



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:07 PM UTC

PDB ID : 1LTL / pdb_00001ltl
Title : THE DODECAMER STRUCTURE OF MCM FROM ARCHAEAL M. THERMOAUTOTROPHICUM
Authors : Fletcher, R.J.; Bishop, B.E.; Leon, R.P.; Sclafani, R.A.; Ogata, C.M.; Chen, X.S.
Deposited on : 2002-05-20
Resolution : 3.00 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

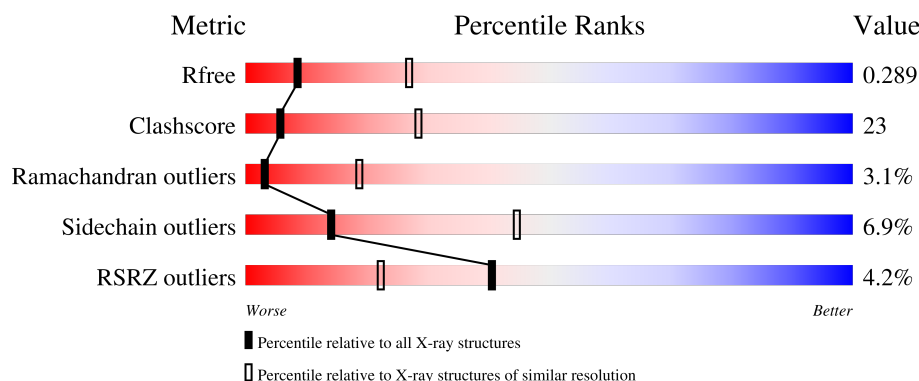
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	279	
1	B	279	
1	C	279	
1	D	279	
1	E	279	

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Mol	Chain	Length	Quality of chain
1	F	279	 53% 27% 5% 14%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11633 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 DNA replication initiator (Cdc21/Cdc54).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	239	Total	C	N	O	S	0	0	0
			1941	1222	341	371	7			
1	B	239	Total	C	N	O	S	0	0	0
			1941	1222	341	371	7			
1	C	239	Total	C	N	O	S	0	0	0
			1926	1209	341	369	7			
1	D	239	Total	C	N	O	S	0	0	0
			1924	1211	341	365	7			
1	E	242	Total	C	N	O	S	0	0	0
			1956	1231	344	374	7			
1	F	239	Total	C	N	O	S	0	0	0
			1939	1220	341	371	7			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	GLY	-	cloning artifact	UNP O27798
A	-3	SER	-	cloning artifact	UNP O27798
A	-2	GLY	-	cloning artifact	UNP O27798
A	-1	SER	-	cloning artifact	UNP O27798
A	0	ARG	-	cloning artifact	UNP O27798
B	-4	GLY	-	cloning artifact	UNP O27798
B	-3	SER	-	cloning artifact	UNP O27798
B	-2	GLY	-	cloning artifact	UNP O27798
B	-1	SER	-	cloning artifact	UNP O27798
B	0	ARG	-	cloning artifact	UNP O27798
C	-4	GLY	-	cloning artifact	UNP O27798
C	-3	SER	-	cloning artifact	UNP O27798
C	-2	GLY	-	cloning artifact	UNP O27798
C	-1	SER	-	cloning artifact	UNP O27798
C	0	ARG	-	cloning artifact	UNP O27798
D	-4	GLY	-	cloning artifact	UNP O27798
D	-3	SER	-	cloning artifact	UNP O27798

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-2	GLY	-	cloning artifact	UNP O27798
D	-1	SER	-	cloning artifact	UNP O27798
D	0	ARG	-	cloning artifact	UNP O27798
E	-4	GLY	-	cloning artifact	UNP O27798
E	-3	SER	-	cloning artifact	UNP O27798
E	-2	GLY	-	cloning artifact	UNP O27798
E	-1	SER	-	cloning artifact	UNP O27798
E	0	ARG	-	cloning artifact	UNP O27798
F	-4	GLY	-	cloning artifact	UNP O27798
F	-3	SER	-	cloning artifact	UNP O27798
F	-2	GLY	-	cloning artifact	UNP O27798
F	-1	SER	-	cloning artifact	UNP O27798
F	0	ARG	-	cloning artifact	UNP O27798

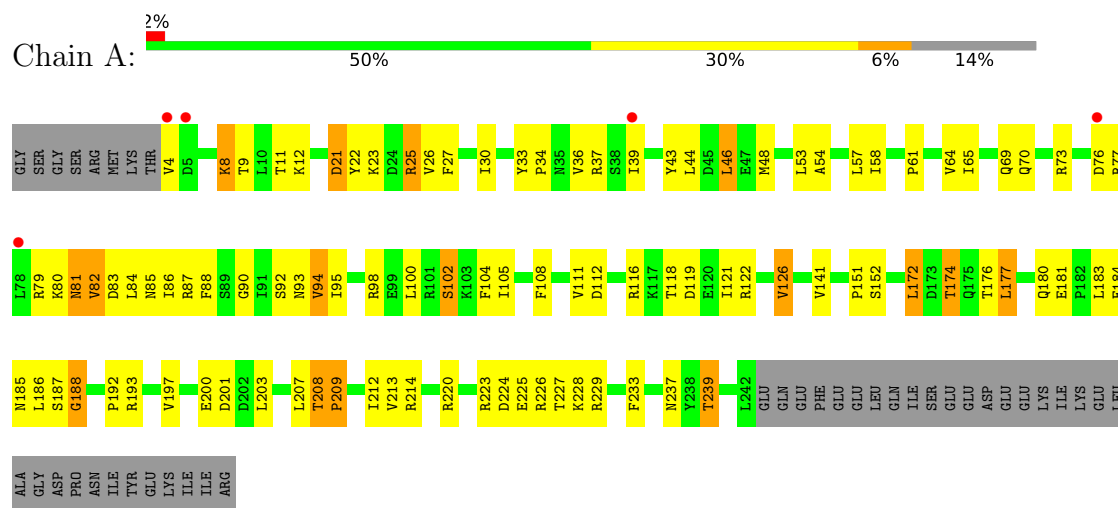
- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		
2	C	1	Total	Zn	0	0
			1	1		
2	D	1	Total	Zn	0	0
			1	1		
2	E	1	Total	Zn	0	0
			1	1		
2	F	1	Total	Zn	0	0
			1	1		

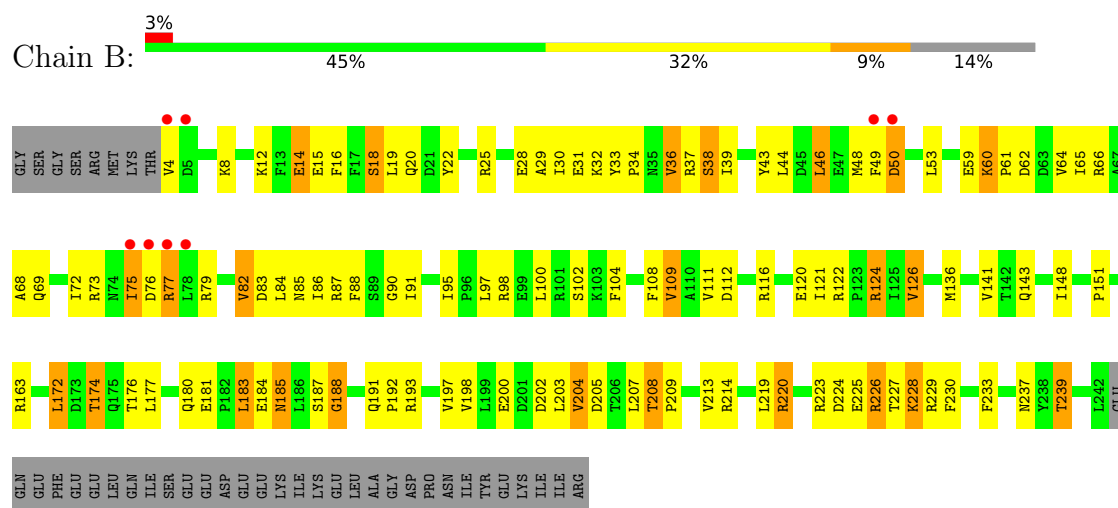
3 Residue-property plots [i](#)

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: DNA replication initiator (Cdc21/Cdc54)

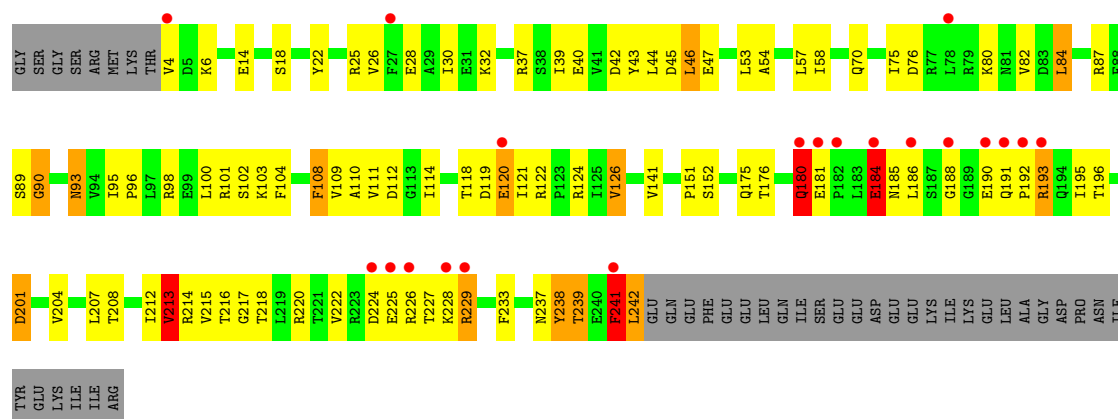


- Molecule 1: DNA replication initiator (Cdc21/Cdc54)

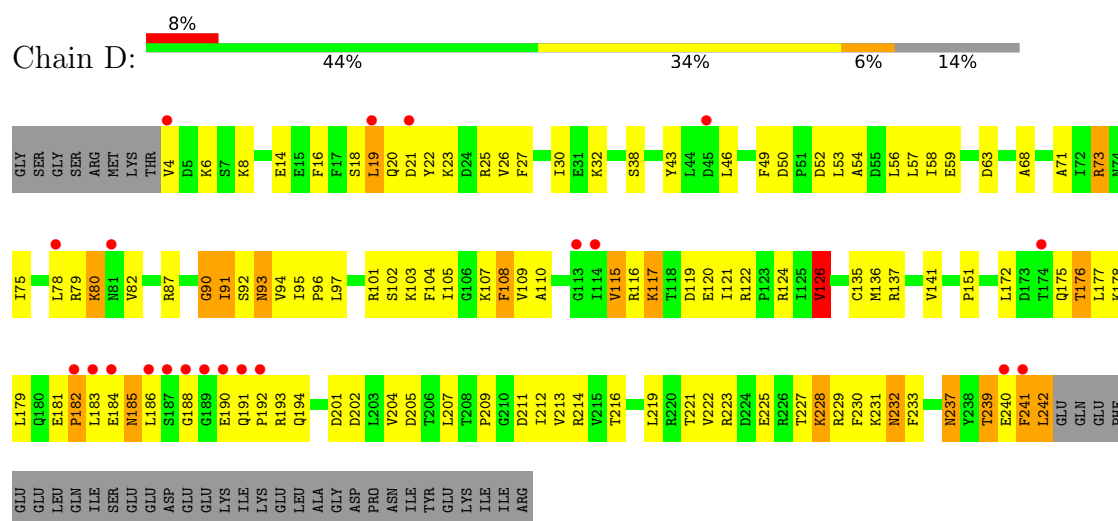


- Molecule 1: DNA replication initiator (Cdc21/Cdc54)

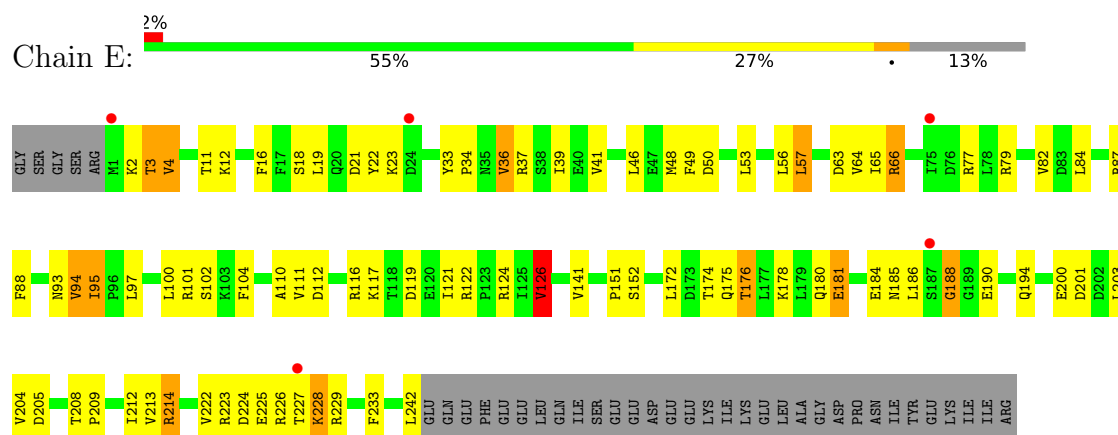




• Molecule 1: DNA replication initiator (Cdc21/Cdc54)

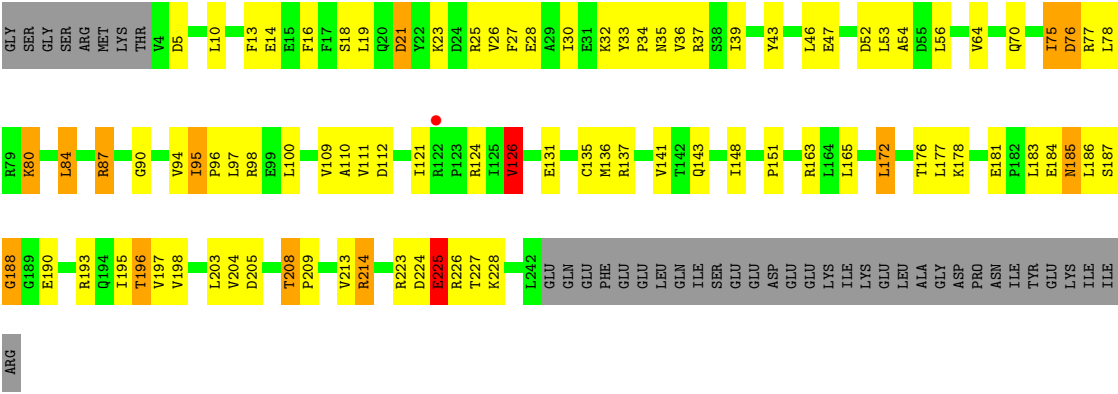


• Molecule 1: DNA replication initiator (Cdc21/Cdc54)



• Molecule 1: DNA replication initiator (Cdc21/Cdc54)





4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	192.32Å 192.32Å 365.54Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 3.00 20.00 – 3.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-3.00) 95.8 (20.00-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.48 (at 2.91Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.245 , 0.295 0.251 , 0.289	Depositor DCC
R_{free} test set	4623 reflections (8.19%)	wwPDB-VP
Wilson B-factor (Å ²)	185.9	Xtriage
Anisotropy	0.037	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 44.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	11633	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.52% 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: ZN

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	0.63	0/1970	0.99	5/2656 (0.2%)
1	B	0.70	0/1970	1.03	5/2656 (0.2%)
1	C	0.88	3/1954 (0.2%)	2.09	15/2635 (0.6%)
1	D	1.64	7/1953 (0.4%)	1.40	18/2633 (0.7%)
1	E	0.67	0/1985	1.00	9/2677 (0.3%)
1	F	0.63	0/1968	0.94	3/2653 (0.1%)
All	All	0.93	10/11800 (0.1%)	1.31	55/15910 (0.3%)

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	C	0	3
1	D	0	1
All	All	0	4

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	237	ASN	C-N	-51.81	0.53	1.33
1	D	182	PRO	C-N	31.21	1.74	1.33
1	D	241	PHE	C-N	25.33	1.68	1.33
1	C	241	PHE	C-N	-23.61	1.00	1.33
1	C	242	LEU	C-O	14.21	1.51	1.23
1	C	193	ARG	C-N	-13.70	1.16	1.33
1	D	239	THR	C-N	-13.37	1.13	1.33
1	D	242	LEU	C-O	12.70	1.49	1.23
1	D	240	GLU	CB-CG	-6.47	1.33	1.52
1	D	192	PRO	C-N	5.64	1.41	1.33

All (55) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	180	GLN	O-C-N	-75.22	38.27	123.27
1	C	241	PHE	O-C-N	-29.97	86.16	123.02
1	D	237	ASN	O-C-N	-23.51	91.33	122.59
1	C	241	PHE	CA-C-N	22.65	162.46	121.70
1	C	241	PHE	C-N-CA	22.65	162.46	121.70
1	C	242	LEU	CA-C-O	21.28	156.98	120.80
1	D	241	PHE	O-C-N	-20.11	94.80	121.95
1	D	237	ASN	CA-C-N	18.16	152.14	122.73
1	D	237	ASN	C-N-CA	18.16	152.14	122.73
1	C	238	TYR	O-C-N	17.68	146.10	122.59
1	D	192	PRO	CA-C-N	17.20	146.48	121.02
1	D	192	PRO	C-N-CA	17.20	146.48	121.02
1	C	238	TYR	CA-C-N	-14.69	101.64	122.93
1	C	238	TYR	C-N-CA	-14.69	101.64	122.93
1	C	180	GLN	CA-C-N	-13.45	82.77	120.97
1	C	180	GLN	C-N-CA	-13.45	82.77	120.97
1	D	239	THR	O-C-N	-10.84	110.63	123.31
1	D	192	PRO	O-C-N	-10.56	108.39	122.64
1	D	241	PHE	CA-C-N	-8.96	105.58	121.70
1	D	241	PHE	C-N-CA	-8.96	105.58	121.70
1	C	213	VAL	N-CA-C	7.59	117.79	108.53
1	B	213	VAL	N-CA-C	7.58	117.77	108.53
1	D	239	THR	CA-C-N	7.46	136.12	122.02
1	D	239	THR	C-N-CA	7.46	136.12	122.02
1	C	241	PHE	CA-C-O	-7.02	112.35	120.49
1	D	240	GLU	CB-CA-C	6.74	126.60	111.78
1	D	241	PHE	CA-C-O	-6.73	113.40	121.34
1	D	38	SER	N-CA-C	6.72	119.94	109.52
1	E	213	VAL	N-CA-C	6.62	118.46	108.80
1	D	242	LEU	CA-C-O	6.62	132.05	120.80
1	A	213	VAL	N-CA-C	6.53	118.33	108.80
1	A	102	SER	N-CA-C	6.47	118.33	111.28
1	E	181	GLU	CA-C-N	6.15	126.33	119.32
1	E	181	GLU	C-N-CA	6.15	126.33	119.32
1	B	203	LEU	N-CA-C	-6.13	105.56	113.16
1	E	102	SER	N-CA-C	6.07	117.97	111.36
1	F	213	VAL	N-CA-C	6.06	117.65	108.80
1	C	18	SER	N-CA-C	-5.92	106.67	114.31
1	E	21	ASP	N-CA-C	5.89	117.70	111.28
1	B	102	SER	N-CA-C	5.88	118.45	111.33
1	A	201	ASP	CB-CA-C	-5.83	109.83	116.54
1	A	122	ARG	CA-C-N	-5.74	114.65	120.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	122	ARG	C-N-CA	-5.74	114.65	120.85
1	F	225	GLU	N-CA-C	5.71	117.58	111.36
1	E	201	ASP	CB-CA-C	-5.55	109.68	117.23
1	C	239	THR	CA-CB-CG2	-5.52	101.11	110.50
1	D	240	GLU	O-C-N	5.28	128.75	122.20
1	B	38	SER	N-CA-C	5.26	118.31	109.95
1	F	126	VAL	CB-CA-C	-5.26	105.50	111.55
1	E	126	VAL	CB-CA-C	-5.25	105.51	111.55
1	E	82	VAL	N-CA-C	5.14	115.26	107.80
1	D	126	VAL	CB-CA-C	-5.11	105.67	111.55
1	E	88	PHE	N-CA-C	5.08	117.89	109.46
1	B	18	SER	N-CA-C	-5.07	108.12	114.56
1	C	126	VAL	CB-CA-C	-5.04	105.76	111.55

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	180	GLN	Mainchain
1	C	241	PHE	Mainchain,Peptide
1	D	237	ASN	Mainchain

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	1941	0	1950	92	0
1	B	1941	0	1950	103	0
1	C	1926	0	1921	117	0
1	D	1924	0	1916	107	0
1	E	1956	0	1959	75	0
1	F	1939	0	1943	71	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	F	1	0	0	0	0
All	All	11633	0	11639	541	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 (541) 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:241:PHE:C	1:D:242:LEU:N	1.68	1.48
1:D:182:PRO:C	1:D:183:LEU:N	1.74	1.42
1:C:98:ARG:HH21	1:C:193:ARG:CZ	1.35	1.38
1:C:227:THR:CG2	1:C:229:ARG:HG3	1.62	1.29
1:C:98:ARG:NE	1:C:193:ARG:HD2	1.43	1.29
1:C:98:ARG:HE	1:C:193:ARG:CD	1.49	1.25
1:C:227:THR:HG22	1:C:229:ARG:CG	1.72	1.18
1:C:225:GLU:O	1:C:226:ARG:HG2	1.41	1.18
1:D:207:LEU:HD11	1:D:239:THR:HG21	1.21	1.15
1:C:98:ARG:NH2	1:C:193:ARG:CZ	2.15	1.09
1:C:98:ARG:HH21	1:C:193:ARG:NE	1.50	1.08
1:A:30:ILE:HD13	1:A:76:ASP:HB3	1.44	0.96
1:B:46:LEU:HD21	1:B:53:LEU:HD23	1.48	0.93
1:C:114:ILE:H	1:C:180:GLN:HB3	1.34	0.93
1:D:126:VAL:HG11	1:D:172:LEU:HD12	1.56	0.87
1:F:27:PHE:HE2	1:F:77:ARG:H	1.22	0.87
1:C:14:GLU:HA	1:C:75:ILE:HD11	1.53	0.87
1:C:98:ARG:HD3	1:C:181:GLU:CD	2.00	0.86
1:E:46:LEU:HD21	1:E:53:LEU:HD23	1.55	0.86
1:D:241:PHE:C	1:D:242:LEU:CA	2.50	0.85
1:C:229:ARG:HD3	1:D:122:ARG:NH1	1.92	0.84
1:B:112:ASP:OD1	1:B:214:ARG:HG3	1.77	0.83
1:C:98:ARG:HE	1:C:193:ARG:HD2	0.68	0.83
1:B:126:VAL:HG11	1:B:172:LEU:HD13	1.59	0.82
1:A:177:LEU:CD2	1:A:197:VAL:HB	2.10	0.82
1:B:88:PHE:H	1:B:237:ASN:HD21	1.27	0.82
1:A:39:ILE:HD11	1:A:84:LEU:HD13	1.63	0.81
1:D:119:ASP:HB2	1:D:176:THR:HG23	1.62	0.80
1:D:241:PHE:O	1:D:242:LEU:N	2.14	0.79
1:C:98:ARG:NH2	1:C:193:ARG:NE	2.25	0.79
1:D:175:GLN:NE2	1:D:205:ASP:H	1.81	0.79
1:F:177:LEU:CD1	1:F:197:VAL:HB	2.13	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:175:GLN:HE22	1:D:205:ASP:H	1.33	0.77
1:C:58:ILE:HA	1:C:108:PHE:HB2	1.67	0.77
1:B:14:GLU:HG3	1:B:75:ILE:HD11	1.66	0.76
1:A:46:LEU:HD21	1:A:53:LEU:HD23	1.67	0.76
1:A:126:VAL:HG11	1:A:172:LEU:HD13	1.66	0.76
1:A:116:ARG:NH2	1:A:180:GLN:HE21	1.85	0.75
1:A:93:ASN:ND2	1:A:95:ILE:HD11	2.02	0.74
1:B:220:ARG:HG3	1:B:220:ARG:HH11	1.51	0.74
1:C:98:ARG:HH21	1:C:193:ARG:NH2	1.84	0.74
1:B:191:GLN:HE21	1:B:192:PRO:HD2	1.52	0.74
1:F:46:LEU:HD11	1:F:53:LEU:HD23	1.68	0.73
1:C:98:ARG:HD3	1:C:181:GLU:OE2	1.86	0.73
1:D:119:ASP:CB	1:D:176:THR:HG23	2.18	0.73
1:C:98:ARG:NE	1:C:193:ARG:CD	2.24	0.73
1:C:225:GLU:O	1:C:226:ARG:CG	2.31	0.73
1:F:177:LEU:HD11	1:F:197:VAL:HB	1.71	0.73
1:B:77:ARG:H	1:B:77:ARG:HD2	1.52	0.73
1:F:112:ASP:OD2	1:F:214:ARG:HG3	1.89	0.73
1:B:227:THR:HG23	1:B:229:ARG:HE	1.55	0.72
1:D:175:GLN:HE22	1:D:205:ASP:N	1.87	0.72
1:C:112:ASP:OD1	1:C:214:ARG:HG3	1.89	0.72
1:E:48:MET:HE2	1:E:48:MET:HA	1.73	0.71
1:A:207:LEU:HD21	1:A:239:THR:HG21	1.73	0.71
1:E:223:ARG:HH22	1:F:148:ILE:HD13	1.55	0.70
1:C:180:GLN:HG3	1:C:181:GLU:C	1.97	0.70
1:A:177:LEU:HD21	1:A:197:VAL:HB	1.73	0.70
1:B:86:ILE:HG13	1:B:86:ILE:O	1.89	0.70
1:A:177:LEU:HD23	1:A:197:VAL:HB	1.74	0.70
1:F:98:ARG:HD2	1:F:193:ARG:HD2	1.73	0.70
1:C:227:THR:CG2	1:C:229:ARG:CG	2.50	0.69
1:A:22:TYR:HD2	1:A:25:ARG:HG3	1.58	0.69
1:A:36:VAL:HG23	1:A:37:ARG:H	1.56	0.69
1:D:227:THR:HB	1:D:229:ARG:HD3	1.72	0.69
1:A:116:ARG:HH22	1:A:180:GLN:HE21	1.40	0.68
1:C:58:ILE:HD13	1:C:93:ASN:HD22	1.58	0.68
1:D:95:ILE:HD13	1:D:109:VAL:HG13	1.76	0.68
1:F:95:ILE:N	1:F:95:ILE:HD12	2.08	0.68
1:B:116:ARG:HD2	1:B:180:GLN:OE1	1.94	0.67
1:D:102:SER:HA	1:D:105:ILE:HG13	1.74	0.67
1:A:87:ARG:CZ	1:A:203:LEU:HD13	2.24	0.67
1:B:95:ILE:HD13	1:B:104:PHE:CZ	2.28	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:112:ASP:OD1	1:A:214:ARG:HG3	1.94	0.67
1:A:73:ARG:HH21	1:A:83:ASP:HA	1.59	0.67
1:D:18:SER:C	1:D:20:GLN:H	2.02	0.67
1:D:213:VAL:HG12	1:D:241:PHE:HA	1.76	0.67
1:A:98:ARG:HD2	1:A:193:ARG:HD2	1.78	0.66
1:B:30:ILE:HD13	1:B:76:ASP:OD1	1.96	0.66
1:C:213:VAL:CG1	1:C:241:PHE:HA	2.25	0.66
1:B:227:THR:CG2	1:B:229:ARG:HE	2.09	0.66
1:D:58:ILE:HA	1:D:108:PHE:HB2	1.77	0.66
1:B:220:ARG:HG3	1:B:220:ARG:NH1	2.08	0.66
1:C:152:SER:HB3	1:D:136:MET:CE	2.25	0.66
1:E:66:ARG:HG2	1:E:66:ARG:HH11	1.59	0.66
1:F:176:THR:HG22	1:F:198:VAL:HG22	1.77	0.65
1:A:100:LEU:HD11	1:A:111:VAL:HG21	1.79	0.65
1:A:39:ILE:CD1	1:A:84:LEU:HD13	2.26	0.65
1:D:95:ILE:HG23	1:D:96:PRO:HD2	1.78	0.64
1:C:180:GLN:HE22	1:C:192:PRO:HB2	1.62	0.64
1:F:39:ILE:HD11	1:F:84:LEU:HD13	1.78	0.64
1:C:220:ARG:HG3	1:C:220:ARG:HH11	1.62	0.64
1:B:220:ARG:HB2	1:B:233:PHE:CZ	2.33	0.64
1:F:126:VAL:HG11	1:F:172:LEU:HD13	1.79	0.63
1:F:223:ARG:NH1	1:F:228:LYS:O	2.31	0.63
1:B:22:TYR:CD2	1:B:25:ARG:HD2	2.34	0.63
1:C:212:ILE:O	1:C:242:LEU:CB	2.46	0.63
1:D:73:ARG:HH11	1:D:73:ARG:HB2	1.64	0.63
1:D:211:ASP:C	1:D:212:ILE:HD12	2.24	0.63
1:F:100:LEU:HD11	1:F:111:VAL:HG21	1.80	0.63
1:B:177:LEU:HD12	1:B:177:LEU:C	2.23	0.63
1:A:225:GLU:O	1:A:226:ARG:HB2	1.97	0.63
1:C:98:ARG:HD3	1:C:181:GLU:OE1	1.99	0.63
1:C:114:ILE:HG12	1:C:212:ILE:HD12	1.80	0.63
1:D:117:LYS:HE3	1:D:178:LYS:HG3	1.80	0.63
1:B:207:LEU:HD21	1:B:239:THR:HG21	1.79	0.62
1:C:141:VAL:HG21	1:C:151:PRO:HG3	1.80	0.62
1:D:141:VAL:HG21	1:D:151:PRO:HG3	1.80	0.62
1:D:232:ASN:HD22	1:D:232:ASN:N	1.96	0.62
1:A:220:ARG:HB2	1:A:233:PHE:CZ	2.34	0.62
1:C:98:ARG:CD	1:C:193:ARG:HD2	2.29	0.62
1:C:222:VAL:HG23	1:C:233:PHE:HE2	1.65	0.62
1:F:36:VAL:HG23	1:F:37:ARG:H	1.64	0.62
1:F:97:LEU:HD11	1:F:195:ILE:HG22	1.82	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:65:ILE:O	1:E:65:ILE:HG22	1.99	0.62
1:B:39:ILE:HD11	1:B:84:LEU:HD13	1.81	0.62
1:E:66:ARG:HG2	1:E:66:ARG:NH1	2.13	0.62
1:B:59:GLU:O	1:B:60:LYS:HD3	2.00	0.62
1:B:100:LEU:HD11	1:B:111:VAL:HG21	1.82	0.62
1:B:79:ARG:HD3	1:B:82:VAL:HG21	1.81	0.62
1:B:183:LEU:HD12	1:B:183:LEU:H	1.65	0.61
1:F:227:THR:O	1:F:228:LYS:HB2	2.00	0.61
1:C:152:SER:HB3	1:D:136:MET:HE2	1.81	0.61
1:F:141:VAL:HG21	1:F:151:PRO:HG3	1.82	0.61
1:C:58:ILE:HG21	1:C:93:ASN:ND2	2.16	0.61
1:E:227:THR:O	1:E:229:ARG:N	2.33	0.61
1:A:39:ILE:HG13	1:A:84:LEU:HD22	1.82	0.61
1:A:119:ASP:OD2	1:A:176:THR:HG22	2.00	0.61
1:A:141:VAL:HG21	1:A:151:PRO:HG3	1.82	0.61
1:E:223:ARG:HD3	1:E:228:LYS:O	2.00	0.61
1:A:30:ILE:HD13	1:A:76:ASP:CB	2.25	0.61
1:C:100:LEU:HD21	1:C:109:VAL:HG11	1.81	0.61
1:D:91:ILE:HD11	1:D:216:THR:HB	1.83	0.61
1:E:141:VAL:HG21	1:E:151:PRO:HG3	1.82	0.60
1:C:180:GLN:HE22	1:C:192:PRO:CB	2.14	0.60
1:C:207:LEU:HD21	1:C:239:THR:CG2	2.31	0.60
1:C:207:LEU:HD21	1:C:239:THR:HG21	1.84	0.60
1:D:80:LYS:HG2	1:D:82:VAL:HG13	1.83	0.60
1:F:121:ILE:HD11	1:F:205:ASP:HB2	1.82	0.60
1:A:93:ASN:HD21	1:A:95:ILE:HD11	1.66	0.60
1:D:116:ARG:HG2	1:D:117:LYS:HD3	1.83	0.60
1:E:119:ASP:OD2	1:E:176:THR:HG22	2.02	0.60
1:B:141:VAL:HG21	1:B:151:PRO:HG3	1.84	0.59
1:C:224:ASP:HB3	1:C:227:THR:HB	1.83	0.59
1:E:152:SER:HB3	1:F:136:MET:CE	2.32	0.59
1:C:84:LEU:HD12	1:C:84:LEU:H	1.67	0.59
1:B:39:ILE:HD12	1:B:72:ILE:HD13	1.85	0.59
1:E:39:ILE:HD11	1:E:84:LEU:HD13	1.85	0.59
1:C:229:ARG:CD	1:D:122:ARG:NH1	2.65	0.58
1:D:110:ALA:HA	1:D:216:THR:HA	1.85	0.58
1:C:180:GLN:NE2	1:C:181:GLU:HG3	2.17	0.58
1:C:213:VAL:HG13	1:C:241:PHE:HA	1.84	0.58
1:C:227:THR:HG22	1:C:229:ARG:HG3	0.75	0.58
1:E:178:LYS:HD2	1:E:194:GLN:NE2	2.19	0.58
1:A:227:THR:HG22	1:A:229:ARG:HG3	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:37:ARG:HH11	1:B:202:ASP:HB3	1.68	0.58
1:F:95:ILE:N	1:F:95:ILE:CD1	2.66	0.58
1:C:114:ILE:N	1:C:180:GLN:HB3	2.12	0.58
1:C:220:ARG:HG3	1:C:220:ARG:NH1	2.18	0.58
1:C:57:LEU:O	1:C:57:LEU:HG	2.04	0.58
1:D:26:VAL:O	1:D:30:ILE:HG12	2.04	0.58
1:D:231:LYS:C	1:D:232:ASN:HD22	2.12	0.58
1:F:177:LEU:HD12	1:F:197:VAL:HB	1.84	0.57
1:E:101:ARG:HG3	1:E:101:ARG:HH11	1.68	0.57
1:F:94:VAL:HG21	1:F:214:ARG:HH21	1.69	0.57
1:A:220:ARG:HH11	1:A:220:ARG:HG3	1.69	0.57
1:A:152:SER:HB3	1:B:136:MET:CE	2.34	0.57
1:C:87:ARG:HD2	1:C:237:ASN:O	2.03	0.57
1:E:224:ASP:OD1	1:E:227:THR:HB	2.04	0.57
1:F:208:THR:HG22	1:F:209:PRO:HD2	1.87	0.57
1:A:87:ARG:HA	1:A:237:ASN:HD21	1.70	0.57
1:C:216:THR:HB	1:C:238:TYR:HB3	1.87	0.57
1:B:49:PHE:C	1:B:50:ASP:CG	2.73	0.56
1:C:225:GLU:C	1:C:226:ARG:HG2	2.27	0.56
1:E:100:LEU:HD11	1:E:111:VAL:HG21	1.87	0.56
1:C:95:ILE:HG23	1:C:96:PRO:HD2	1.86	0.56
1:C:201:ASP:O	1:C:204:VAL:HG22	2.05	0.56
1:E:112:ASP:OD1	1:E:214:ARG:HG3	2.05	0.56
1:E:180:GLN:HG2	1:E:181:GLU:N	2.19	0.56
1:F:109:VAL:HG12	1:F:110:ALA:N	2.19	0.56
1:A:44:LEU:O	1:A:48:MET:HG2	2.05	0.56
1:B:97:LEU:C	1:B:97:LEU:HD23	2.30	0.56
1:B:208:THR:HG23	1:B:209:PRO:HD2	1.86	0.56
1:E:126:VAL:HG11	1:E:172:LEU:HD22	1.87	0.56
1:D:18:SER:HA	1:D:23:LYS:HB2	1.88	0.55
1:E:185:ASN:O	1:E:186:LEU:HD23	2.07	0.55
1:E:77:ARG:HG3	1:E:77:ARG:HH11	1.70	0.55
1:B:37:ARG:HA	1:B:84:LEU:HD23	1.87	0.55
1:B:37:ARG:NH1	1:B:202:ASP:HB3	2.22	0.55
1:B:176:THR:HG22	1:B:198:VAL:HG22	1.87	0.55
1:C:229:ARG:HH11	1:D:122:ARG:HH12	1.53	0.55
1:F:10:LEU:HD13	1:F:70:GLN:HB3	1.87	0.55
1:C:98:ARG:NH2	1:C:193:ARG:NH2	2.49	0.55
1:E:225:GLU:O	1:E:226:ARG:HB3	2.06	0.55
1:B:177:LEU:HD11	1:B:197:VAL:HB	1.88	0.55
1:A:36:VAL:O	1:A:37:ARG:HB2	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:88:PHE:N	1:B:237:ASN:HD21	2.02	0.54
1:E:126:VAL:CG1	1:E:172:LEU:HD22	2.37	0.54
1:B:46:LEU:CD2	1:B:53:LEU:HD23	2.32	0.54
1:A:26:VAL:O	1:A:30:ILE:HG13	2.06	0.54
1:C:224:ASP:OD1	1:C:225:GLU:N	2.40	0.54
1:D:97:LEU:O	1:D:97:LEU:HD23	2.08	0.54
1:A:152:SER:HA	1:B:163:ARG:NH2	2.23	0.54
1:D:115:VAL:CG1	1:D:211:ASP:H	2.20	0.54
1:A:27:PHE:HE2	1:A:77:ARG:HG3	1.73	0.54
1:C:186:LEU:HB3	1:C:190:GLU:HB2	1.90	0.54
1:A:37:ARG:O	1:A:84:LEU:HD23	2.07	0.54
1:A:121:ILE:HD12	1:A:121:ILE:H	1.72	0.54
1:E:152:SER:HB3	1:F:136:MET:HE2	1.90	0.54
1:C:213:VAL:HG12	1:C:241:PHE:CA	2.34	0.53
1:D:135:CYS:O	1:F:137:ARG:NH2	2.42	0.53
1:B:69:GLN:NE2	1:B:85:ASN:HA	2.24	0.53
1:E:208:THR:CG2	1:E:209:PRO:HD2	2.38	0.53
1:F:98:ARG:HD3	1:F:181:GLU:OE1	2.09	0.53
1:A:88:PHE:N	1:A:88:PHE:CD1	2.77	0.53
1:B:104:PHE:O	1:B:219:LEU:HB3	2.09	0.53
1:A:229:ARG:HD3	1:B:122:ARG:NE	2.24	0.53
1:D:22:TYR:O	1:D:26:VAL:HG23	2.09	0.53
1:D:214:ARG:HG3	1:D:214:ARG:HH11	1.74	0.53
1:A:212:ILE:N	1:A:212:ILE:HD12	2.24	0.53
1:C:228:LYS:NZ	1:D:225:GLU:HG3	2.23	0.53
1:F:21:ASP:HB3	1:F:25:ARG:HH22	1.73	0.53
1:A:121:ILE:HD12	1:A:121:ILE:N	2.23	0.53
1:D:227:THR:C	1:D:229:ARG:H	2.17	0.53
1:E:36:VAL:HG23	1:E:37:ARG:H	1.71	0.53
1:C:180:GLN:NE2	1:C:181:GLU:CG	2.72	0.53
1:D:50:ASP:HB3	1:D:53:LEU:HB3	1.91	0.53
1:D:186:LEU:HB3	1:D:190:GLU:HB2	1.90	0.52
1:A:223:ARG:HH22	1:B:148:ILE:HD13	1.75	0.52
1:A:58:ILE:HD12	1:A:58:ILE:N	2.23	0.52
1:C:229:ARG:NH1	1:D:122:ARG:HH12	2.07	0.52
1:B:16:PHE:CZ	1:B:22:TYR:CE1	2.98	0.52
1:B:33:TYR:CD1	1:B:34:PRO:HA	2.45	0.52
1:B:37:ARG:O	1:B:85:ASN:N	2.41	0.52
1:D:137:ARG:NH2	1:F:135:CYS:O	2.43	0.52
1:C:180:GLN:NE2	1:C:181:GLU:N	2.57	0.52
1:A:95:ILE:N	1:A:95:ILE:HD12	2.25	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:180:GLN:NE2	1:C:192:PRO:HB2	2.24	0.52
1:C:227:THR:C	1:C:229:ARG:H	2.16	0.52
1:F:225:GLU:OE1	1:F:225:GLU:HA	2.08	0.52
1:F:87:ARG:HE	1:F:203:LEU:HD11	1.75	0.51
1:E:97:LEU:HD23	1:E:97:LEU:O	2.10	0.51
1:E:57:LEU:HD23	1:E:64:VAL:CG1	2.40	0.51
1:A:94:VAL:C	1:A:95:ILE:HD12	2.35	0.51
1:A:220:ARG:HG3	1:A:220:ARG:NH1	2.25	0.51
1:C:109:VAL:HG12	1:C:110:ALA:N	2.26	0.51
1:D:95:ILE:HD12	1:D:95:ILE:N	2.26	0.51
1:C:89:SER:HA	1:C:238:TYR:CE2	2.46	0.51
1:A:82:VAL:HG12	1:A:84:LEU:HG	1.92	0.51
1:C:98:ARG:CD	1:C:181:GLU:OE2	2.56	0.51
1:C:100:LEU:HD11	1:C:111:VAL:HG21	1.93	0.51
1:F:178:LYS:CE	1:F:196:THR:HG23	2.41	0.51
1:A:226:ARG:H	1:A:228:LYS:NZ	2.09	0.51
1:A:36:VAL:HG23	1:A:37:ARG:N	2.22	0.51
1:A:21:ASP:OD2	1:A:21:ASP:N	2.43	0.51
1:B:121:ILE:HD11	1:B:205:ASP:HB2	1.92	0.51
1:C:226:ARG:O	1:C:228:LYS:HG3	2.10	0.51
1:B:223:ARG:NH1	1:B:228:LYS:O	2.42	0.50
1:C:152:SER:HB3	1:D:136:MET:HE1	1.93	0.50
1:C:118:THR:HG22	1:C:120:GLU:H	1.77	0.50
1:B:120:GLU:HG3	1:B:121:ILE:N	2.25	0.50
1:C:216:THR:HG22	1:C:217:GLY:N	2.26	0.50
1:F:30:ILE:HG21	1:F:76:ASP:CG	2.36	0.50
1:B:223:ARG:HA	1:B:230:PHE:HA	1.92	0.50
1:C:22:TYR:O	1:C:26:VAL:HG23	2.11	0.50
1:B:68:ALA:HB1	1:B:86:ILE:HD13	1.93	0.50
1:C:39:ILE:HD11	1:C:84:LEU:HD13	1.93	0.50
1:F:112:ASP:OD2	1:F:214:ARG:NH1	2.45	0.50
1:A:152:SER:HB3	1:B:136:MET:HE2	1.94	0.50
1:C:98:ARG:HE	1:C:193:ARG:NE	2.07	0.49
1:B:4:VAL:HG12	1:B:4:VAL:O	2.12	0.49
1:D:121:ILE:HD13	1:D:204:VAL:CG1	2.41	0.49
1:D:175:GLN:HE21	1:D:204:VAL:HG13	1.77	0.49
1:E:37:ARG:C	1:E:84:LEU:HD23	2.38	0.49
1:E:46:LEU:CD2	1:E:53:LEU:HD23	2.37	0.49
1:B:15:GLU:O	1:B:18:SER:HB2	2.13	0.49
1:D:116:ARG:NH1	1:D:194:GLN:HE22	2.09	0.49
1:E:227:THR:C	1:E:229:ARG:H	2.20	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:98:ARG:HD3	1:F:181:GLU:CD	2.37	0.49
1:B:177:LEU:CD1	1:B:197:VAL:HB	2.41	0.49
1:B:183:LEU:O	1:B:185:ASN:N	2.38	0.49
1:E:33:TYR:CD1	1:E:34:PRO:HA	2.47	0.49
1:A:9:THR:HG23	1:A:53:LEU:HD11	1.95	0.49
1:D:20:GLN:HA	1:D:23:LYS:HB3	1.94	0.49
1:C:121:ILE:HD12	1:C:175:GLN:NE2	2.28	0.49
1:D:181:GLU:HG3	1:D:193:ARG:HB2	1.95	0.49
1:D:228:LYS:HE3	1:D:230:PHE:CE1	2.48	0.49
1:E:228:LYS:N	1:E:228:LYS:HE2	2.28	0.49
1:B:126:VAL:HG13	1:B:172:LEU:HB3	1.94	0.48
1:C:37:ARG:C	1:C:84:LEU:HB3	2.38	0.48
1:D:231:LYS:C	1:D:232:ASN:ND2	2.70	0.48
1:E:212:ILE:HG22	1:E:242:LEU:HB2	1.95	0.48
1:D:94:VAL:HG11	1:D:214:ARG:NH1	2.28	0.48
1:C:119:ASP:OD2	1:C:176:THR:HB	2.13	0.48
1:D:227:THR:O	1:D:229:ARG:HD2	2.13	0.48
1:D:232:ASN:N	1:D:232:ASN:ND2	2.60	0.48
1:D:53:LEU:O	1:D:53:LEU:HD23	2.14	0.48
1:F:16:PHE:O	1:F:19:LEU:HG	2.14	0.48
1:F:121:ILE:HD13	1:F:204:VAL:HG22	1.95	0.48
1:E:46:LEU:HD21	1:E:53:LEU:CD2	2.38	0.48
1:C:98:ARG:CG	1:C:181:GLU:OE2	2.62	0.48
1:E:97:LEU:HD23	1:E:97:LEU:C	2.39	0.48
1:E:212:ILE:N	1:E:212:ILE:HD12	2.28	0.48
1:A:23:LYS:O	1:A:27:PHE:HD1	1.97	0.48
1:B:37:ARG:CA	1:B:84:LEU:HD23	2.44	0.48
1:B:61:PRO:HG3	1:B:108:PHE:HB2	1.96	0.48
1:B:227:THR:O	1:B:229:ARG:N	2.45	0.48
1:E:87:ARG:HG2	1:E:87:ARG:HH11	1.78	0.48
1:F:18:SER:O	1:F:23:LYS:HD2	2.14	0.48
1:A:65:ILE:O	1:A:69:GLN:HG3	2.14	0.47
1:B:227:THR:C	1:B:229:ARG:H	2.22	0.47
1:C:14:GLU:CA	1:C:75:ILE:HD11	2.36	0.47
1:D:227:THR:CB	1:D:229:ARG:HD3	2.41	0.47
1:B:30:ILE:O	1:B:30:ILE:HG22	2.12	0.47
1:D:94:VAL:C	1:D:95:ILE:HD12	2.39	0.47
1:D:184:GLU:HG3	1:D:184:GLU:O	2.14	0.47
1:A:227:THR:O	1:A:228:LYS:HB2	2.14	0.47
1:B:62:ASP:O	1:B:66:ARG:HG3	2.14	0.47
1:B:95:ILE:HD13	1:B:104:PHE:CE1	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:26:VAL:HG12	1:A:30:ILE:HD11	1.96	0.47
1:A:87:ARG:CZ	1:A:203:LEU:CD1	2.93	0.47
1:D:116:ARG:HH11	1:D:194:GLN:HE22	1.61	0.47
1:E:208:THR:HG23	1:E:209:PRO:HD2	1.96	0.47
1:F:94:VAL:HG21	1:F:214:ARG:NH2	2.28	0.47
1:D:25:ARG:O	1:D:25:ARG:NH1	2.47	0.47
1:F:36:VAL:HG23	1:F:37:ARG:N	2.28	0.47
1:C:95:ILE:HG12	1:C:104:PHE:CZ	2.50	0.47
1:C:101:ARG:HH11	1:C:101:ARG:HG3	1.78	0.47
1:C:213:VAL:HG12	1:C:241:PHE:HA	1.94	0.47
1:E:95:ILE:O	1:E:95:ILE:HG13	2.07	0.47
1:B:225:GLU:O	1:B:226:ARG:HB2	2.14	0.47
1:C:195:ILE:HG12	1:C:196:THR:N	2.30	0.47
1:D:57:LEU:O	1:D:57:LEU:HG	2.14	0.47
1:D:176:THR:O	1:D:177:LEU:HD12	2.14	0.47
1:E:152:SER:HB3	1:F:136:MET:HE1	1.96	0.47
1:F:121:ILE:CD1	1:F:205:ASP:HB2	2.45	0.47
1:F:26:VAL:O	1:F:30:ILE:HG13	2.15	0.47
1:A:37:ARG:C	1:A:84:LEU:HD23	2.39	0.47
1:D:46:LEU:HD11	1:D:53:LEU:CD2	2.45	0.47
1:F:109:VAL:CG1	1:F:110:ALA:N	2.78	0.47
1:B:183:LEU:H	1:B:183:LEU:CD1	2.25	0.47
1:A:86:ILE:N	1:A:86:ILE:HD12	2.30	0.46
1:E:112:ASP:OD1	1:E:214:ARG:NH1	2.47	0.46
1:F:186:LEU:HB3	1:F:190:GLU:HB2	1.98	0.46
1:A:152:SER:HB3	1:B:136:MET:HE1	1.96	0.46
1:D:18:SER:C	1:D:20:GLN:N	2.67	0.46
1:D:43:TYR:HB3	1:D:90:GLY:O	2.15	0.46
1:C:180:GLN:OE1	1:C:192:PRO:HB2	2.15	0.46
1:D:52:ASP:C	1:D:54:ALA:H	2.22	0.46
1:D:119:ASP:CG	1:D:176:THR:HG23	2.41	0.46
1:E:188:GLY:C	1:E:190:GLU:H	2.23	0.46
1:C:228:LYS:HZ2	1:D:225:GLU:HG3	1.79	0.46
1:E:18:SER:O	1:E:19:LEU:C	2.59	0.46
1:E:124:ARG:NH1	1:E:200:GLU:OE1	2.48	0.46
1:B:33:TYR:CG	1:B:34:PRO:HA	2.51	0.46
1:C:42:ASP:HA	1:C:89:SER:OG	2.16	0.46
1:E:186:LEU:HB3	1:E:190:GLU:HB2	1.97	0.46
1:B:98:ARG:HD3	1:B:181:GLU:OE1	2.15	0.46
1:D:222:VAL:HG13	1:D:233:PHE:HE2	1.81	0.46
1:E:16:PHE:CZ	1:E:22:TYR:CE1	3.03	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:94:VAL:HA	1:E:110:ALA:O	2.15	0.46
1:E:101:ARG:HG3	1:E:101:ARG:NH1	2.31	0.46
1:E:116:ARG:O	1:E:117:LYS:HB3	2.15	0.46
1:C:84:LEU:HD12	1:C:84:LEU:N	2.30	0.46
1:D:241:PHE:C	1:D:242:LEU:HA	2.39	0.45
1:A:4:VAL:HG12	1:A:8:LYS:HZ3	1.82	0.45
1:B:86:ILE:O	1:B:86:ILE:CG1	2.63	0.45
1:C:114:ILE:HG23	1:C:212:ILE:HD13	1.98	0.45
1:D:214:ARG:HG3	1:D:214:ARG:NH1	2.31	0.45
1:E:95:ILE:HD11	1:E:100:LEU:HD23	1.98	0.45
1:B:112:ASP:OD1	1:B:214:ARG:CG	2.59	0.45
1:B:223:ARG:HH11	1:B:223:ARG:HB2	1.80	0.45
1:D:14:GLU:HA	1:D:75:ILE:HD11	1.98	0.45
1:D:16:PHE:HB2	1:D:49:PHE:CE2	2.51	0.45
1:B:43:TYR:HB3	1:B:90:GLY:O	2.16	0.45
1:C:95:ILE:N	1:C:95:ILE:HD12	2.31	0.45
1:F:121:ILE:HG23	1:F:204:VAL:HG21	1.98	0.45
1:C:119:ASP:HB2	1:C:176:THR:HB	1.98	0.45
1:D:101:ARG:C	1:D:103:LYS:H	2.25	0.45
1:F:13:PHE:HE1	1:F:64:VAL:HG13	1.82	0.45
1:A:208:THR:O	1:A:209:PRO:C	2.59	0.45
1:E:174:THR:HG22	1:E:175:GLN:N	2.32	0.45
1:C:102:SER:C	1:C:104:PHE:N	2.75	0.45
1:C:213:VAL:HA	1:C:242:LEU:CB	2.47	0.45
1:A:93:ASN:ND2	1:A:95:ILE:CD1	2.77	0.44
1:B:121:ILE:HD13	1:B:204:VAL:CG1	2.48	0.44
1:E:124:ARG:NH1	1:E:200:GLU:OE2	2.50	0.44
1:A:4:VAL:HG12	1:A:8:LYS:NZ	2.33	0.44
1:B:95:ILE:CD1	1:B:104:PHE:CZ	3.00	0.44
1:D:179:LEU:O	1:D:194:GLN:HA	2.17	0.44
1:E:124:ARG:NH1	1:E:200:GLU:CD	2.76	0.44
1:F:124:ARG:NH2	1:F:143:GLN:O	2.50	0.44
1:C:227:THR:CG2	1:C:229:ARG:CD	2.95	0.44
1:F:32:LYS:O	1:F:35:ASN:N	2.48	0.44
1:A:92:SER:O	1:A:94:VAL:HG12	2.17	0.44
1:D:4:VAL:HA	1:D:8:LYS:HD3	2.00	0.44
1:A:174:THR:OG1	1:A:200:GLU:HG2	2.16	0.44
1:B:37:ARG:O	1:B:84:LEU:HA	2.18	0.44
1:D:176:THR:C	1:D:177:LEU:HD12	2.43	0.44
1:F:126:VAL:HG13	1:F:172:LEU:HB3	1.98	0.44
1:B:25:ARG:HA	1:B:28:GLU:HG3	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:93:ASN:O	1:D:93:ASN:CG	2.61	0.44
1:D:223:ARG:HG2	1:D:223:ARG:HH11	1.83	0.44
1:A:187:SER:O	1:A:188:GLY:O	2.36	0.44
1:B:98:ARG:HD3	1:B:181:GLU:OE2	2.18	0.44
1:B:220:ARG:HH11	1:B:220:ARG:CG	2.23	0.44
1:C:25:ARG:HG3	1:C:25:ARG:NH1	2.32	0.44
1:D:175:GLN:NE2	1:D:204:VAL:HG13	2.33	0.44
1:C:110:ALA:HA	1:C:215:VAL:O	2.18	0.44
1:D:120:GLU:CD	1:D:120:GLU:H	2.26	0.44
1:E:95:ILE:HD11	1:E:100:LEU:HG	1.99	0.44
1:B:220:ARG:HB2	1:B:233:PHE:CE2	2.53	0.44
1:C:212:ILE:C	1:C:242:LEU:CB	2.91	0.44
1:F:43:TYR:HB3	1:F:90:GLY:O	2.18	0.44
1:A:102:SER:HA	1:A:105:ILE:HG13	2.00	0.43
1:B:16:PHE:HZ	1:B:22:TYR:CE1	2.35	0.43
1:B:44:LEU:HD23	1:B:44:LEU:O	2.18	0.43
1:F:224:ASP:OD2	1:F:225:GLU:N	2.51	0.43
1:C:43:TYR:HB3	1:C:90:GLY:O	2.18	0.43
1:B:121:ILE:HD13	1:B:204:VAL:HG13	1.99	0.43
1:C:28:GLU:O	1:C:32:LYS:HG3	2.18	0.43
1:F:36:VAL:O	1:F:37:ARG:HB2	2.19	0.43
1:A:22:TYR:O	1:A:26:VAL:HG23	2.18	0.43
1:B:19:LEU:O	1:B:20:GLN:C	2.61	0.43
1:B:73:ARG:HD3	1:B:83:ASP:OD1	2.18	0.43
1:F:14:GLU:HA	1:F:75:ILE:HD11	2.01	0.43
1:A:98:ARG:HD3	1:A:181:GLU:OE2	2.17	0.43
1:B:122:ARG:O	1:B:174:THR:HG22	2.19	0.43
1:D:19:LEU:O	1:D:21:ASP:N	2.46	0.43
1:E:11:THR:O	1:E:12:LYS:C	2.59	0.43
1:A:79:ARG:C	1:A:81:ASN:H	2.26	0.43
1:A:22:TYR:HA	1:A:25:ARG:CG	2.49	0.43
1:B:225:GLU:OE1	1:B:225:GLU:HA	2.18	0.43
1:E:33:TYR:HA	1:E:34:PRO:HA	1.77	0.43
1:C:184:GLU:H	1:C:184:GLU:HG2	1.53	0.43
1:E:124:ARG:HH11	1:E:200:GLU:CD	2.27	0.43
1:E:50:ASP:OD1	1:E:50:ASP:C	2.62	0.43
1:E:223:ARG:CD	1:E:228:LYS:O	2.65	0.43
1:D:126:VAL:HG13	1:D:172:LEU:HB2	2.01	0.43
1:A:224:ASP:C	1:A:224:ASP:OD1	2.61	0.42
1:C:98:ARG:CD	1:C:193:ARG:HB2	2.49	0.42
1:D:19:LEU:O	1:D:20:GLN:HB2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:213:VAL:HG12	1:D:241:PHE:CA	2.46	0.42
1:F:14:GLU:O	1:F:18:SER:OG	2.32	0.42
1:F:131:GLU:HB2	1:F:165:LEU:HD11	2.00	0.42
1:A:36:VAL:O	1:A:37:ARG:CB	2.63	0.42
1:D:115:VAL:HG13	1:D:211:ASP:H	1.82	0.42
1:E:203:LEU:HD23	1:E:203:LEU:HA	1.71	0.42
1:E:223:ARG:HH22	1:F:148:ILE:CD1	2.30	0.42
1:A:33:TYR:CD1	1:A:34:PRO:HA	2.54	0.42
1:B:98:ARG:HD3	1:B:181:GLU:CD	2.44	0.42
1:C:47:GLU:HA	1:C:54:ALA:HB2	2.01	0.42
1:E:225:GLU:O	1:E:227:THR:N	2.52	0.42
1:F:78:LEU:C	1:F:80:LYS:H	2.27	0.42
1:F:121:ILE:CD1	1:F:204:VAL:HG22	2.49	0.42
1:B:104:PHE:C	1:B:219:LEU:HD23	2.44	0.42
1:E:39:ILE:CD1	1:E:84:LEU:HD13	2.50	0.42
1:A:43:TYR:CE1	1:A:54:ALA:HB1	2.54	0.42
1:D:78:LEU:HB2	1:D:80:LYS:HE2	2.01	0.42
1:E:57:LEU:HD23	1:E:64:VAL:HG11	2.00	0.42
1:B:29:ALA:C	1:B:31:GLU:N	2.76	0.42
1:D:32:LYS:N	1:D:32:LYS:HE2	2.35	0.42
1:A:98:ARG:HD3	1:A:181:GLU:CD	2.45	0.42
1:B:69:GLN:HE21	1:B:85:ASN:HA	1.83	0.42
1:C:98:ARG:CZ	1:C:193:ARG:NE	2.83	0.42
1:F:47:GLU:HA	1:F:54:ALA:HB2	2.00	0.42
1:B:64:VAL:O	1:B:65:ILE:C	2.62	0.42
1:C:30:ILE:HD13	1:C:76:ASP:OD1	2.20	0.42
1:D:183:LEU:O	1:D:185:ASN:N	2.48	0.42
1:A:4:VAL:HB	1:A:8:LYS:HB2	2.00	0.42
1:A:9:THR:HG23	1:A:53:LEU:CD1	2.50	0.42
1:A:57:LEU:HD23	1:A:64:VAL:HG11	2.01	0.42
1:B:95:ILE:HD12	1:B:109:VAL:HG22	2.02	0.42
1:C:102:SER:C	1:C:104:PHE:H	2.28	0.42
1:E:95:ILE:HD13	1:E:104:PHE:CE1	2.55	0.42
1:F:226:ARG:HG3	1:F:227:THR:N	2.35	0.42
1:C:102:SER:O	1:C:104:PHE:N	2.53	0.41
1:D:59:GLU:HA	1:D:107:LYS:HD2	2.02	0.41
1:D:172:LEU:HD23	1:D:172:LEU:HA	1.80	0.41
1:E:152:SER:HA	1:F:163:ARG:NH2	2.35	0.41
1:F:95:ILE:HD13	1:F:109:VAL:HG13	2.02	0.41
1:F:183:LEU:O	1:F:185:ASN:N	2.46	0.41
1:F:187:SER:O	1:F:188:GLY:O	2.38	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:43:TYR:HB3	1:A:90:GLY:O	2.20	0.41
1:E:63:ASP:O	1:E:66:ARG:HB2	2.20	0.41
1:C:58:ILE:HA	1:C:108:PHE:CB	2.45	0.41
1:A:61:PRO:HB3	1:A:108:PHE:HB2	2.03	0.41
1:A:69:GLN:HG2	1:A:86:ILE:HD13	2.02	0.41
1:A:119:ASP:OD2	1:A:176:THR:CG2	2.67	0.41
1:E:36:VAL:HG23	1:E:37:ARG:N	2.35	0.41
1:B:124:ARG:HH11	1:B:200:GLU:CD	2.27	0.41
1:F:178:LYS:HE2	1:F:196:THR:HG23	2.03	0.41
1:C:80:LYS:HB3	1:C:82:VAL:HG13	2.03	0.41
1:B:44:LEU:HD23	1:B:44:LEU:C	2.45	0.41
1:B:88:PHE:H	1:B:237:ASN:ND2	2.07	0.41
1:C:98:ARG:NE	1:C:193:ARG:NE	2.69	0.41
1:C:98:ARG:CZ	1:C:193:ARG:CZ	2.95	0.41
1:C:109:VAL:CG1	1:C:110:ALA:N	2.83	0.41
1:D:227:THR:C	1:D:229:ARG:N	2.78	0.41
1:A:126:VAL:CG1	1:A:172:LEU:HD22	2.51	0.41
1:A:229:ARG:HD3	1:B:122:ARG:CZ	2.51	0.41
1:D:183:LEU:HD23	1:D:183:LEU:HA	1.92	0.41
1:A:95:ILE:HG12	1:A:104:PHE:CE1	2.55	0.41
1:A:181:GLU:OE1	1:A:192:PRO:HA	2.21	0.41
1:B:104:PHE:O	1:B:219:LEU:HD23	2.20	0.41
1:B:124:ARG:NH2	1:B:143:GLN:O	2.54	0.41
1:B:187:SER:O	1:B:188:GLY:O	2.39	0.41
1:D:52:ASP:C	1:D:54:ALA:N	2.78	0.41
1:D:87:ARG:NH2	1:D:202:ASP:HB3	2.36	0.41
1:E:22:TYR:O	1:E:23:LYS:C	2.64	0.41
1:B:121:ILE:CD1	1:B:204:VAL:HG13	2.51	0.41
1:D:228:LYS:HA	1:D:228:LYS:HD2	1.91	0.41
1:E:95:ILE:HD11	1:E:100:LEU:CD2	2.50	0.41
1:C:42:ASP:HB3	1:C:45:ASP:OD2	2.21	0.40
1:C:46:LEU:HD11	1:C:53:LEU:HD23	2.02	0.40
1:C:121:ILE:HD12	1:C:121:ILE:N	2.36	0.40
1:D:56:LEU:HD23	1:D:56:LEU:O	2.21	0.40
1:D:104:PHE:C	1:D:219:LEU:HD23	2.46	0.40
1:E:19:LEU:HD11	1:E:49:PHE:CD1	2.57	0.40
1:E:121:ILE:CD1	1:E:205:ASP:HB2	2.51	0.40
1:E:222:VAL:HG23	1:E:233:PHE:HE2	1.86	0.40
1:A:22:TYR:HA	1:A:25:ARG:HG2	2.02	0.40
1:B:36:VAL:C	1:B:38:SER:H	2.29	0.40
1:D:68:ALA:O	1:D:71:ALA:HB3	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:11:THR:O	1:A:12:LYS:C	2.65	0.40
1:B:8:LYS:O	1:B:12:LYS:HG3	2.22	0.40
1:B:87:ARG:HG2	1:B:87:ARG:HH11	1.86	0.40
1:C:4:VAL:C	1:C:6:LYS:H	2.29	0.40
1:D:14:GLU:CA	1:D:75:ILE:HD11	2.51	0.40
1:D:108:PHE:HD2	1:D:108:PHE:HA	1.73	0.40
1:F:18:SER:O	1:F:19:LEU:C	2.64	0.40
1:F:95:ILE:HG22	1:F:96:PRO:O	2.21	0.40
1:A:37:ARG:O	1:A:85:ASN:N	2.42	0.40
1:A:183:LEU:O	1:A:186:LEU:HG	2.21	0.40
1:C:93:ASN:OD1	1:C:95:ILE:HD11	2.21	0.40
1:C:112:ASP:OD1	1:C:214:ARG:NH1	2.54	0.40
1:D:221:THR:HG22	1:D:232:ASN:CG	2.46	0.40
1:D:172:LEU:HD22	1:D:201:ASP:OD1	2.22	0.40
1:E:3:THR:O	1:E:4:VAL:HB	2.21	0.40
1:F:33:TYR:HA	1:F:34:PRO:HA	1.81	0.40
1:F:100:LEU:HD21	1:F:109:VAL:HG11	2.04	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	237/279 (85%)	214 (90%)	18 (8%)	5 (2%)	5	27
1	B	237/279 (85%)	208 (88%)	21 (9%)	8 (3%)	3	17
1	C	237/279 (85%)	209 (88%)	21 (9%)	7 (3%)	3	19
1	D	237/279 (85%)	199 (84%)	28 (12%)	10 (4%)	2	13
1	E	240/279 (86%)	213 (89%)	19 (8%)	8 (3%)	3	17
1	F	237/279 (85%)	207 (87%)	24 (10%)	6 (2%)	4	23
All	All	1425/1674 (85%)	1250 (88%)	131 (9%)	44 (3%)	3	19

All (44) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	188	GLY
1	B	228	LYS
1	C	188	GLY
1	D	92	SER
1	D	188	GLY
1	E	2	LYS
1	E	3	THR
1	E	93	ASN
1	E	228	LYS
1	F	188	GLY
1	A	185	ASN
1	B	184	GLU
1	B	188	GLY
1	C	90	GLY
1	C	184	GLU
1	D	90	GLY
1	E	4	VAL
1	E	79	ARG
1	E	188	GLY
1	F	80	LYS
1	A	80	LYS
1	A	82	VAL
1	B	82	VAL
1	C	120	GLU
1	D	6	LYS
1	D	19	LEU
1	D	185	ASN
1	F	184	GLU
1	F	185	ASN
1	A	81	ASN
1	B	48	MET
1	B	185	ASN
1	B	226	ARG
1	C	103	LYS
1	C	185	ASN
1	D	93	ASN
1	F	76	ASP
1	C	40	GLU
1	D	228	LYS
1	E	94	VAL
1	D	79	ARG
1	B	75	ILE

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Mol	Chain	Res	Type
1	F	75	ILE
1	D	209	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	219/256 (86%)	204 (93%)	15 (7%)	14	45
1	B	219/256 (86%)	199 (91%)	20 (9%)	9	33
1	C	215/256 (84%)	199 (93%)	16 (7%)	13	42
1	D	213/256 (83%)	200 (94%)	13 (6%)	17	49
1	E	219/256 (86%)	207 (94%)	12 (6%)	19	53
1	F	218/256 (85%)	204 (94%)	14 (6%)	16	48
All	All	1303/1536 (85%)	1213 (93%)	90 (7%)	14	45

All (90) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	LYS
1	A	21	ASP
1	A	25	ARG
1	A	46	LEU
1	A	70	GLN
1	A	94	VAL
1	A	118	THR
1	A	126	VAL
1	A	172	LEU
1	A	174	THR
1	A	177	LEU
1	A	184	GLU
1	A	208	THR
1	A	209	PRO
1	A	239	THR
1	B	14	GLU

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Mol	Chain	Res	Type
1	B	32	LYS
1	B	36	VAL
1	B	46	LEU
1	B	50	ASP
1	B	60	LYS
1	B	77	ARG
1	B	91	ILE
1	B	109	VAL
1	B	124	ARG
1	B	126	VAL
1	B	172	LEU
1	B	174	THR
1	B	183	LEU
1	B	193	ARG
1	B	204	VAL
1	B	208	THR
1	B	220	ARG
1	B	224	ASP
1	B	239	THR
1	C	44	LEU
1	C	46	LEU
1	C	70	GLN
1	C	84	LEU
1	C	93	ASN
1	C	108	PHE
1	C	122	ARG
1	C	124	ARG
1	C	126	VAL
1	C	184	GLU
1	C	191	GLN
1	C	201	ASP
1	C	208	THR
1	C	213	VAL
1	C	218	THR
1	C	229	ARG
1	D	27	PHE
1	D	63	ASP
1	D	73	ARG
1	D	80	LYS
1	D	91	ILE
1	D	108	PHE
1	D	115	VAL

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Mol	Chain	Res	Type
1	D	117	LYS
1	D	124	ARG
1	D	126	VAL
1	D	176	THR
1	D	191	GLN
1	D	232	ASN
1	E	36	VAL
1	E	41	VAL
1	E	56	LEU
1	E	57	LEU
1	E	66	ARG
1	E	95	ILE
1	E	122	ARG
1	E	126	VAL
1	E	176	THR
1	E	184	GLU
1	E	204	VAL
1	E	214	ARG
1	F	5	ASP
1	F	21	ASP
1	F	28	GLU
1	F	52	ASP
1	F	56	LEU
1	F	84	LEU
1	F	87	ARG
1	F	95	ILE
1	F	126	VAL
1	F	172	LEU
1	F	196	THR
1	F	208	THR
1	F	214	ARG
1	F	225	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (34) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	81	ASN
1	A	85	ASN
1	A	93	ASN
1	A	139	HIS
1	A	146	ASN
1	A	180	GLN

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Mol	Chain	Res	Type
1	A	237	ASN
1	B	69	GLN
1	B	93	ASN
1	B	139	HIS
1	B	191	GLN
1	B	237	ASN
1	C	69	GLN
1	C	70	GLN
1	C	93	ASN
1	C	139	HIS
1	C	185	ASN
1	D	20	GLN
1	D	70	GLN
1	D	93	ASN
1	D	139	HIS
1	D	175	GLN
1	D	194	GLN
1	D	232	ASN
1	E	20	GLN
1	E	81	ASN
1	E	93	ASN
1	E	139	HIS
1	E	180	GLN
1	F	69	GLN
1	F	81	ASN
1	F	139	HIS
1	F	185	ASN
1	F	194	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 [i](#)

Of 6 ligands modelled in this entry, 6 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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	D	4
1	C	2

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	D	182:PRO	C	183:LEU	N	1.74
1	D	241:PHE	C	242:LEU	N	1.68
1	C	193:ARG	C	194:GLN	N	1.16
1	D	239:THR	C	240:GLU	N	1.13
1	C	241:PHE	C	242:LEU	N	1.00
1	D	237:ASN	C	238:TYR	N	0.53

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	239/279 (85%)	0.04	5 (2%)	63 40	15, 46, 107, 117	0
1	B	239/279 (85%)	-0.06	8 (3%)	49 28	13, 40, 94, 110	0
1	C	239/279 (85%)	0.62	20 (8%)	17 9	25, 81, 114, 124	0
1	D	239/279 (85%)	0.82	21 (8%)	15 8	23, 88, 116, 127	0
1	E	242/279 (86%)	-0.08	5 (2%)	63 40	17, 46, 100, 117	0
1	F	239/279 (85%)	-0.11	1 (0%)	88 76	17, 47, 95, 108	0
All	All	1437/1674 (85%)	0.20	60 (4%)	40 22	13, 55, 112, 127	0

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	182	PRO	4.6
1	C	225	GLU	4.5
1	D	4	VAL	4.2
1	C	226	ARG	4.1
1	D	191	GLN	4.0
1	C	224	ASP	4.0
1	D	182	PRO	3.9
1	B	49	PHE	3.7
1	A	4	VAL	3.6
1	C	191	GLN	3.6
1	D	187	SER	3.6
1	D	184	GLU	3.5
1	C	228	LYS	3.4
1	D	189	GLY	3.2
1	C	181	GLU	3.1
1	B	77	ARG	3.1
1	D	192	PRO	3.1
1	D	183	LEU	3.1
1	D	114	ILE	3.0

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Mol	Chain	Res	Type	RSRZ
1	C	229	ARG	3.0
1	D	241	PHE	3.0
1	D	190	GLU	3.0
1	D	188	GLY	2.9
1	B	75	ILE	2.8
1	D	240	GLU	2.7
1	E	1	MET	2.7
1	C	241	PHE	2.6
1	C	188	GLY	2.6
1	C	186	LEU	2.6
1	D	78	LEU	2.6
1	B	78	LEU	2.5
1	C	180	GLN	2.5
1	C	120	GLU	2.5
1	D	174	THR	2.5
1	A	78	LEU	2.5
1	C	184	GLU	2.5
1	C	27	PHE	2.5
1	D	186	LEU	2.4
1	E	227	THR	2.4
1	D	21	ASP	2.4
1	C	193	ARG	2.4
1	A	39	ILE	2.4
1	C	192	PRO	2.4
1	B	4	VAL	2.3
1	C	190	GLU	2.3
1	C	4	VAL	2.3
1	B	50	ASP	2.3
1	D	19	LEU	2.3
1	A	5	ASP	2.3
1	D	113	GLY	2.2
1	C	78	LEU	2.2
1	E	75	ILE	2.2
1	E	187	SER	2.2
1	B	76	ASP	2.2
1	A	76	ASP	2.1
1	B	5	ASP	2.1
1	D	45	ASP	2.1
1	F	122	ARG	2.0
1	D	81	ASN	2.0
1	E	24	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($\text{\AA}^2$)	Q<0.9
2	ZN	D	275	1/1	0.97	0.28	86,86,86,86	0
2	ZN	B	275	1/1	0.98	0.27	66,66,66,66	0
2	ZN	C	275	1/1	0.98	0.23	77,77,77,77	0
2	ZN	A	275	1/1	0.98	0.26	56,56,56,56	0
2	ZN	E	275	1/1	0.99	0.17	55,55,55,55	0
2	ZN	F	275	1/1	0.99	0.16	50,50,50,50	0

6.5 Other polymers [i](#)

There are no such residues in this entry.