934
B	2	MSE	MET	modified residue	UNP P43934
B	33	MSE	MET	modified residue	UNP P43934
B	60	MSE	MET	modified residue	UNP P43934
B	61	MSE	MET	modified residue	UNP P43934
B	96	MSE	MET	modified residue	UNP P43934
B	105	MSE	MET	modified residue	UNP P43934
B	119	MSE	MET	modified residue	UNP P43934
C	1	MSE	MET	modified residue	UNP P43934
C	2	MSE	MET	modified residue	UNP P43934
C	33	MSE	MET	modified residue	UNP P43934
C	60	MSE	MET	modified residue	UNP P43934
C	61	MSE	MET	modified residue	UNP P43934

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Chain	Residue	Modelled	Actual	Comment	Reference
C	96	MSE	MET	modified residue	UNP P43934
C	105	MSE	MET	modified residue	UNP P43934
C	119	MSE	MET	modified residue	UNP P43934
D	1	MSE	MET	modified residue	UNP P43934
D	2	MSE	MET	modified residue	UNP P43934
D	33	MSE	MET	modified residue	UNP P43934
D	60	MSE	MET	modified residue	UNP P43934
D	61	MSE	MET	modified residue	UNP P43934
D	96	MSE	MET	modified residue	UNP P43934
D	105	MSE	MET	modified residue	UNP P43934
D	119	MSE	MET	modified residue	UNP P43934

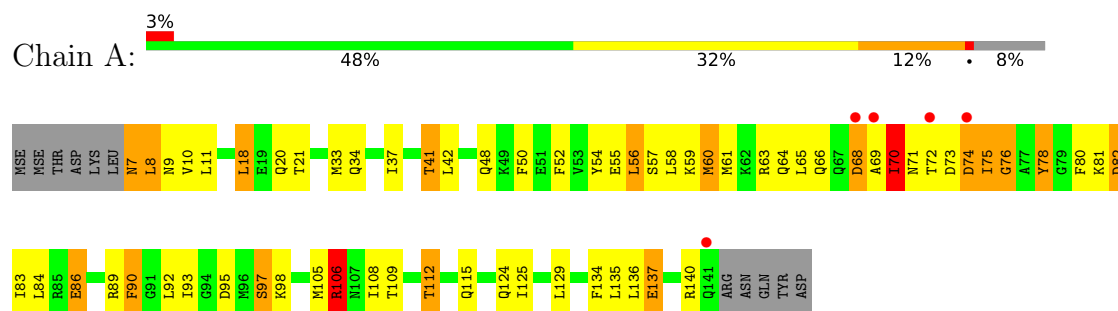
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	45	Total O 45 45	0	0
2	B	40	Total O 40 40	0	0
2	C	12	Total O 12 12	0	0
2	D	43	Total O 43 43	0	0

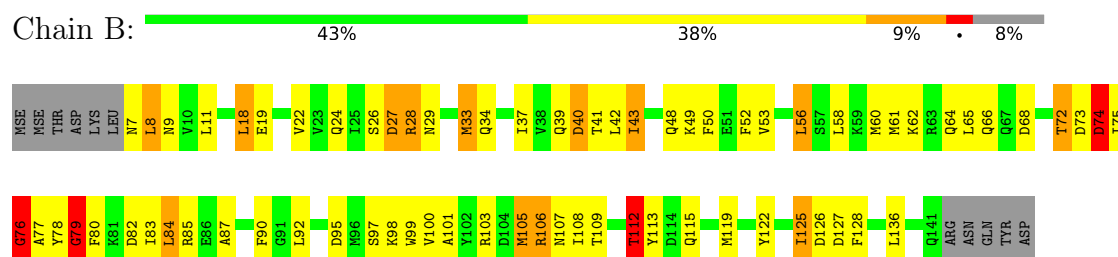
3 Residue-property plots

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

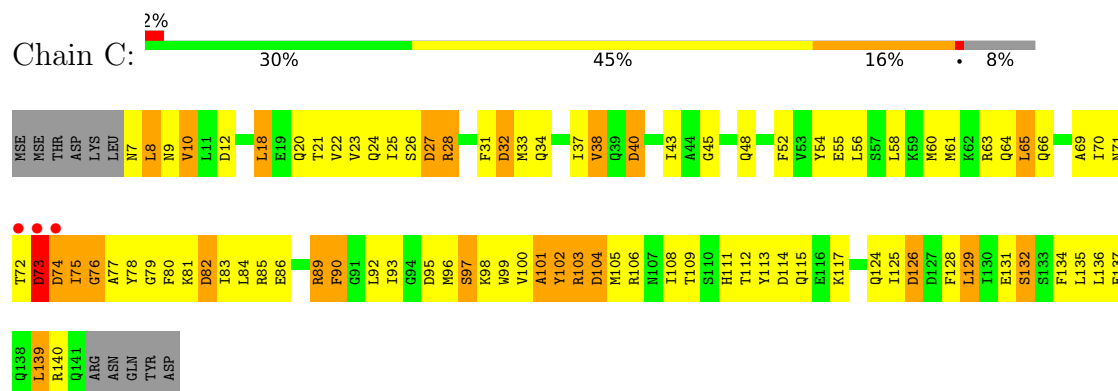
• Molecule 1: HYPOTHETICAL PROTEIN HI0074



• Molecule 1: HYPOTHETICAL PROTEIN HI0074

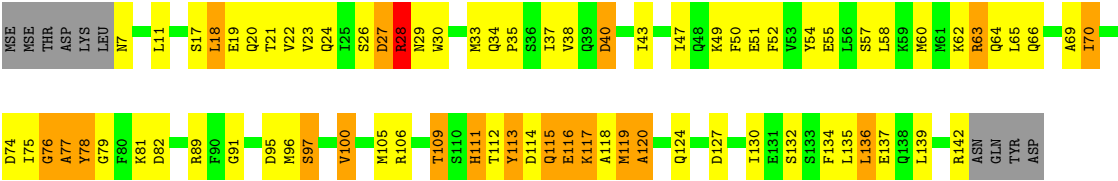


• Molecule 1: HYPOTHETICAL PROTEIN HI0074



• Molecule 1: HYPOTHETICAL PROTEIN HI0074





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	38.30Å 87.10Å 165.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.40 20.00 – 2.40	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-2.40) 96.2 (20.00-2.40)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.69 (at 2.39Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.211 , 0.282 0.232 , 0.297	Depositor DCC
R_{free} test set	1131 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	42.8	Xtriage
Anisotropy	0.672	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 58.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4575	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.83% 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

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	1.29	2/1119 (0.2%)	1.52	19/1499 (1.3%)
1	B	1.17	1/1119 (0.1%)	1.39	13/1499 (0.9%)
1	C	1.03	0/1119	1.40	15/1499 (1.0%)
1	D	1.07	0/1130	1.41	11/1513 (0.7%)
All	All	1.15	3/4487 (0.1%)	1.43	58/6010 (1.0%)

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	B	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	60	MSE	SE-CE	-6.88	1.74	1.95
1	B	43	ILE	CA-CB	6.61	1.61	1.54
1	A	70	ILE	CA-CB	5.87	1.60	1.54

All (58) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	8	LEU	N-CA-C	-10.58	99.77	112.89
1	C	73	ASP	N-CA-C	10.24	123.72	111.33
1	A	34	GLN	CA-C-N	-8.93	111.19	120.03
1	A	34	GLN	C-N-CA	-8.93	111.19	120.03
1	C	23	VAL	CB-CA-C	-8.69	100.75	111.88
1	C	102	TYR	N-CA-C	-8.57	101.88	111.82
1	B	100	VAL	N-CA-C	-8.56	101.89	110.62
1	B	34	GLN	N-CA-C	8.05	118.85	109.60
1	C	89	ARG	N-CA-C	-8.04	102.60	111.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	125	ILE	N-CA-C	7.91	118.71	110.72
1	C	126	ASP	N-CA-C	-7.75	101.95	111.33
1	B	77	ALA	N-CA-C	-7.67	101.77	113.89
1	D	116	GLU	N-CA-C	-7.61	102.98	112.72
1	D	34	GLN	CA-C-N	7.57	127.33	119.76
1	D	34	GLN	C-N-CA	7.57	127.33	119.76
1	C	92	LEU	N-CA-C	-7.54	103.02	112.90
1	A	71	ASN	N-CA-C	-7.40	101.15	111.81
1	A	8	LEU	N-CA-C	-7.24	103.39	111.71
1	A	69	ALA	N-CA-C	-7.21	105.01	113.88
1	A	78	TYR	N-CA-C	7.19	120.39	108.96
1	B	83	ILE	N-CA-C	-6.99	103.96	110.53
1	B	99	TRP	N-CA-C	6.92	119.85	111.82
1	B	125	ILE	N-CA-C	6.92	117.71	110.72
1	B	127	ASP	N-CA-C	-6.85	103.75	111.14
1	D	117	LYS	N-CA-C	-6.69	103.68	110.97
1	C	129	LEU	N-CA-C	6.57	119.28	111.33
1	A	41	THR	N-CA-C	-6.53	104.24	111.36
1	D	77	ALA	CA-C-N	-6.47	111.94	122.81
1	D	77	ALA	C-N-CA	-6.47	111.94	122.81
1	C	104	ASP	N-CA-C	-6.39	104.56	113.37
1	B	115	GLN	N-CA-C	6.36	118.21	111.28
1	A	61	MSE	N-CA-C	-6.26	104.46	111.28
1	A	81	LYS	N-CA-C	6.23	119.98	112.38
1	D	76	GLY	N-CA-C	-6.11	106.01	115.73
1	B	79	GLY	N-CA-C	6.07	127.57	113.18
1	C	81	LYS	N-CA-C	-5.70	104.27	111.11
1	A	115	GLN	N-CA-C	5.67	119.01	111.75
1	A	82	ASP	N-CA-C	-5.65	104.51	111.40
1	D	74	ASP	N-CA-C	5.57	119.58	112.34
1	C	32	ASP	N-CA-C	5.55	117.33	111.28
1	A	76	GLY	N-CA-C	-5.50	100.15	113.18
1	A	97	SER	N-CA-C	-5.48	105.41	111.71
1	A	70	ILE	N-CA-C	5.45	116.73	109.37
1	D	78	TYR	N-CA-C	5.45	118.46	109.85
1	C	76	GLY	N-CA-C	-5.39	100.40	113.18
1	C	65	LEU	N-CA-C	-5.37	106.76	113.15
1	A	61	MSE	CG-SE-CE	-5.34	87.17	98.92
1	A	106	ARG	CA-CB-CG	-5.33	103.44	114.10
1	D	66	GLN	N-CA-C	-5.27	105.45	111.14
1	D	109	THR	N-CA-C	-5.23	106.49	112.92
1	B	76	GLY	N-CA-C	-5.18	100.90	113.18

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	101	ALA	N-CA-C	-5.18	106.11	112.90
1	C	38	VAL	CB-CA-C	-5.18	105.15	112.14
1	B	84	LEU	N-CA-C	5.17	116.61	110.97
1	B	64	GLN	N-CA-C	-5.17	105.54	111.07
1	A	68	ASP	N-CA-C	5.08	121.61	110.80
1	A	93	ILE	N-CA-C	-5.06	98.82	109.34
1	B	113	TYR	N-CA-C	-5.04	106.64	112.89

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	122	TYR	Sidechain

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	1106	0	1085	80	0
1	B	1106	0	1085	67	0
1	C	1106	0	1085	143	0
1	D	1117	0	1098	96	0
2	A	45	0	0	0	0
2	B	40	0	0	0	0
2	C	12	0	0	2	0
2	D	43	0	0	4	0
All	All	4575	0	4353	346	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 39.

All (346) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:75:ILE:HG23	1:B:78:TYR:HB3	1.18	1.11

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:7:ASN:ND2	1:D:63:ARG:HH22	1.52	1.07
1:A:106:ARG:HG2	1:A:106:ARG:HH11	1.11	1.05
1:C:20:GLN:HB3	1:D:37:ILE:HD12	1.35	1.04
1:C:75:ILE:HG23	1:C:78:TYR:HD1	1.19	1.04
1:C:65:LEU:HD22	1:C:86:GLU:HG3	1.39	1.02
1:B:75:ILE:HD13	1:B:82:ASP:HB3	1.38	1.01
1:D:78:TYR:HD2	1:D:79:GLY:N	1.61	0.98
1:D:28:ARG:HH11	1:D:28:ARG:HG2	1.29	0.95
1:C:22:VAL:HG21	1:C:129:LEU:HD11	1.45	0.95
1:B:33:MSE:HA	1:C:33:MSE:HE1	1.50	0.94
1:C:106:ARG:HD2	1:C:106:ARG:O	1.71	0.90
1:B:66:GLN:HG2	1:B:72:THR:HG23	1.53	0.89
1:D:64:GLN:HE22	1:D:142:ARG:HG2	1.38	0.88
1:D:7:ASN:HD22	1:D:63:ARG:HH22	1.21	0.88
1:D:23:VAL:O	1:D:26:SER:HB2	1.74	0.88
1:D:70:ILE:HG13	2:D:1139:HOH:O	1.72	0.87
1:D:105:MSE:O	1:D:109:THR:HG23	1.74	0.87
1:A:20:GLN:HB3	1:B:37:ILE:HD12	1.58	0.86
1:B:106:ARG:HG2	1:B:106:ARG:HH11	1.40	0.86
1:D:49:LYS:HA	1:D:52:PHE:HD1	1.41	0.85
1:C:20:GLN:HB3	1:D:37:ILE:CD1	2.06	0.85
1:C:63:ARG:O	1:C:66:GLN:HB3	1.77	0.85
1:D:33:MSE:HG3	2:D:1046:HOH:O	1.76	0.85
1:C:95:ASP:HB3	1:C:98:LYS:HG2	1.58	0.84
1:D:7:ASN:HD22	1:D:63:ARG:NH2	1.76	0.84
1:D:105:MSE:CE	1:D:124:GLN:HB2	2.07	0.83
1:A:106:ARG:HH11	1:A:106:ARG:CG	1.92	0.82
1:C:114:ASP:HB3	1:C:117:LYS:HB2	1.61	0.82
1:D:7:ASN:ND2	1:D:63:ARG:NH2	2.27	0.82
1:C:75:ILE:HG23	1:C:78:TYR:CD1	2.10	0.82
1:D:78:TYR:CD2	1:D:79:GLY:N	2.41	0.81
1:C:25:ILE:HG23	1:C:31:PHE:CE1	2.16	0.80
1:C:37:ILE:HD12	1:D:20:GLN:HB3	1.63	0.80
1:C:70:ILE:HA	1:C:73:ASP:OD2	1.82	0.79
1:D:115:GLN:H	1:D:115:GLN:NE2	1.81	0.79
1:A:106:ARG:HG2	1:A:106:ARG:NH1	1.90	0.78
1:B:62:LYS:HE3	1:B:76:GLY:HA2	1.65	0.77
1:C:10:VAL:HG12	1:C:56:LEU:HD21	1.67	0.77
1:C:7:ASN:HD21	1:C:60:MSE:SE	2.17	0.76
1:D:51:GLU:OE1	1:D:109:THR:OG1	2.02	0.76
1:C:95:ASP:HB3	1:C:98:LYS:CG	2.17	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:50:PHE:CZ	1:B:105:MSE:HE2	2.22	0.75
1:A:64:GLN:O	1:A:68:ASP:HB2	1.87	0.74
1:C:75:ILE:CG2	1:C:78:TYR:HD1	1.99	0.74
1:B:75:ILE:HG23	1:B:78:TYR:CB	2.08	0.74
1:C:7:ASN:HD21	1:C:60:MSE:CE	2.01	0.74
1:C:95:ASP:O	1:C:98:LYS:HB2	1.88	0.73
1:A:70:ILE:HD12	1:A:75:ILE:HD11	1.70	0.73
1:C:65:LEU:HD21	1:C:86:GLU:O	1.88	0.73
1:D:28:ARG:HG2	1:D:28:ARG:NH1	1.92	0.73
1:C:111:HIS:HA	1:C:113:TYR:CZ	2.25	0.72
1:B:60:MSE:HE2	1:B:60:MSE:HA	1.72	0.71
1:A:7:ASN:ND2	1:A:7:ASN:O	2.24	0.71
1:C:7:ASN:ND2	1:C:60:MSE:SE	2.73	0.71
1:A:64:GLN:HG2	1:A:92:LEU:HD11	1.73	0.71
1:D:114:ASP:HB3	1:D:117:LYS:HB2	1.73	0.70
1:C:104:ASP:O	1:C:108:ILE:HG12	1.91	0.70
1:C:112:THR:HB	1:D:52:PHE:HZ	1.57	0.70
1:C:20:GLN:C	1:D:37:ILE:HD11	2.17	0.69
1:C:78:TYR:CD2	1:C:82:ASP:HB3	2.26	0.69
1:C:37:ILE:CD1	1:D:20:GLN:HB3	2.22	0.69
1:C:75:ILE:O	1:C:75:ILE:HG22	1.91	0.69
1:A:20:GLN:HB3	1:B:37:ILE:CD1	2.24	0.68
1:A:70:ILE:HD12	1:A:75:ILE:CD1	2.23	0.68
1:C:65:LEU:HD22	1:C:86:GLU:CG	2.19	0.68
1:D:115:GLN:H	1:D:115:GLN:HE21	1.43	0.66
1:B:78:TYR:CD2	1:B:79:GLY:N	2.62	0.66
1:C:102:TYR:HA	1:C:105:MSE:HE2	1.77	0.66
1:A:7:ASN:O	1:A:10:VAL:HG22	1.95	0.66
1:C:26:SER:O	1:C:28:ARG:N	2.29	0.66
1:B:103:ARG:HG2	1:B:107:ASN:ND2	2.11	0.66
1:D:115:GLN:NE2	1:D:115:GLN:N	2.44	0.65
1:A:136:LEU:HD22	1:A:140:ARG:HH21	1.61	0.65
1:C:90:PHE:HD1	1:C:90:PHE:N	1.94	0.65
1:B:26:SER:O	1:B:28:ARG:N	2.28	0.65
1:C:95:ASP:CB	1:C:98:LYS:HG2	2.26	0.65
1:A:65:LEU:HD23	1:A:90:PHE:CE2	2.31	0.65
1:A:75:ILE:O	1:A:75:ILE:HG22	1.94	0.65
1:C:112:THR:HB	1:D:52:PHE:CZ	2.31	0.65
1:A:70:ILE:HG23	1:A:73:ASP:H	1.62	0.65
1:A:95:ASP:OD2	1:A:97:SER:OG	2.11	0.65
1:C:24:GLN:HG2	2:C:1034:HOH:O	1.97	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:127:ASP:O	1:D:130:ILE:HG22	1.97	0.64
1:B:19:GLU:O	1:B:22:VAL:HG22	1.97	0.64
1:A:75:ILE:HG23	1:A:78:TYR:CD1	2.32	0.64
1:D:26:SER:O	1:D:28:ARG:N	2.28	0.64
1:C:85:ARG:HB3	1:C:89:ARG:HH12	1.64	0.63
1:A:70:ILE:HD12	1:A:75:ILE:CG1	2.27	0.63
1:C:10:VAL:HG12	1:C:56:LEU:CD2	2.29	0.63
1:B:66:GLN:HE21	1:B:72:THR:CG2	2.12	0.63
1:C:90:PHE:N	1:C:90:PHE:CD1	2.66	0.62
1:D:62:LYS:CE	1:D:76:GLY:HA2	2.29	0.62
1:A:106:ARG:CG	1:A:106:ARG:NH1	2.55	0.62
1:C:75:ILE:HG22	1:C:78:TYR:HB2	1.81	0.62
1:C:103:ARG:C	1:C:105:MSE:H	2.07	0.62
1:C:78:TYR:HD2	1:C:82:ASP:HB3	1.62	0.62
1:B:66:GLN:CG	1:B:72:THR:HG23	2.27	0.62
1:B:28:ARG:HG2	1:B:28:ARG:HH11	1.64	0.62
1:C:79:GLY:O	1:C:82:ASP:HB2	1.99	0.62
1:A:73:ASP:O	1:A:74:ASP:C	2.43	0.61
1:A:78:TYR:CD2	1:A:82:ASP:HB3	2.35	0.61
1:D:105:MSE:HE3	1:D:124:GLN:HB2	1.82	0.61
1:C:135:LEU:CD1	1:C:139:LEU:HD12	2.31	0.61
1:D:114:ASP:OD2	1:D:116:GLU:HB2	1.99	0.61
1:C:74:ASP:C	1:C:76:GLY:H	2.08	0.61
1:C:7:ASN:HD21	1:C:60:MSE:HE1	1.65	0.60
1:C:9:ASN:O	1:C:12:ASP:HB2	2.00	0.60
1:D:117:LYS:O	1:D:120:ALA:N	2.32	0.60
1:C:66:GLN:HG3	1:C:73:ASP:CG	2.26	0.60
1:D:19:GLU:O	1:D:22:VAL:HG22	2.00	0.60
1:A:112:THR:HG22	1:B:52:PHE:CZ	2.37	0.60
1:A:70:ILE:HD12	1:A:75:ILE:HG13	1.83	0.60
1:A:137:GLU:OE2	1:A:140:ARG:CZ	2.50	0.60
1:C:37:ILE:CG2	1:D:24:GLN:HG3	2.31	0.60
1:D:113:TYR:HD1	1:D:114:ASP:N	2.00	0.60
1:A:109:THR:O	1:A:112:THR:HB	2.02	0.59
1:D:117:LYS:O	1:D:118:ALA:C	2.42	0.59
1:A:20:GLN:C	1:B:37:ILE:HD11	2.27	0.59
1:D:78:TYR:HD2	1:D:79:GLY:H	0.79	0.59
1:C:102:TYR:CD1	1:C:128:PHE:HD1	2.21	0.59
1:C:73:ASP:O	1:C:75:ILE:N	2.32	0.59
1:A:112:THR:HG22	1:B:52:PHE:HZ	1.67	0.59
1:C:18:LEU:HD22	1:C:18:LEU:O	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:103:ARG:HG2	1:B:107:ASN:HD21	1.67	0.59
1:D:81:LYS:HZ2	1:D:100:VAL:HG22	1.68	0.59
1:D:47:ILE:HD13	1:D:112:THR:HG21	1.84	0.58
1:C:100:VAL:O	1:C:103:ARG:HB3	2.03	0.58
1:D:40:ASP:OD2	1:D:40:ASP:C	2.46	0.58
1:D:11:LEU:CD2	1:D:136:LEU:HG	2.33	0.58
1:A:18:LEU:HD13	1:A:129:LEU:HD21	1.85	0.58
1:A:137:GLU:OE2	1:A:140:ARG:NH1	2.37	0.58
1:C:102:TYR:CE1	1:C:128:PHE:HD1	2.22	0.58
1:C:126:ASP:O	1:C:129:LEU:HB2	2.04	0.58
1:B:106:ARG:HG2	1:B:106:ARG:NH1	2.06	0.58
1:C:75:ILE:CG2	1:C:78:TYR:CD1	2.81	0.58
1:C:52:PHE:CZ	1:D:112:THR:HB	2.39	0.58
1:C:135:LEU:HD11	1:C:139:LEU:HD12	1.87	0.57
1:D:105:MSE:SE	1:D:124:GLN:OE1	2.73	0.57
1:A:78:TYR:HD2	1:A:82:ASP:HB3	1.69	0.57
1:C:61:MSE:SE	1:C:93:ILE:HD11	2.54	0.57
1:C:22:VAL:HG21	1:C:129:LEU:CD1	2.27	0.57
1:C:69:ALA:C	1:C:71:ASN:H	2.11	0.57
1:A:70:ILE:HD13	1:A:73:ASP:HB2	1.86	0.57
1:B:109:THR:O	1:B:112:THR:HB	2.05	0.57
1:C:61:MSE:CE	1:C:93:ILE:HD11	2.35	0.57
1:B:49:LYS:O	1:B:53:VAL:HG23	2.04	0.57
1:D:78:TYR:CD1	1:D:82:ASP:OD2	2.58	0.56
1:D:105:MSE:HE1	1:D:124:GLN:HB2	1.84	0.56
1:C:75:ILE:HD12	1:C:78:TYR:CD1	2.40	0.56
1:A:52:PHE:CZ	1:B:112:THR:HG22	2.41	0.56
1:C:48:GLN:HG3	1:C:52:PHE:CE1	2.41	0.56
1:D:113:TYR:CD1	1:D:114:ASP:N	2.73	0.56
1:C:105:MSE:O	1:C:109:THR:HG23	2.06	0.56
1:C:128:PHE:O	1:C:132:SER:OG	2.25	0.55
1:A:37:ILE:CG2	1:B:24:GLN:HG3	2.37	0.55
1:C:20:GLN:CB	1:D:37:ILE:CD1	2.82	0.55
1:D:69:ALA:HB2	1:D:75:ILE:HD12	1.87	0.55
1:C:21:THR:N	1:D:37:ILE:HD11	2.21	0.55
1:A:50:PHE:CZ	1:A:105:MSE:HE2	2.42	0.55
1:B:40:ASP:C	1:B:40:ASP:OD2	2.49	0.55
1:B:66:GLN:HE21	1:B:72:THR:HG22	1.72	0.55
1:A:60:MSE:HE1	1:A:63:ARG:CZ	2.37	0.54
1:B:98:LYS:O	1:B:101:ALA:HB3	2.06	0.54
1:D:21:THR:OG1	1:D:49:LYS:NZ	2.35	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:66:GLN:NE2	1:B:72:THR:CG2	2.69	0.54
1:B:82:ASP:OD2	1:B:85:ARG:NH1	2.40	0.54
1:D:35:PRO:HG2	1:D:38:VAL:HG23	1.90	0.54
1:C:75:ILE:HD12	1:C:78:TYR:HD1	1.73	0.54
1:C:101:ALA:O	1:C:105:MSE:HB2	2.06	0.54
1:A:75:ILE:HG23	1:A:78:TYR:CG	2.42	0.54
1:D:49:LYS:HA	1:D:52:PHE:CD1	2.32	0.53
1:C:137:GLU:OE2	2:C:1027:HOH:O	2.19	0.53
1:A:10:VAL:HG23	1:A:11:LEU:N	2.23	0.53
1:D:20:GLN:OE1	1:D:20:GLN:HA	2.07	0.53
1:A:75:ILE:HG23	1:A:78:TYR:HB2	1.90	0.53
1:C:102:TYR:CD1	1:C:128:PHE:CD1	2.97	0.53
1:C:75:ILE:CG2	1:C:78:TYR:HB2	2.38	0.53
1:C:99:TRP:HA	1:C:102:TYR:CD1	2.43	0.53
1:C:32:ASP:C	1:C:34:GLN:N	2.67	0.53
1:C:111:HIS:CD2	1:C:117:LYS:HZ3	2.25	0.53
1:B:28:ARG:HG2	1:B:28:ARG:NH1	2.24	0.52
1:C:37:ILE:HG21	1:D:24:GLN:CG	2.39	0.52
1:C:52:PHE:HZ	1:D:112:THR:HB	1.74	0.52
1:C:58:LEU:HD11	1:C:80:PHE:HE1	1.74	0.52
1:C:84:LEU:O	1:C:85:ARG:C	2.50	0.52
1:A:7:ASN:ND2	1:A:7:ASN:C	2.67	0.52
1:A:48:GLN:OE1	1:B:48:GLN:OE1	2.27	0.52
1:C:20:GLN:C	1:D:37:ILE:CD1	2.82	0.52
1:C:79:GLY:O	1:C:83:ILE:HG12	2.10	0.52
1:C:37:ILE:HG23	1:C:38:VAL:N	2.25	0.52
1:D:89:ARG:C	1:D:91:GLY:H	2.18	0.52
1:A:41:THR:HG22	1:B:42:LEU:HA	1.90	0.52
1:A:76:GLY:C	1:A:78:TYR:N	2.65	0.52
1:B:75:ILE:CG2	1:B:78:TYR:HB3	2.13	0.52
1:C:37:ILE:CG2	1:C:38:VAL:N	2.73	0.52
1:A:42:LEU:HD23	1:B:41:THR:CG2	2.40	0.51
1:C:74:ASP:O	1:C:76:GLY:N	2.43	0.51
1:D:134:PHE:O	1:D:137:GLU:HB2	2.10	0.51
1:C:78:TYR:HD2	1:C:82:ASP:CB	2.23	0.51
1:D:35:PRO:HG2	1:D:38:VAL:CG2	2.40	0.51
1:D:95:ASP:OD2	1:D:97:SER:HB3	2.10	0.51
1:C:80:PHE:CE2	1:C:103:ARG:HB2	2.45	0.51
1:A:74:ASP:O	1:A:76:GLY:N	2.42	0.51
1:B:33:MSE:CA	1:C:33:MSE:HE1	2.32	0.51
1:B:80:PHE:CE2	1:B:84:LEU:HD11	2.46	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:95:ASP:CG	1:C:98:LYS:HG2	2.36	0.51
1:D:43:ILE:HD13	1:D:119:MSE:CE	2.41	0.51
1:A:134:PHE:O	1:A:137:GLU:HB2	2.10	0.51
1:D:111:HIS:HB3	1:D:117:LYS:NZ	2.26	0.51
1:B:61:MSE:O	1:B:65:LEU:HG	2.10	0.50
1:C:24:GLN:HG3	1:D:37:ILE:HG21	1.91	0.50
1:D:54:TYR:CD2	1:D:106:ARG:HG3	2.47	0.50
1:C:131:GLU:O	1:C:134:PHE:HB3	2.12	0.50
1:B:7:ASN:O	1:B:8:LEU:C	2.53	0.50
1:A:75:ILE:HG12	1:A:78:TYR:CE1	2.46	0.50
1:A:20:GLN:CB	1:B:37:ILE:CD1	2.89	0.49
1:C:37:ILE:HG21	1:D:24:GLN:HG3	1.94	0.49
1:C:66:GLN:NE2	1:C:73:ASP:O	2.45	0.49
1:A:33:MSE:HE1	1:D:33:MSE:HA	1.95	0.49
1:A:54:TYR:CD2	1:A:106:ARG:HD2	2.47	0.49
1:C:103:ARG:C	1:C:105:MSE:N	2.66	0.49
1:D:47:ILE:HD13	1:D:112:THR:CG2	2.43	0.49
1:D:62:LYS:HE3	1:D:76:GLY:HA2	1.93	0.49
1:A:70:ILE:C	1:A:72:THR:H	2.20	0.49
1:B:95:ASP:OD2	1:B:97:SER:HB2	2.12	0.49
1:C:32:ASP:C	1:C:34:GLN:H	2.19	0.49
1:B:73:ASP:O	1:B:74:ASP:O	2.31	0.49
1:B:18:LEU:O	1:B:22:VAL:HG13	2.13	0.49
1:B:39:GLN:O	1:B:43:ILE:HD12	2.13	0.49
1:A:7:ASN:HD21	1:A:60:MSE:HE3	1.77	0.48
1:A:65:LEU:HD22	1:A:86:GLU:HB3	1.95	0.48
1:A:70:ILE:CD1	1:A:75:ILE:HG13	2.43	0.48
1:C:20:GLN:CB	1:D:37:ILE:HD12	2.25	0.48
1:C:38:VAL:HG21	1:D:30:TRP:CH2	2.49	0.48
1:C:37:ILE:HD11	1:D:21:THR:N	2.29	0.48
1:C:40:ASP:HB2	1:C:115:GLN:NE2	2.28	0.48
1:C:111:HIS:HA	1:C:113:TYR:CE2	2.48	0.48
1:A:7:ASN:C	1:A:7:ASN:HD22	2.21	0.48
1:C:75:ILE:C	1:C:77:ALA:H	2.21	0.48
1:A:7:ASN:ND2	1:A:60:MSE:CE	2.77	0.48
1:C:64:GLN:C	1:C:66:GLN:H	2.21	0.48
1:C:55:GLU:OE2	1:C:106:ARG:NH1	2.45	0.47
1:C:82:ASP:O	1:C:83:ILE:C	2.57	0.47
1:C:58:LEU:HD11	1:C:80:PHE:CE1	2.49	0.47
1:A:90:PHE:HD1	1:A:90:PHE:N	2.13	0.47
1:A:10:VAL:CG2	1:A:11:LEU:N	2.77	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:125:ILE:O	1:C:126:ASP:C	2.58	0.47
1:D:28:ARG:HD2	2:D:1095:HOH:O	2.14	0.47
1:A:95:ASP:CG	1:A:98:LYS:HZ2	2.22	0.47
1:B:103:ARG:O	1:B:107:ASN:ND2	2.48	0.47
1:C:10:VAL:CG1	1:C:56:LEU:HD21	2.42	0.47
1:C:69:ALA:C	1:C:71:ASN:N	2.72	0.47
1:D:112:THR:C	1:D:114:ASP:N	2.69	0.47
1:A:8:LEU:HD12	1:A:8:LEU:HA	1.76	0.47
1:B:68:ASP:HB3	1:B:90:PHE:HE2	1.79	0.47
1:C:74:ASP:C	1:C:76:GLY:N	2.73	0.47
1:A:76:GLY:C	1:A:78:TYR:H	2.23	0.46
1:C:61:MSE:HE1	1:C:93:ILE:HD11	1.97	0.46
1:D:78:TYR:CG	1:D:82:ASP:OD2	2.68	0.46
1:A:20:GLN:C	1:B:37:ILE:CD1	2.88	0.46
1:D:78:TYR:CE1	1:D:82:ASP:OD2	2.68	0.46
1:A:90:PHE:N	1:A:90:PHE:CD1	2.83	0.46
1:C:106:ARG:HD2	1:C:106:ARG:C	2.30	0.46
1:C:134:PHE:O	1:C:137:GLU:N	2.48	0.46
1:C:106:ARG:HH11	1:C:106:ARG:HG2	1.79	0.46
1:A:42:LEU:HD23	1:B:41:THR:HG22	1.98	0.46
1:D:135:LEU:HD12	1:D:139:LEU:HG	1.96	0.46
1:D:51:GLU:OE1	1:D:109:THR:CB	2.64	0.46
1:A:20:GLN:CB	1:B:37:ILE:HD12	2.37	0.46
1:A:80:PHE:CZ	1:A:84:LEU:HD11	2.51	0.46
1:C:111:HIS:CD2	1:C:117:LYS:NZ	2.83	0.46
1:C:105:MSE:SE	1:C:124:GLN:OE1	2.85	0.45
1:D:57:SER:HB3	1:D:135:LEU:CD2	2.46	0.45
1:A:52:PHE:HZ	1:B:112:THR:HG22	1.81	0.45
1:B:58:LEU:HD23	1:B:58:LEU:HA	1.80	0.45
1:A:57:SER:HB3	1:A:135:LEU:CD2	2.47	0.45
1:C:103:ARG:HH11	1:C:103:ARG:CG	2.30	0.45
1:B:7:ASN:O	1:B:9:ASN:N	2.50	0.45
1:A:80:PHE:CE2	1:A:84:LEU:HD11	2.52	0.45
1:A:70:ILE:CG2	1:A:72:THR:HA	2.47	0.45
1:D:60:MSE:HE1	1:D:63:ARG:NH1	2.31	0.45
1:B:87:ALA:O	1:B:92:LEU:HB2	2.17	0.44
1:D:55:GLU:CG	1:D:106:ARG:HH21	2.30	0.44
1:B:27:ASP:OD1	1:B:27:ASP:O	2.34	0.44
1:D:18:LEU:O	1:D:19:GLU:C	2.60	0.44
1:B:112:THR:O	1:B:112:THR:CG2	2.65	0.44
1:B:68:ASP:HB3	1:B:90:PHE:CE2	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:27:ASP:O	1:C:28:ARG:C	2.59	0.44
1:C:125:ILE:HA	1:C:128:PHE:HB3	1.98	0.44
1:C:78:TYR:HB3	1:C:83:ILE:CD1	2.48	0.44
1:D:112:THR:O	1:D:113:TYR:C	2.59	0.43
1:C:58:LEU:HA	1:C:58:LEU:HD23	1.78	0.43
1:D:81:LYS:NZ	1:D:100:VAL:HG22	2.33	0.43
1:B:50:PHE:CE2	1:B:105:MSE:HE2	2.53	0.43
1:C:45:GLY:O	1:C:48:GLN:HB3	2.18	0.43
1:B:50:PHE:CE1	1:B:105:MSE:HE2	2.51	0.43
1:A:52:PHE:O	1:A:56:LEU:HD22	2.19	0.43
1:B:11:LEU:HD13	1:B:56:LEU:HB3	1.99	0.43
1:C:72:THR:HG22	1:C:72:THR:O	2.18	0.43
1:D:65:LEU:HD23	2:D:1080:HOH:O	2.18	0.43
1:C:75:ILE:HG23	1:C:75:ILE:HD12	1.70	0.43
1:C:97:SER:HA	1:C:100:VAL:HB	2.00	0.43
1:B:7:ASN:C	1:B:9:ASN:N	2.76	0.43
1:C:75:ILE:C	1:C:77:ALA:N	2.76	0.43
1:A:11:LEU:HD22	1:A:60:MSE:HG3	2.00	0.42
1:A:66:GLN:O	1:A:70:ILE:HG22	2.19	0.42
1:B:27:ASP:OD1	1:B:27:ASP:C	2.62	0.42
1:C:126:ASP:HA	1:C:129:LEU:HD12	2.01	0.42
1:C:24:GLN:HG3	1:D:37:ILE:CG2	2.48	0.42
1:C:80:PHE:CE2	1:C:84:LEU:HD11	2.54	0.42
1:D:75:ILE:C	1:D:77:ALA:H	2.25	0.42
1:D:81:LYS:HZ2	1:D:100:VAL:CG2	2.31	0.42
1:C:65:LEU:CD2	1:C:86:GLU:O	2.64	0.42
1:D:62:LYS:HE2	1:D:76:GLY:HA2	2.02	0.42
1:B:60:MSE:HE2	1:B:60:MSE:CA	2.44	0.42
1:C:96:MSE:HG3	1:C:97:SER:N	2.35	0.42
1:C:102:TYR:HA	1:C:105:MSE:CE	2.48	0.42
1:D:57:SER:HB3	1:D:135:LEU:HD23	2.00	0.42
1:B:125:ILE:O	1:B:128:PHE:HB3	2.19	0.42
1:C:37:ILE:HG21	1:D:24:GLN:HG2	2.02	0.42
1:C:40:ASP:HB2	1:C:115:GLN:HE22	1.85	0.42
1:D:27:ASP:O	1:D:29:ASN:N	2.52	0.42
1:B:33:MSE:HA	1:C:33:MSE:CE	2.34	0.42
1:C:112:THR:C	1:C:114:ASP:N	2.77	0.42
1:C:134:PHE:HA	1:C:137:GLU:CD	2.45	0.42
1:D:95:ASP:OD2	1:D:97:SER:N	2.52	0.42
1:A:18:LEU:O	1:A:21:THR:HB	2.19	0.42
1:A:80:PHE:O	1:A:83:ILE:HB	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:54:TYR:C	1:C:54:TYR:CD2	2.98	0.42
1:A:50:PHE:CE2	1:A:105:MSE:HE2	2.55	0.41
1:A:70:ILE:C	1:A:72:THR:N	2.77	0.41
1:A:7:ASN:HB3	1:A:63:ARG:HH22	1.85	0.41
1:D:26:SER:C	1:D:28:ARG:N	2.78	0.41
1:D:111:HIS:C	1:D:113:TYR:N	2.78	0.41
1:C:56:LEU:HD12	1:C:56:LEU:HA	1.72	0.41
1:A:105:MSE:O	1:A:108:ILE:HG12	2.20	0.41
1:C:106:ARG:NH1	1:C:106:ARG:HG2	2.35	0.41
1:A:10:VAL:O	1:A:11:LEU:C	2.63	0.41
1:C:66:GLN:CG	1:C:73:ASP:OD1	2.69	0.41
1:C:137:GLU:O	1:C:140:ARG:HB3	2.21	0.41
1:A:105:MSE:SE	1:A:124:GLN:OE1	2.89	0.40
1:D:69:ALA:HB2	1:D:75:ILE:CD1	2.51	0.40
1:C:114:ASP:OD2	1:C:117:LYS:HG3	2.21	0.40
1:B:105:MSE:O	1:B:108:ILE:HG12	2.22	0.40
1:C:18:LEU:HD23	1:C:18:LEU:HA	1.93	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	133/146 (91%)	123 (92%)	8 (6%)	2 (2%)	8	12
1	B	133/146 (91%)	121 (91%)	5 (4%)	7 (5%)	1	1
1	C	133/146 (91%)	112 (84%)	17 (13%)	4 (3%)	3	3
1	D	134/146 (92%)	121 (90%)	9 (7%)	4 (3%)	3	3
All	All	533/584 (91%)	477 (90%)	39 (7%)	17 (3%)	3	3

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	27	ASP
1	B	74	ASP
1	C	27	ASP
1	D	27	ASP
1	D	28	ARG
1	A	74	ASP
1	B	28	ARG
1	C	74	ASP
1	C	75	ILE
1	D	111	HIS
1	B	8	LEU
1	B	79	GLY
1	B	112	THR
1	B	76	GLY
1	D	120	ALA
1	C	82	ASP
1	A	75	ILE

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	119/122 (98%)	105 (88%)	14 (12%)	5	7
1	B	119/122 (98%)	106 (89%)	13 (11%)	6	9
1	C	119/122 (98%)	106 (89%)	13 (11%)	6	9
1	D	120/122 (98%)	104 (87%)	16 (13%)	4	5
All	All	477/488 (98%)	421 (88%)	56 (12%)	5	7

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	ASN
1	A	9	ASN
1	A	18	LEU
1	A	55	GLU
1	A	56	LEU

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Mol	Chain	Res	Type
1	A	58	LEU
1	A	59	LYS
1	A	70	ILE
1	A	86	GLU
1	A	89	ARG
1	A	90	PHE
1	A	106	ARG
1	A	112	THR
1	A	137	GLU
1	B	18	LEU
1	B	29	ASN
1	B	33	MSE
1	B	40	ASP
1	B	56	LEU
1	B	72	THR
1	B	74	ASP
1	B	105	MSE
1	B	106	ARG
1	B	112	THR
1	B	119	MSE
1	B	126	ASP
1	B	136	LEU
1	C	8	LEU
1	C	10	VAL
1	C	18	LEU
1	C	28	ARG
1	C	40	ASP
1	C	43	ILE
1	C	73	ASP
1	C	90	PHE
1	C	97	SER
1	C	103	ARG
1	C	132	SER
1	C	136	LEU
1	C	139	LEU
1	D	17	SER
1	D	18	LEU
1	D	28	ARG
1	D	40	ASP
1	D	50	PHE
1	D	58	LEU
1	D	63	ARG

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Mol	Chain	Res	Type
1	D	70	ILE
1	D	96	MSE
1	D	97	SER
1	D	100	VAL
1	D	113	TYR
1	D	115	GLN
1	D	119	MSE
1	D	132	SER
1	D	136	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	7	ASN
1	A	9	ASN
1	A	24	GLN
1	A	71	ASN
1	A	115	GLN
1	B	20	GLN
1	B	66	GLN
1	B	107	ASN
1	B	115	GLN
1	C	7	ASN
1	C	39	GLN
1	C	66	GLN
1	C	67	GLN
1	C	111	HIS
1	C	115	GLN
1	D	7	ASN
1	D	48	GLN
1	D	64	GLN
1	D	115	GLN

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

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

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	129/146 (88%)	-0.15	5 (3%) 43 39	23, 42, 82, 85	0
1	B	129/146 (88%)	-0.23	0 100 100	21, 50, 80, 83	0
1	C	129/146 (88%)	0.25	3 (2%) 61 57	38, 72, 83, 85	0
1	D	130/146 (89%)	0.01	0 100 100	34, 59, 82, 84	0
All	All	517/584 (88%)	-0.03	8 (1%) 72 68	21, 56, 82, 85	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	72	THR	3.0
1	A	72	THR	2.7
1	A	141	GLN	2.5
1	A	69	ALA	2.4
1	A	68	ASP	2.3
1	C	74	ASP	2.3
1	A	74	ASP	2.3
1	C	73	ASP	2.3

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.