



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 04:24 AM UTC

PDB ID : 2CV9 / pdb_00002cv9
Title : Crystal structure of a hypothetical protein from *Thermus thermophilus* HB8
Authors : Kanagawa, M.; Yokoyama, S.; Kuramitsu, S.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2005-06-01
Resolution : 2.50 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

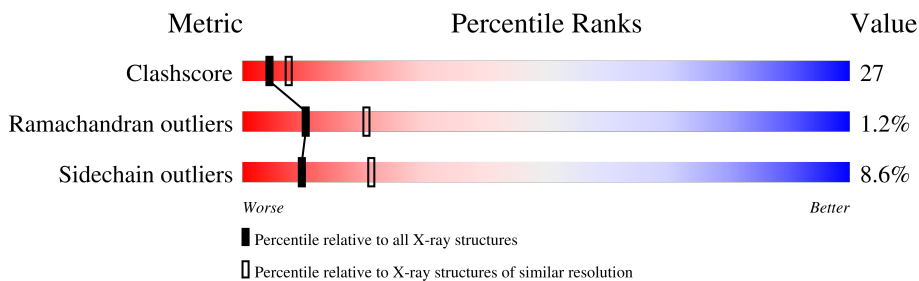
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	252	
1	B	252	
1	C	252	
1	D	252	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8338 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 TTHA0625.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	252	Total 1977	C 1259	N 351	O 361	Se 6	0	0	0
1	B	252	Total 1977	C 1259	N 351	O 361	Se 6	0	0	0
1	C	252	Total 1977	C 1259	N 351	O 361	Se 6	0	0	0
1	D	252	Total 1977	C 1259	N 351	O 361	Se 6	0	0	0

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	modified residue	UNP Q5SKL8
A	10	MSE	MET	modified residue	UNP Q5SKL8
A	113	MSE	MET	modified residue	UNP Q5SKL8
A	118	MSE	MET	modified residue	UNP Q5SKL8
A	153	MSE	MET	modified residue	UNP Q5SKL8
A	193	MSE	MET	modified residue	UNP Q5SKL8
B	1	MSE	MET	modified residue	UNP Q5SKL8
B	10	MSE	MET	modified residue	UNP Q5SKL8
B	113	MSE	MET	modified residue	UNP Q5SKL8
B	118	MSE	MET	modified residue	UNP Q5SKL8
B	153	MSE	MET	modified residue	UNP Q5SKL8
B	193	MSE	MET	modified residue	UNP Q5SKL8
C	1	MSE	MET	modified residue	UNP Q5SKL8
C	10	MSE	MET	modified residue	UNP Q5SKL8
C	113	MSE	MET	modified residue	UNP Q5SKL8
C	118	MSE	MET	modified residue	UNP Q5SKL8
C	153	MSE	MET	modified residue	UNP Q5SKL8
C	193	MSE	MET	modified residue	UNP Q5SKL8
D	1	MSE	MET	modified residue	UNP Q5SKL8
D	10	MSE	MET	modified residue	UNP Q5SKL8
D	113	MSE	MET	modified residue	UNP Q5SKL8

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Chain	Residue	Modelled	Actual	Comment	Reference
D	118	MSE	MET	modified residue	UNP Q5SKL8
D	153	MSE	MET	modified residue	UNP Q5SKL8
D	193	MSE	MET	modified residue	UNP Q5SKL8

- Molecule 2 is water.

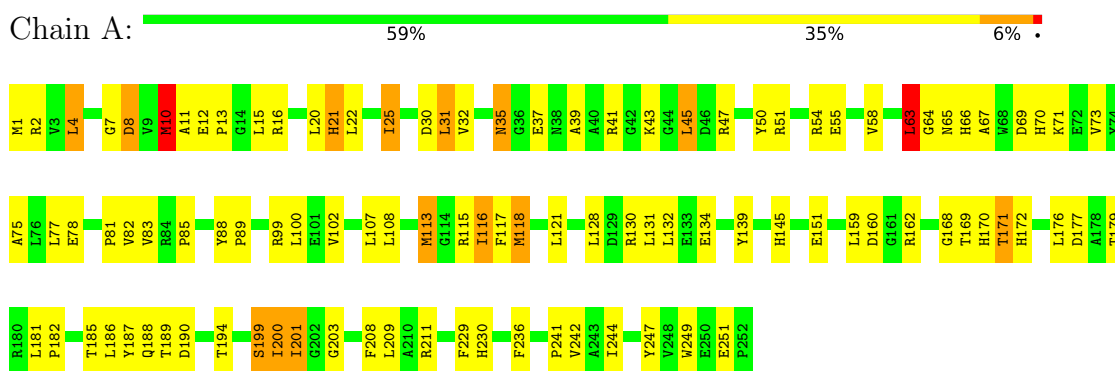
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	120	Total 120	O 120	0	0
2	B	108	Total 108	O 108	0	0
2	C	100	Total 100	O 100	0	0
2	D	102	Total 102	O 102	0	0

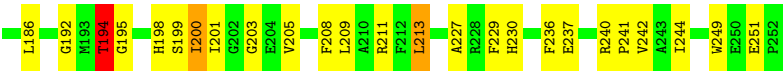
3 Residue-property plots [i](#)

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

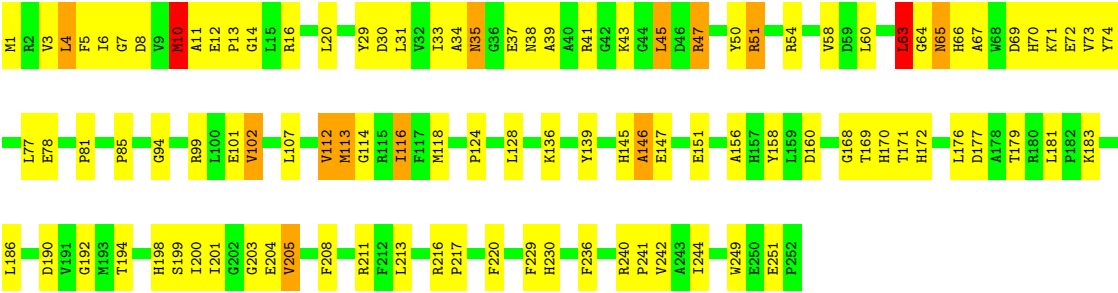
Note EDS was not executed.

- Molecule 1: hypothetical protein TTHA0625





● Molecule 1: hypothetical protein TTHA0625



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	76.48Å 76.48Å 193.94Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.78 – 2.50	Depositor
% Data completeness (in resolution range)	97.7 (19.78-2.50)	Depositor
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.197 , 0.247	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	8338	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

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	0.47	0/2021	0.95	3/2735 (0.1%)
1	B	0.45	0/2021	1.00	6/2735 (0.2%)
1	C	0.45	0/2021	1.02	8/2735 (0.3%)
1	D	0.46	0/2021	1.02	5/2735 (0.2%)
All	All	0.46	0/8084	1.00	22/10940 (0.2%)

There are no bond length outliers.

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	21	HIS	N-CA-C	8.51	121.78	111.40
1	B	242	VAL	N-CA-C	7.46	119.13	111.00
1	D	242	VAL	N-CA-C	7.21	118.15	111.45
1	D	63	LEU	N-CA-C	7.04	115.51	108.75
1	C	242	VAL	N-CA-C	6.79	117.76	111.45
1	A	242	VAL	N-CA-C	6.18	118.14	111.58
1	C	21	HIS	N-CA-C	6.14	120.79	113.12
1	D	205	VAL	N-CA-C	6.13	116.60	110.23
1	D	198	HIS	N-CA-C	-6.05	99.76	109.39
1	B	39	ALA	N-CA-C	5.93	118.51	111.33
1	C	195	GLY	N-CA-C	-5.74	102.88	110.74
1	B	11	ALA	N-CA-C	5.67	114.20	108.75
1	A	21	HIS	N-CA-C	5.51	119.97	112.04
1	C	30	ASP	N-CA-C	-5.48	106.27	113.12
1	C	198	HIS	N-CA-C	-5.48	100.67	109.39
1	C	148	ALA	N-CA-C	5.23	117.25	109.25
1	A	63	LEU	CB-CA-C	-5.18	108.98	116.54
1	C	194	THR	N-CA-C	-5.16	100.11	108.52
1	B	88	TYR	N-CA-C	-5.15	102.78	110.20
1	B	198	HIS	N-CA-C	-5.14	102.73	110.70
1	C	213	LEU	N-CA-C	5.10	117.97	111.69
1	D	146	ALA	N-CA-C	5.09	116.03	108.60

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	1977	0	1958	103	0
1	B	1977	0	1958	109	0
1	C	1977	0	1958	114	0
1	D	1977	0	1958	104	0
2	A	120	0	0	7	0
2	B	108	0	0	7	0
2	C	100	0	0	4	0
2	D	102	0	0	4	0
All	All	8338	0	7832	419	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 27.

All (419) 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:171:THR:HB	1:B:190:ASP:OD1	1.59	1.00
1:D:194:THR:HG22	1:D:230:HIS:H	1.27	0.99
1:D:240:ARG:HD3	1:D:241:PRO:HD2	1.46	0.96
1:D:200:ILE:HD12	1:D:205:VAL:HA	1.49	0.94
1:D:1:MSE:HE1	1:D:107:LEU:HD22	1.53	0.90
1:A:171:THR:HB	1:A:190:ASP:OD1	1.72	0.88
1:B:113:MSE:HE2	1:B:145:HIS:HB3	1.54	0.88
1:B:200:ILE:HD12	1:B:205:VAL:HA	1.56	0.88
1:C:172:HIS:ND1	1:C:201:ILE:HG23	1.90	0.86
1:D:113:MSE:HE2	1:D:145:HIS:HB3	1.58	0.85
1:B:150:SER:HA	1:C:201:ILE:HD13	1.57	0.84
1:D:1:MSE:HE3	1:D:3:VAL:HG21	1.58	0.83
1:D:1:MSE:HE2	1:D:236:PHE:HD2	1.43	0.83
1:A:113:MSE:HE2	1:A:145:HIS:HB3	1.61	0.83
1:B:30:ASP:HB3	1:B:102:VAL:HG11	1.59	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:63:LEU:HD23	1:C:67:ALA:HB1	1.62	0.81
1:D:200:ILE:N	1:D:200:ILE:HD13	1.95	0.81
1:C:1:MSE:HE1	1:C:107:LEU:HD22	1.62	0.81
1:D:64:GLY:HA2	1:D:113:MSE:HE1	1.62	0.80
1:D:200:ILE:CD1	1:D:205:VAL:HA	2.13	0.79
1:C:1:MSE:HE3	1:C:3:VAL:HG21	1.65	0.79
1:D:186:LEU:HD13	1:D:244:ILE:HB	1.65	0.78
1:A:30:ASP:HB3	1:A:102:VAL:HG11	1.67	0.77
1:C:65:ASN:H	1:C:65:ASN:HD22	1.29	0.77
1:C:16:ARG:HD2	1:C:251:GLU:OE1	1.87	0.75
1:A:12:GLU:CD	1:A:12:GLU:H	1.96	0.74
1:C:200:ILE:HD12	1:C:205:VAL:HA	1.68	0.74
1:B:200:ILE:H	1:B:200:ILE:HD13	1.53	0.73
1:C:64:GLY:HA2	1:C:113:MSE:HE1	1.71	0.73
1:C:11:ALA:N	2:C:309:HOH:O	2.21	0.73
1:C:112:VAL:HG21	1:C:128:LEU:CD2	2.18	0.73
1:D:71:LYS:HD2	1:D:71:LYS:C	2.14	0.72
1:D:200:ILE:HD13	1:D:200:ILE:H	1.55	0.72
1:D:11:ALA:N	2:D:305:HOH:O	2.21	0.72
1:D:200:ILE:O	1:D:201:ILE:HG22	1.90	0.72
1:C:201:ILE:O	1:C:201:ILE:HG22	1.89	0.71
1:D:65:ASN:HD22	1:D:65:ASN:H	1.39	0.71
1:A:43:LYS:HE3	1:A:69:ASP:CG	2.16	0.71
1:C:71:LYS:C	1:C:71:LYS:HD2	2.16	0.70
1:C:1:MSE:HE2	1:C:236:PHE:HD2	1.55	0.70
1:C:113:MSE:HE2	1:C:145:HIS:HB3	1.71	0.70
1:D:16:ARG:HD2	1:D:251:GLU:OE1	1.91	0.70
1:B:12:GLU:H	1:B:12:GLU:CD	1.98	0.70
1:B:47:ARG:HD2	1:B:51:ARG:NH2	2.07	0.70
1:B:16:ARG:O	1:B:20:LEU:HD13	1.92	0.69
1:A:43:LYS:HB3	1:A:69:ASP:CG	2.17	0.69
1:D:200:ILE:HG13	1:D:208:PHE:HB2	1.74	0.69
1:A:200:ILE:HD13	1:A:200:ILE:H	1.56	0.69
1:B:31:LEU:HD13	1:B:102:VAL:HG21	1.75	0.69
1:B:7:GLY:O	1:B:194:THR:HG23	1.93	0.68
1:A:4:LEU:HB3	1:A:32:VAL:HG22	1.74	0.68
1:B:1:MSE:HE1	1:B:107:LEU:HD22	1.76	0.68
1:A:63:LEU:HB3	1:A:67:ALA:HB2	1.76	0.68
1:A:47:ARG:HD2	1:A:51:ARG:NH2	2.09	0.68
1:A:70:HIS:O	1:A:73:VAL:HG22	1.93	0.68
1:C:7:GLY:O	1:C:194:THR:HG23	1.95	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:200:ILE:HG13	1:C:208:PHE:CB	2.25	0.67
1:A:16:ARG:O	1:A:20:LEU:HD13	1.95	0.67
1:C:200:ILE:HD13	1:C:200:ILE:N	2.09	0.67
1:D:1:MSE:HE3	1:D:3:VAL:CG2	2.24	0.67
1:B:200:ILE:O	1:B:201:ILE:HG22	1.94	0.66
1:D:112:VAL:HG21	1:D:128:LEU:CD2	2.26	0.66
1:B:1:MSE:HE1	1:B:107:LEU:HB2	1.78	0.66
1:C:200:ILE:HG13	1:C:208:PHE:HB2	1.78	0.66
1:B:200:ILE:O	1:B:200:ILE:HG12	1.96	0.66
1:D:200:ILE:HG13	1:D:208:PHE:CB	2.26	0.66
1:D:112:VAL:HG21	1:D:128:LEU:HD21	1.79	0.65
1:D:30:ASP:HB3	1:D:102:VAL:HG11	1.77	0.65
1:B:200:ILE:HD13	1:B:200:ILE:N	2.12	0.65
1:D:65:ASN:HD22	1:D:65:ASN:N	1.94	0.65
1:C:200:ILE:CD1	1:C:205:VAL:HA	2.26	0.65
1:C:240:ARG:HD3	1:C:241:PRO:HD2	1.79	0.65
1:B:63:LEU:HB3	1:B:67:ALA:HB2	1.80	0.64
1:C:194:THR:HG22	1:C:230:HIS:H	1.63	0.64
1:B:37:GLU:HG2	1:B:66:HIS:HB2	1.80	0.63
1:A:21:HIS:HE1	1:A:247:TYR:OH	1.81	0.63
1:C:200:ILE:HG12	1:C:200:ILE:O	1.99	0.63
1:B:21:HIS:HE1	1:B:247:TYR:OH	1.81	0.63
1:C:12:GLU:HB2	1:C:13:PRO:HD3	1.81	0.63
1:D:77:LEU:HG	1:D:94:GLY:HA3	1.80	0.63
1:D:1:MSE:HE2	1:D:236:PHE:CD2	2.30	0.63
1:D:16:ARG:O	1:D:20:LEU:HB2	1.99	0.62
1:C:15:LEU:HD21	1:C:52:LEU:HD12	1.81	0.62
1:D:116:ILE:HG23	1:D:151:GLU:OE1	1.98	0.62
1:B:168:GLY:HA3	1:B:189:THR:O	1.99	0.62
1:B:22:LEU:HA	1:B:25:ILE:HD11	1.81	0.62
1:D:124:PRO:O	1:D:128:LEU:HD23	1.98	0.62
1:C:1:MSE:HE3	1:C:3:VAL:CG2	2.30	0.62
1:A:22:LEU:HA	1:A:25:ILE:HD11	1.81	0.61
1:A:1:MSE:HE1	1:A:107:LEU:HD13	1.81	0.61
1:B:71:LYS:HD2	1:B:71:LYS:C	2.25	0.61
1:C:47:ARG:HH11	1:C:47:ARG:HB3	1.64	0.61
1:C:201:ILE:O	1:C:201:ILE:CG2	2.48	0.61
1:B:43:LYS:HB3	1:B:69:ASP:CG	2.26	0.61
1:B:139:TYR:CE1	1:B:236:PHE:HB3	2.36	0.61
1:C:101:GLU:O	1:C:102:VAL:HG23	2.01	0.61
1:A:43:LYS:HE3	1:A:69:ASP:HB2	1.83	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:118:MSE:HE1	1:C:146:ALA:HA	1.83	0.61
1:D:60:LEU:HD12	1:D:81:PRO:O	2.01	0.60
1:C:114:GLY:C	1:C:118:MSE:HE3	2.26	0.60
1:C:30:ASP:HB3	1:C:102:VAL:HG11	1.84	0.60
1:D:145:HIS:HD2	1:D:169:THR:OG1	1.85	0.60
1:B:113:MSE:CE	1:B:145:HIS:HB3	2.29	0.60
1:D:200:ILE:O	1:D:200:ILE:HG12	2.02	0.59
1:C:71:LYS:HD2	1:C:71:LYS:O	2.01	0.59
1:A:43:LYS:HB3	1:A:69:ASP:OD2	2.02	0.59
1:B:1:MSE:HE2	1:B:236:PHE:HD2	1.68	0.59
1:B:41:ARG:HG2	1:B:41:ARG:HH21	1.68	0.59
1:B:70:HIS:O	1:B:73:VAL:HG22	2.03	0.59
1:A:168:GLY:HA3	1:A:189:THR:O	2.01	0.59
1:A:115:ARG:HG2	2:A:289:HOH:O	2.02	0.59
1:C:63:LEU:HD11	1:C:82:VAL:CG1	2.33	0.59
1:B:153:MSE:HE3	1:C:201:ILE:CD1	2.32	0.59
1:B:43:LYS:HB3	1:B:69:ASP:OD2	2.03	0.58
1:D:63:LEU:HB3	1:D:67:ALA:HB2	1.85	0.58
1:A:7:GLY:O	1:A:194:THR:HG23	2.03	0.58
1:C:65:ASN:HD22	1:C:65:ASN:N	2.01	0.58
1:D:171:THR:HB	1:D:190:ASP:OD1	2.03	0.58
1:C:112:VAL:HG21	1:C:128:LEU:HD21	1.83	0.58
1:D:194:THR:HG22	1:D:230:HIS:N	2.08	0.58
1:A:22:LEU:HA	1:A:25:ILE:CD1	2.34	0.58
1:C:41:ARG:HG2	1:C:41:ARG:HH21	1.68	0.58
1:A:37:GLU:HG2	1:A:66:HIS:HB2	1.85	0.58
1:B:1:MSE:HE3	1:B:3:VAL:HG21	1.84	0.58
1:D:50:TYR:OH	1:D:81:PRO:HD2	2.03	0.58
1:D:139:TYR:CE1	1:D:236:PHE:HB3	2.38	0.58
1:A:31:LEU:HG	1:A:100:LEU:HD13	1.85	0.57
1:A:1:MSE:HE1	1:A:107:LEU:HB2	1.86	0.57
1:A:194:THR:HG22	1:A:230:HIS:H	1.69	0.57
1:D:1:MSE:CE	1:D:107:LEU:HD22	2.32	0.57
1:A:47:ARG:O	1:A:51:ARG:HG3	2.04	0.57
1:A:65:ASN:HD21	1:A:170:HIS:CE1	2.22	0.57
1:C:139:TYR:CE1	1:C:236:PHE:HB3	2.39	0.57
1:B:10:MSE:HB3	2:B:254:HOH:O	2.05	0.57
1:C:199:SER:HB2	1:C:203:GLY:O	2.04	0.57
1:A:201:ILE:HG22	1:A:201:ILE:O	2.04	0.56
1:D:172:HIS:ND1	1:D:201:ILE:HG13	2.19	0.56
1:B:25:ILE:HD13	1:B:25:ILE:H	1.70	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:169:THR:O	1:C:170:HIS:HB3	2.05	0.56
1:A:43:LYS:HE3	1:A:69:ASP:CB	2.36	0.56
1:A:200:ILE:HD13	1:A:200:ILE:N	2.21	0.56
1:C:229:PHE:HB3	1:C:249:TRP:HB3	1.88	0.56
1:D:101:GLU:O	1:D:102:VAL:HG23	2.05	0.56
1:D:118:MSE:HE2	1:D:147:GLU:H	1.71	0.56
1:C:22:LEU:HA	1:C:25:ILE:CD1	2.35	0.56
1:D:118:MSE:HE1	1:D:146:ALA:HA	1.88	0.56
1:A:229:PHE:HB3	1:A:249:TRP:HB3	1.87	0.55
2:A:350:HOH:O	1:D:213:LEU:HA	2.05	0.55
1:A:201:ILE:O	1:A:201:ILE:CG2	2.54	0.55
1:D:177:ASP:O	1:D:179:THR:HG23	2.06	0.55
1:A:117:PHE:C	1:A:118:MSE:HG2	2.31	0.54
1:D:41:ARG:HD3	2:D:321:HOH:O	2.06	0.54
1:A:108:LEU:HD21	1:A:131:LEU:HD11	1.89	0.54
1:C:62:SER:C	1:C:63:LEU:HD12	2.33	0.54
1:C:177:ASP:O	1:C:179:THR:HG23	2.07	0.54
1:C:145:HIS:HD2	1:C:169:THR:OG1	1.90	0.54
1:D:7:GLY:HA3	1:D:192:GLY:O	2.07	0.54
1:D:216:ARG:HB3	1:D:217:PRO:HD2	1.90	0.54
1:D:201:ILE:HG23	1:D:201:ILE:O	2.07	0.54
1:B:31:LEU:HG	1:B:100:LEU:HD13	1.90	0.53
1:C:16:ARG:O	1:C:20:LEU:HB2	2.07	0.53
1:D:160:ASP:CG	1:D:183:LYS:HB2	2.33	0.53
1:B:12:GLU:OE2	1:B:41:ARG:HA	2.08	0.53
1:B:201:ILE:O	1:B:201:ILE:HG23	2.07	0.53
1:A:2:ARG:HD3	2:A:322:HOH:O	2.08	0.53
1:B:26:ARG:HD3	1:B:57:GLY:HA3	1.90	0.53
1:D:229:PHE:HB3	1:D:249:TRP:HB3	1.89	0.53
1:A:131:LEU:HD13	1:A:131:LEU:C	2.34	0.53
1:A:199:SER:HB2	1:A:203:GLY:O	2.07	0.53
1:D:74:TYR:O	1:D:78:GLU:HG2	2.08	0.53
1:A:45:LEU:O	1:A:73:VAL:HG13	2.09	0.53
1:D:113:MSE:HE2	1:D:145:HIS:CB	2.36	0.53
1:C:134:GLU:HG3	1:C:135:GLU:HG2	1.91	0.53
1:C:1:MSE:HE2	1:C:236:PHE:CD2	2.40	0.53
1:C:50:TYR:OH	1:C:81:PRO:HD2	2.09	0.53
1:D:1:MSE:HE1	1:D:107:LEU:CD2	2.34	0.53
1:A:43:LYS:HE3	1:A:69:ASP:OD1	2.09	0.53
1:B:194:THR:HG22	1:B:230:HIS:H	1.73	0.53
1:B:229:PHE:HB3	1:B:249:TRP:HB3	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:41:ARG:HG2	1:B:41:ARG:NH2	2.25	0.52
1:C:87:ASN:ND2	1:C:121:LEU:HB2	2.24	0.52
1:B:45:LEU:O	1:B:73:VAL:HG13	2.09	0.52
1:B:47:ARG:O	1:B:51:ARG:HG3	2.10	0.52
1:D:169:THR:O	1:D:170:HIS:HB3	2.09	0.52
1:B:70:HIS:ND1	1:B:72:GLU:HB2	2.24	0.52
1:C:108:LEU:HD21	1:C:131:LEU:CD1	2.40	0.52
1:C:112:VAL:HG21	1:C:128:LEU:HD22	1.89	0.52
1:A:10:MSE:HB3	2:A:253:HOH:O	2.10	0.51
1:B:16:ARG:NH2	2:B:266:HOH:O	2.43	0.51
1:A:172:HIS:ND1	1:A:201:ILE:HG23	2.25	0.51
1:C:155:LEU:O	1:C:159:LEU:HD13	2.09	0.51
1:B:108:LEU:HD21	1:B:131:LEU:HD11	1.92	0.51
1:B:168:GLY:O	1:B:190:ASP:HA	2.11	0.51
1:D:35:ASN:ND2	1:D:37:GLU:H	2.09	0.51
1:A:200:ILE:HG12	1:A:200:ILE:O	2.11	0.51
1:D:199:SER:HB2	1:D:203:GLY:O	2.10	0.51
1:A:113:MSE:HG2	1:A:121:LEU:CD1	2.40	0.51
1:B:10:MSE:O	1:B:38:ASN:O	2.28	0.51
1:A:35:ASN:C	1:A:35:ASN:HD22	2.19	0.51
1:B:1:MSE:CE	1:B:107:LEU:HD22	2.40	0.51
1:A:71:LYS:HD2	1:A:71:LYS:C	2.36	0.50
1:B:130:ARG:O	1:B:134:GLU:HG2	2.11	0.50
1:B:117:PHE:C	1:B:118:MSE:HG2	2.36	0.50
1:B:158:TYR:HB2	1:C:209:LEU:HD11	1.93	0.50
1:B:22:LEU:HA	1:B:25:ILE:CD1	2.41	0.50
1:C:113:MSE:HE2	1:C:145:HIS:ND1	2.26	0.50
1:A:139:TYR:CE1	1:A:236:PHE:HB3	2.46	0.50
1:A:176:LEU:HD13	1:D:176:LEU:HD13	1.93	0.50
1:C:236:PHE:C	1:C:237:GLU:HG3	2.36	0.50
1:B:1:MSE:HE3	1:B:3:VAL:CG2	2.41	0.50
1:B:21:HIS:CE1	1:B:247:TYR:OH	2.61	0.50
1:B:35:ASN:C	1:B:35:ASN:HD22	2.20	0.50
1:B:70:HIS:CE1	1:B:72:GLU:HB2	2.46	0.50
1:B:153:MSE:CE	1:C:201:ILE:HD12	2.42	0.49
1:A:7:GLY:O	1:A:8:ASP:C	2.55	0.49
1:C:186:LEU:HD13	1:C:244:ILE:HB	1.94	0.49
1:A:43:LYS:NZ	1:A:69:ASP:HB2	2.27	0.49
1:A:51:ARG:O	1:A:55:GLU:HG3	2.12	0.49
1:C:1:MSE:HE1	1:C:107:LEU:HB2	1.93	0.49
1:D:172:HIS:HB2	1:D:201:ILE:CD1	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:63:LEU:HD11	1:C:82:VAL:HG13	1.94	0.49
1:C:240:ARG:CD	1:C:241:PRO:HD2	2.41	0.49
1:C:236:PHE:O	1:C:237:GLU:HG3	2.12	0.49
1:D:7:GLY:O	1:D:194:THR:HG23	2.12	0.49
1:A:10:MSE:O	1:A:39:ALA:HA	2.13	0.49
1:B:153:MSE:HE3	1:C:201:ILE:HD11	1.92	0.49
1:C:71:LYS:HG3	2:C:305:HOH:O	2.12	0.49
1:A:31:LEU:HD13	1:A:102:VAL:HG21	1.94	0.49
1:A:50:TYR:OH	1:A:81:PRO:HD2	2.13	0.49
1:B:196:THR:HG22	2:B:277:HOH:O	2.12	0.49
1:D:12:GLU:HB2	1:D:13:PRO:HD3	1.94	0.49
1:A:21:HIS:CE1	1:A:247:TYR:OH	2.64	0.48
1:C:63:LEU:HD23	1:C:67:ALA:CB	2.37	0.48
1:C:63:LEU:CD1	1:C:63:LEU:N	2.76	0.48
1:B:224:GLN:HB3	2:B:266:HOH:O	2.14	0.48
1:A:43:LYS:CE	1:A:69:ASP:HB2	2.42	0.48
1:A:99:ARG:NH1	1:A:99:ARG:HB2	2.28	0.48
1:C:10:MSE:HB3	1:C:11:ALA:H	1.47	0.48
1:C:160:ASP:CG	1:C:183:LYS:HB2	2.39	0.48
1:C:244:ILE:O	1:C:244:ILE:HG23	2.12	0.48
1:D:65:ASN:H	1:D:65:ASN:ND2	2.10	0.48
1:C:47:ARG:NE	1:C:51:ARG:HH12	2.12	0.48
1:C:208:PHE:O	1:C:211:ARG:HB3	2.13	0.48
1:B:127:ALA:HA	1:B:130:ARG:NH2	2.29	0.48
1:B:131:LEU:C	1:B:131:LEU:HD13	2.38	0.48
1:B:153:MSE:HE3	1:C:201:ILE:HD12	1.94	0.48
1:C:1:MSE:CE	1:C:107:LEU:HD22	2.39	0.48
1:B:169:THR:O	1:B:170:HIS:HB3	2.14	0.48
1:D:139:TYR:CD1	1:D:236:PHE:HB3	2.48	0.48
1:C:7:GLY:HA3	1:C:192:GLY:O	2.13	0.47
1:B:108:LEU:HD21	1:B:131:LEU:CD1	2.45	0.47
1:C:43:LYS:HB3	1:C:69:ASP:OD2	2.13	0.47
1:A:169:THR:O	1:A:170:HIS:HB3	2.15	0.47
1:A:1:MSE:HE3	1:A:236:PHE:HD2	1.78	0.47
1:B:51:ARG:O	1:B:55:GLU:HG3	2.13	0.47
1:A:55:GLU:HB3	2:A:367:HOH:O	2.15	0.47
1:A:11:ALA:HB1	1:A:12:GLU:OE2	2.15	0.47
1:D:10:MSE:O	1:D:39:ALA:HA	2.14	0.47
1:D:114:GLY:C	1:D:118:MSE:HE3	2.40	0.47
1:D:43:LYS:HE3	1:D:69:ASP:CG	2.39	0.47
1:A:16:ARG:HD2	1:A:251:GLU:OE1	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:12:GLU:CD	1:C:12:GLU:H	2.22	0.47
1:C:22:LEU:O	1:C:25:ILE:HD13	2.14	0.47
1:D:6:ILE:HB	1:D:34:ALA:CB	2.45	0.47
1:A:35:ASN:C	1:A:35:ASN:ND2	2.73	0.47
1:A:160:ASP:HA	1:A:185:THR:OG1	2.15	0.47
1:C:22:LEU:HD12	1:C:25:ILE:HD11	1.97	0.47
1:A:30:ASP:CB	1:A:102:VAL:HG11	2.42	0.46
1:B:35:ASN:ND2	1:B:37:GLU:H	2.13	0.46
1:A:113:MSE:CE	1:A:145:HIS:HB3	2.36	0.46
1:C:84:ARG:O	1:C:96:GLY:HA2	2.15	0.46
1:D:1:MSE:HE1	1:D:107:LEU:HB2	1.95	0.46
1:A:54:ARG:HD2	1:A:58:VAL:O	2.14	0.46
1:C:71:LYS:C	1:C:71:LYS:CD	2.85	0.46
1:D:13:PRO:HB2	2:D:254:HOH:O	2.14	0.46
1:B:127:ALA:HA	1:B:130:ARG:HH21	1.80	0.46
1:C:70:HIS:O	1:C:73:VAL:HG22	2.14	0.46
1:B:200:ILE:HG13	1:B:208:PHE:CB	2.46	0.46
1:C:63:LEU:HB3	1:C:67:ALA:HB2	1.96	0.46
1:D:156:ALA:HB1	1:D:181:LEU:HD11	1.97	0.46
1:C:116:ILE:HG23	1:C:151:GLU:OE1	2.16	0.46
1:C:160:ASP:O	1:C:183:LYS:HD2	2.15	0.46
1:A:181:LEU:HB3	1:A:182:PRO:HD2	1.98	0.46
1:C:63:LEU:HD13	1:C:83:VAL:O	2.16	0.46
1:D:43:LYS:HZ2	1:D:66:HIS:HB3	1.81	0.46
1:C:77:LEU:HG	1:C:94:GLY:HA3	1.98	0.46
1:A:85:PRO:CG	1:A:113:MSE:HE3	2.46	0.45
1:B:12:GLU:CD	1:B:12:GLU:N	2.70	0.45
1:A:43:LYS:HZ1	1:A:69:ASP:HB2	1.82	0.45
1:A:177:ASP:O	1:A:179:THR:HG23	2.16	0.45
1:A:200:ILE:O	1:A:200:ILE:CG1	2.64	0.45
1:C:63:LEU:HD12	1:C:63:LEU:N	2.31	0.45
1:D:208:PHE:O	1:D:211:ARG:HB3	2.17	0.45
1:A:209:LEU:HD11	1:D:158:TYR:HB2	1.98	0.45
1:B:181:LEU:HB3	1:B:182:PRO:HD2	1.98	0.45
1:D:85:PRO:HG2	1:D:113:MSE:HE3	1.98	0.45
1:C:47:ARG:HH11	1:C:47:ARG:CB	2.29	0.45
1:C:118:MSE:HE2	2:C:253:HOH:O	2.16	0.45
1:D:43:LYS:HZ2	1:D:66:HIS:CG	2.34	0.45
1:A:1:MSE:CE	1:A:236:PHE:HD2	2.30	0.45
1:B:156:ALA:HB1	1:B:181:LEU:HD11	1.99	0.45
1:C:194:THR:HB	1:C:229:PHE:HA	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:4:LEU:HB2	1:D:29:TYR:CD2	2.52	0.45
1:A:116:ILE:HG22	1:A:151:GLU:OE1	2.17	0.45
1:A:186:LEU:HG	1:A:241:PRO:HB2	1.98	0.45
1:C:16:ARG:HG2	1:C:20:LEU:HD22	1.99	0.45
1:B:200:ILE:HD13	1:B:203:GLY:O	2.17	0.45
1:C:45:LEU:O	1:C:73:VAL:HG12	2.17	0.45
1:B:200:ILE:N	1:B:200:ILE:CD1	2.80	0.45
1:C:1:MSE:HE1	1:C:107:LEU:CD2	2.41	0.45
1:D:168:GLY:O	1:D:190:ASP:HA	2.16	0.44
1:B:2:ARG:HB2	1:B:29:TYR:HA	1.99	0.44
1:D:47:ARG:NE	1:D:51:ARG:HH12	2.14	0.44
1:D:220:PHE:CD1	1:D:220:PHE:C	2.95	0.44
1:C:31:LEU:HD22	1:C:102:VAL:HG21	1.99	0.44
1:D:37:GLU:HG2	1:D:66:HIS:HB2	1.99	0.44
1:B:37:GLU:CG	1:B:66:HIS:HB2	2.45	0.44
1:C:13:PRO:HB2	2:C:254:HOH:O	2.16	0.44
1:A:12:GLU:N	1:A:13:PRO:CD	2.81	0.44
1:A:187:TYR:CG	1:A:188:GLN:N	2.85	0.44
1:B:1:MSE:CE	1:B:236:PHE:HD2	2.30	0.44
1:B:200:ILE:O	1:B:200:ILE:CG1	2.65	0.44
1:A:1:MSE:HA	1:A:30:ASP:OD2	2.18	0.44
1:B:43:LYS:HE3	1:B:69:ASP:OD1	2.18	0.44
1:D:35:ASN:C	1:D:35:ASN:HD22	2.26	0.44
1:D:85:PRO:CG	1:D:113:MSE:HE3	2.48	0.44
1:D:145:HIS:CD2	1:D:169:THR:OG1	2.69	0.44
1:B:150:SER:CA	1:C:201:ILE:HD13	2.39	0.44
1:D:113:MSE:HE2	1:D:145:HIS:ND1	2.32	0.44
1:D:200:ILE:N	1:D:200:ILE:CD1	2.69	0.44
1:A:108:LEU:HD21	1:A:131:LEU:CD1	2.48	0.43
1:C:200:ILE:HD13	1:C:200:ILE:H	1.81	0.43
1:B:31:LEU:HD13	1:B:102:VAL:CG2	2.46	0.43
1:B:64:GLY:HA2	1:B:113:MSE:HE1	2.00	0.43
1:A:85:PRO:HG3	1:A:113:MSE:HE3	2.00	0.43
1:B:114:GLY:O	1:B:118:MSE:HE3	2.18	0.43
1:D:160:ASP:OD1	1:D:183:LYS:HB2	2.18	0.43
1:B:11:ALA:HB1	1:B:12:GLU:OE2	2.18	0.43
1:D:31:LEU:CD2	1:D:102:VAL:HG21	2.49	0.43
1:A:168:GLY:O	1:A:190:ASP:HA	2.18	0.43
1:B:13:PRO:HD2	2:B:254:HOH:O	2.17	0.43
1:A:64:GLY:HA2	1:A:113:MSE:HE1	1.99	0.43
1:B:200:ILE:HD11	1:B:208:PHE:HB2	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:252:PRO:HG2	2:B:330:HOH:O	2.17	0.43
1:C:169:THR:O	1:C:170:HIS:CB	2.67	0.43
1:C:186:LEU:HG	1:C:241:PRO:HB2	2.00	0.43
1:B:127:ALA:CA	1:B:130:ARG:HH21	2.31	0.43
1:C:102:VAL:O	1:C:102:VAL:HG12	2.19	0.43
1:D:43:LYS:HZ2	1:D:66:HIS:CB	2.32	0.43
1:D:200:ILE:HD11	1:D:204:GLU:C	2.44	0.43
1:B:65:ASN:HD21	1:B:170:HIS:CE1	2.37	0.43
1:A:12:GLU:CD	1:A:12:GLU:N	2.73	0.42
1:A:75:ALA:O	1:A:78:GLU:HG2	2.19	0.42
1:B:1:MSE:HE1	1:B:107:LEU:HD13	2.00	0.42
1:C:31:LEU:CD2	1:C:102:VAL:HG21	2.49	0.42
1:C:88:TYR:HD1	1:C:89:PRO:HD2	1.83	0.42
1:D:10:MSE:HE2	1:D:38:ASN:HD22	1.84	0.42
1:A:186:LEU:HD13	1:A:244:ILE:HB	2.01	0.42
1:C:6:ILE:HB	1:C:34:ALA:CB	2.49	0.42
1:D:6:ILE:HB	1:D:34:ALA:HB2	2.01	0.42
1:D:45:LEU:O	1:D:73:VAL:HG12	2.20	0.42
1:D:99:ARG:NH2	1:D:136:LYS:O	2.44	0.42
1:A:63:LEU:HD21	1:A:82:VAL:CG1	2.48	0.42
1:D:77:LEU:HG	1:D:94:GLY:CA	2.47	0.42
1:B:35:ASN:C	1:B:35:ASN:ND2	2.78	0.42
1:B:99:ARG:NH1	1:B:99:ARG:HB2	2.34	0.42
1:A:200:ILE:HD11	1:A:208:PHE:CD2	2.54	0.42
1:D:5:PHE:HA	1:D:33:ILE:O	2.19	0.42
1:A:200:ILE:HD11	1:A:208:PHE:CG	2.55	0.42
1:B:77:LEU:HG	1:B:94:GLY:HA3	2.02	0.42
1:B:158:TYR:CE1	1:C:213:LEU:HD11	2.54	0.42
1:A:12:GLU:OE2	1:A:41:ARG:HA	2.20	0.42
1:A:145:HIS:CD2	1:A:170:HIS:HB2	2.55	0.42
1:B:88:TYR:HD1	1:B:89:PRO:HD2	1.85	0.42
1:C:116:ILE:O	1:C:117:PHE:HB2	2.19	0.42
1:D:11:ALA:HB1	1:D:12:GLU:OE1	2.20	0.42
1:D:118:MSE:HE1	1:D:146:ALA:CA	2.49	0.42
1:A:171:THR:O	1:A:172:HIS:HB2	2.19	0.42
1:B:45:LEU:HD13	1:B:76:LEU:CD1	2.50	0.42
1:B:209:LEU:HD12	1:B:209:LEU:HA	1.91	0.42
1:C:37:GLU:HG2	1:C:66:HIS:HB2	2.01	0.42
1:D:70:HIS:O	1:D:73:VAL:HG22	2.20	0.42
1:A:37:GLU:CG	1:A:66:HIS:HB2	2.50	0.41
1:A:45:LEU:HD11	1:A:50:TYR:HB2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:63:LEU:HD22	1:A:83:VAL:O	2.20	0.41
1:C:4:LEU:HB2	1:C:29:TYR:CD2	2.56	0.41
1:C:145:HIS:CD2	1:C:170:HIS:HB2	2.54	0.41
1:D:14:GLY:N	2:D:305:HOH:O	2.16	0.41
1:D:54:ARG:HD2	1:D:58:VAL:O	2.21	0.41
1:C:13:PRO:HB3	1:C:227:ALA:HB2	2.03	0.41
1:D:200:ILE:HD13	1:D:203:GLY:O	2.21	0.41
1:A:200:ILE:N	1:A:200:ILE:CD1	2.84	0.41
1:B:63:LEU:HB3	1:B:64:GLY:H	1.60	0.41
1:A:25:ILE:H	1:A:25:ILE:HD13	1.86	0.41
1:A:162:ARG:HG3	2:A:316:HOH:O	2.19	0.41
1:B:7:GLY:O	1:B:8:ASP:C	2.63	0.41
1:D:172:HIS:HB2	1:D:201:ILE:HD12	2.01	0.41
1:B:62:SER:C	1:B:63:LEU:HD13	2.46	0.41
1:D:64:GLY:CA	1:D:113:MSE:HE1	2.43	0.41
1:B:30:ASP:CB	1:B:102:VAL:HG11	2.42	0.41
1:B:236:PHE:C	1:B:237:GLU:HG3	2.46	0.41
1:C:35:ASN:C	1:C:35:ASN:ND2	2.79	0.41
1:C:108:LEU:HD21	1:C:131:LEU:HD11	2.02	0.41
1:B:128:LEU:HD12	1:B:128:LEU:HA	1.97	0.40
1:B:132:LEU:HD12	1:B:132:LEU:HA	1.88	0.40
1:B:171:THR:HG22	1:B:173:VAL:H	1.87	0.40
1:A:13:PRO:HD2	2:A:253:HOH:O	2.20	0.40
1:A:30:ASP:O	1:A:31:LEU:HD13	2.21	0.40
1:A:211:ARG:HD2	1:A:211:ARG:O	2.22	0.40
1:B:241:PRO:HD3	2:B:337:HOH:O	2.20	0.40
1:D:35:ASN:ND2	1:D:35:ASN:C	2.79	0.40
1:B:10:MSE:HB3	1:B:11:ALA:H	1.45	0.40
1:A:43:LYS:HZ2	1:A:66:HIS:CB	2.34	0.40
1:A:130:ARG:O	1:A:134:GLU:HG2	2.21	0.40
1:B:209:LEU:HD11	1:C:158:TYR:HB2	2.03	0.40
1:C:118:MSE:CE	1:C:146:ALA:HA	2.49	0.40
1:A:88:TYR:HD1	1:A:89:PRO:HD2	1.86	0.40
1:B:205:VAL:O	1:B:209:LEU:HB2	2.21	0.40
1:D:43:LYS:HB3	1:D:69:ASP:OD2	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	250/252 (99%)	238 (95%)	8 (3%)	4 (2%)	7	14
1	B	250/252 (99%)	240 (96%)	7 (3%)	3 (1%)	10	20
1	C	250/252 (99%)	232 (93%)	15 (6%)	3 (1%)	10	20
1	D	250/252 (99%)	236 (94%)	12 (5%)	2 (1%)	16	31
All	All	1000/1008 (99%)	946 (95%)	42 (4%)	12 (1%)	10	20

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	10	MSE
1	B	10	MSE
1	C	10	MSE
1	D	10	MSE
1	A	8	ASP
1	B	8	ASP
1	C	8	ASP
1	D	8	ASP
1	A	199	SER
1	B	201	ILE
1	C	102	VAL
1	A	201	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	165/198 (83%)	148 (90%)	17 (10%)	7	14
1	B	52/198 (26%)	47 (90%)	5 (10%)	8	17
1	C	88/198 (44%)	85 (97%)	3 (3%)	32	60
1	D	139/198 (70%)	126 (91%)	13 (9%)	8	18
All	All	444/792 (56%)	406 (91%)	38 (9%)	10	21

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	LEU
1	A	10	MSE
1	A	15	LEU
1	A	25	ILE
1	A	31	LEU
1	A	35	ASN
1	A	45	LEU
1	A	63	LEU
1	A	77	LEU
1	A	113	MSE
1	A	116	ILE
1	A	118	MSE
1	A	128	LEU
1	A	132	LEU
1	A	159	LEU
1	A	171	THR
1	A	200	ILE
1	B	112	VAL
1	B	113	MSE
1	B	116	ILE
1	B	118	MSE
1	B	132	LEU
1	C	171	THR
1	C	194	THR
1	C	200	ILE
1	D	4	LEU
1	D	10	MSE
1	D	35	ASN
1	D	45	LEU
1	D	47	ARG
1	D	51	ARG
1	D	63	LEU
1	D	65	ASN

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Mol	Chain	Res	Type
1	D	72	GLU
1	D	102	VAL
1	D	112	VAL
1	D	113	MSE
1	D	116	ILE

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

Mol	Chain	Res	Type
1	A	21	HIS
1	A	35	ASN
1	A	66	HIS
1	A	111	GLN
1	A	145	HIS
1	A	170	HIS
1	A	188	GLN
1	B	111	GLN
1	B	145	HIS
1	C	145	HIS
1	C	157	HIS
1	C	170	HIS
1	C	218	GLN
1	D	21	HIS
1	D	35	ASN
1	D	38	ASN
1	D	65	ASN
1	D	218	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides 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

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.