



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:36 AM UTC

PDB ID : 3SO4 / pdb_00003so4
Title : Methionine-adenosyltransferase from Entamoeba histolytica
Authors : Merritt, E.A.; Medical Structural Genomics of Pathogenic Protozoa (MSGPP)
Deposited on : 2011-06-29
Resolution : 3.18 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

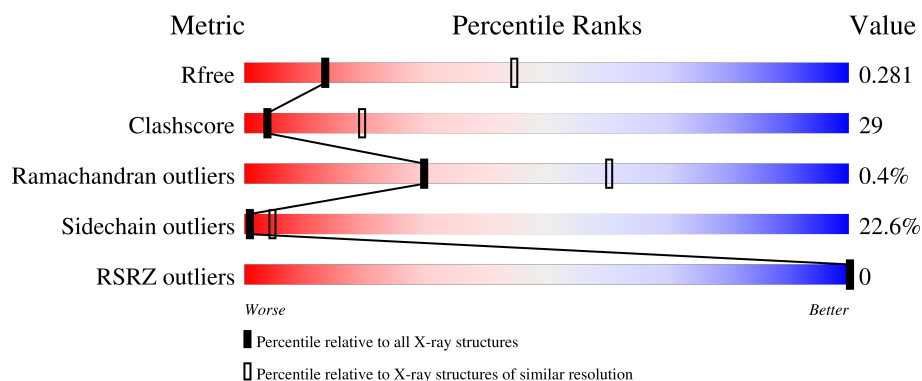
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.18 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	2001 (3.20-3.16)
Clashscore	190562	2119 (3.20-3.16)
Ramachandran outliers	187476	2070 (3.20-3.16)
Sidechain outliers	187428	2069 (3.20-3.16)
RSRZ outliers	180081	2001 (3.20-3.16)

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	415	
1	B	415	
1	C	415	
1	D	415	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ACT	C	400	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11498 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 Methionine-adenosyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	379	Total	C	N	O	S	0	0	0
			2858	1804	487	549	18			
1	B	379	Total	C	N	O	S	0	0	0
			2881	1818	491	554	18			
1	C	379	Total	C	N	O	S	0	0	0
			2865	1807	490	550	18			
1	D	379	Total	C	N	O	S	0	0	0
			2874	1815	491	550	18			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-17	MET	-	expression tag	UNP C4M272
A	-16	ALA	-	expression tag	UNP C4M272
A	-15	HIS	-	expression tag	UNP C4M272
A	-14	HIS	-	expression tag	UNP C4M272
A	-13	HIS	-	expression tag	UNP C4M272
A	-12	HIS	-	expression tag	UNP C4M272
A	-11	HIS	-	expression tag	UNP C4M272
A	-10	HIS	-	expression tag	UNP C4M272
A	-9	MET	-	expression tag	UNP C4M272
A	-8	GLY	-	expression tag	UNP C4M272
A	-7	THR	-	expression tag	UNP C4M272
A	-6	LEU	-	expression tag	UNP C4M272
A	-5	GLU	-	expression tag	UNP C4M272
A	-4	ALA	-	expression tag	UNP C4M272
A	-3	GLN	-	expression tag	UNP C4M272
A	-2	THR	-	expression tag	UNP C4M272
A	-1	GLN	-	expression tag	UNP C4M272
A	0	GLY	-	expression tag	UNP C4M272
A	1	PRO	-	expression tag	UNP C4M272
A	2	GLY	-	expression tag	UNP C4M272
A	3	SER	-	expression tag	UNP C4M272

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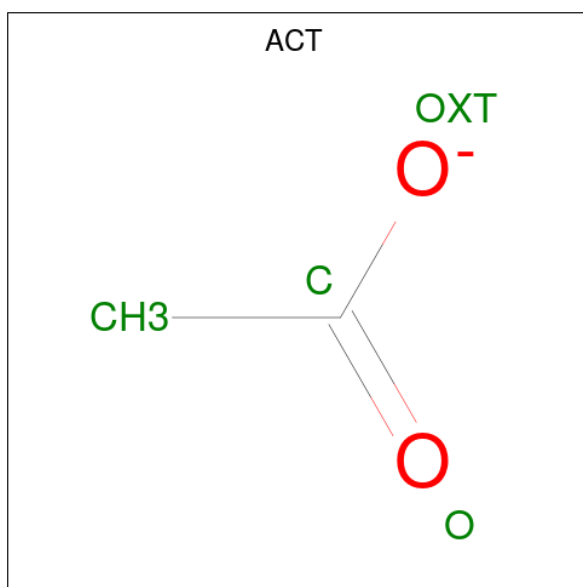
Chain	Residue	Modelled	Actual	Comment	Reference
B	-17	MET	-	expression tag	UNP C4M272
B	-16	ALA	-	expression tag	UNP C4M272
B	-15	HIS	-	expression tag	UNP C4M272
B	-14	HIS	-	expression tag	UNP C4M272
B	-13	HIS	-	expression tag	UNP C4M272
B	-12	HIS	-	expression tag	UNP C4M272
B	-11	HIS	-	expression tag	UNP C4M272
B	-10	HIS	-	expression tag	UNP C4M272
B	-9	MET	-	expression tag	UNP C4M272
B	-8	GLY	-	expression tag	UNP C4M272
B	-7	THR	-	expression tag	UNP C4M272
B	-6	LEU	-	expression tag	UNP C4M272
B	-5	GLU	-	expression tag	UNP C4M272
B	-4	ALA	-	expression tag	UNP C4M272
B	-3	GLN	-	expression tag	UNP C4M272
B	-2	THR	-	expression tag	UNP C4M272
B	-1	GLN	-	expression tag	UNP C4M272
B	0	GLY	-	expression tag	UNP C4M272
B	1	PRO	-	expression tag	UNP C4M272
B	2	GLY	-	expression tag	UNP C4M272
B	3	SER	-	expression tag	UNP C4M272
C	-17	MET	-	expression tag	UNP C4M272
C	-16	ALA	-	expression tag	UNP C4M272
C	-15	HIS	-	expression tag	UNP C4M272
C	-14	HIS	-	expression tag	UNP C4M272
C	-13	HIS	-	expression tag	UNP C4M272
C	-12	HIS	-	expression tag	UNP C4M272
C	-11	HIS	-	expression tag	UNP C4M272
C	-10	HIS	-	expression tag	UNP C4M272
C	-9	MET	-	expression tag	UNP C4M272
C	-8	GLY	-	expression tag	UNP C4M272
C	-7	THR	-	expression tag	UNP C4M272
C	-6	LEU	-	expression tag	UNP C4M272
C	-5	GLU	-	expression tag	UNP C4M272
C	-4	ALA	-	expression tag	UNP C4M272
C	-3	GLN	-	expression tag	UNP C4M272
C	-2	THR	-	expression tag	UNP C4M272
C	-1	GLN	-	expression tag	UNP C4M272
C	0	GLY	-	expression tag	UNP C4M272
C	1	PRO	-	expression tag	UNP C4M272
C	2	GLY	-	expression tag	UNP C4M272
C	3	SER	-	expression tag	UNP C4M272

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-17	MET	-	expression tag	UNP C4M272
D	-16	ALA	-	expression tag	UNP C4M272
D	-15	HIS	-	expression tag	UNP C4M272
D	-14	HIS	-	expression tag	UNP C4M272
D	-13	HIS	-	expression tag	UNP C4M272
D	-12	HIS	-	expression tag	UNP C4M272
D	-11	HIS	-	expression tag	UNP C4M272
D	-10	HIS	-	expression tag	UNP C4M272
D	-9	MET	-	expression tag	UNP C4M272
D	-8	GLY	-	expression tag	UNP C4M272
D	-7	THR	-	expression tag	UNP C4M272
D	-6	LEU	-	expression tag	UNP C4M272
D	-5	GLU	-	expression tag	UNP C4M272
D	-4	ALA	-	expression tag	UNP C4M272
D	-3	GLN	-	expression tag	UNP C4M272
D	-2	THR	-	expression tag	UNP C4M272
D	-1	GLN	-	expression tag	UNP C4M272
D	0	GLY	-	expression tag	UNP C4M272
D	1	PRO	-	expression tag	UNP C4M272
D	2	GLY	-	expression tag	UNP C4M272
D	3	SER	-	expression tag	UNP C4M272

- Molecule 2 is ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		

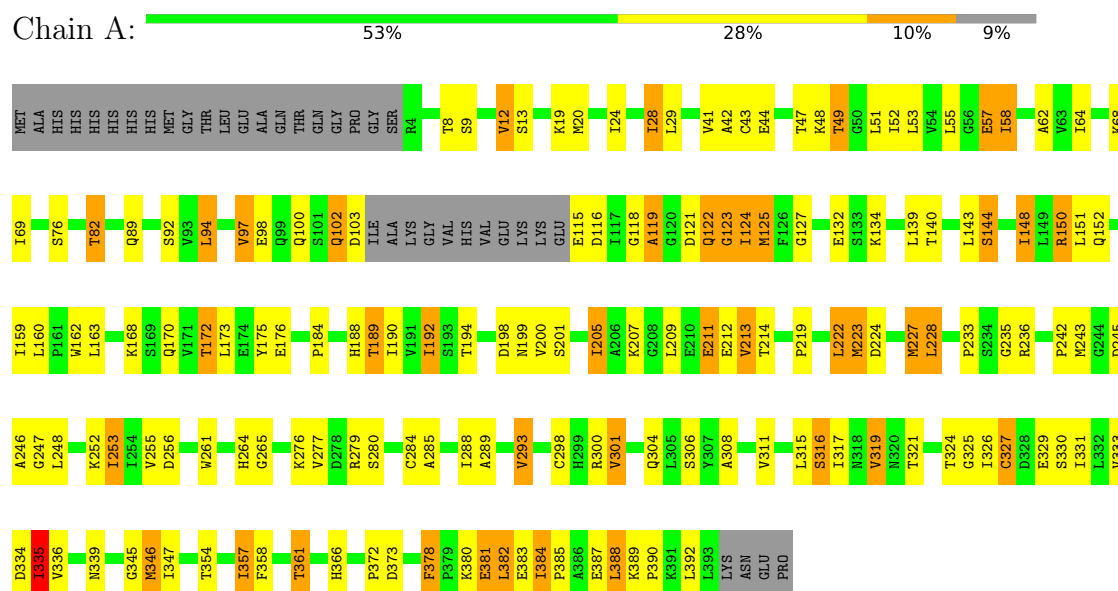
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	O	0	0
			1	1		
3	B	1	Total	O	0	0
			1	1		
3	C	1	Total	O	0	0
			1	1		
3	D	1	Total	O	0	0
			1	1		

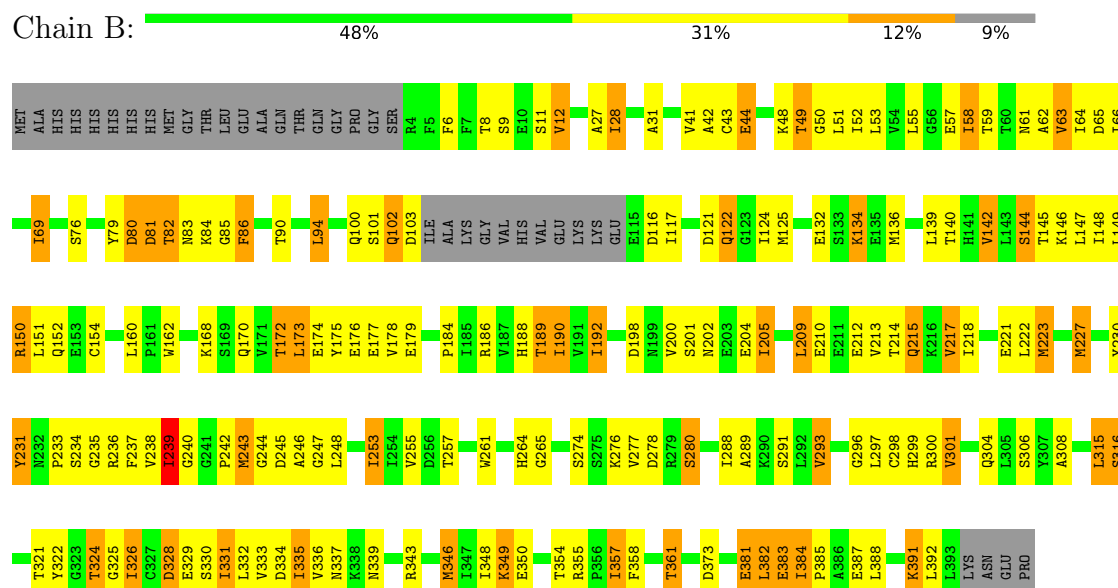
3 Residue-property plots

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: Methionine-adenosyltransferase

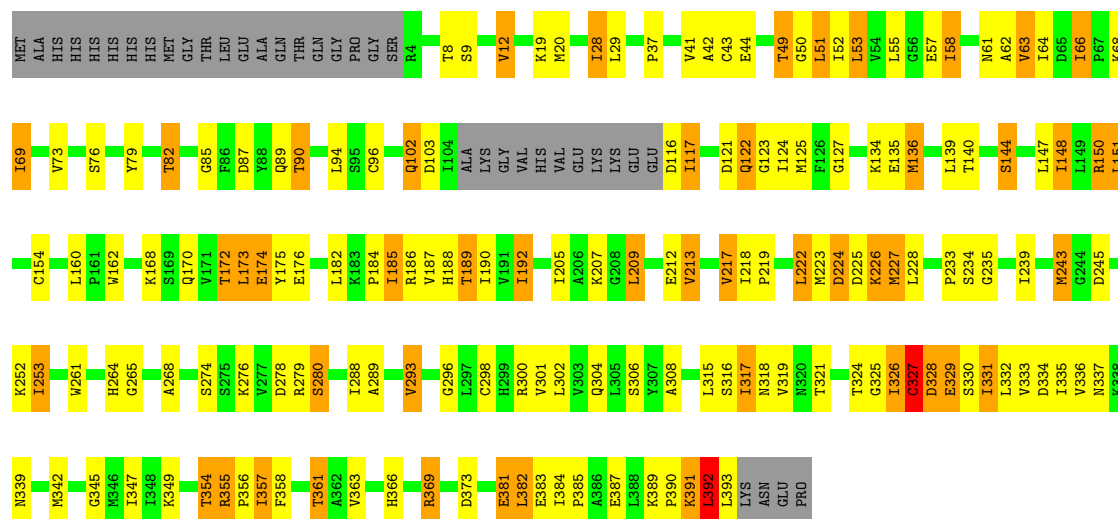


• Molecule 1: Methionine-adenosyltransferase



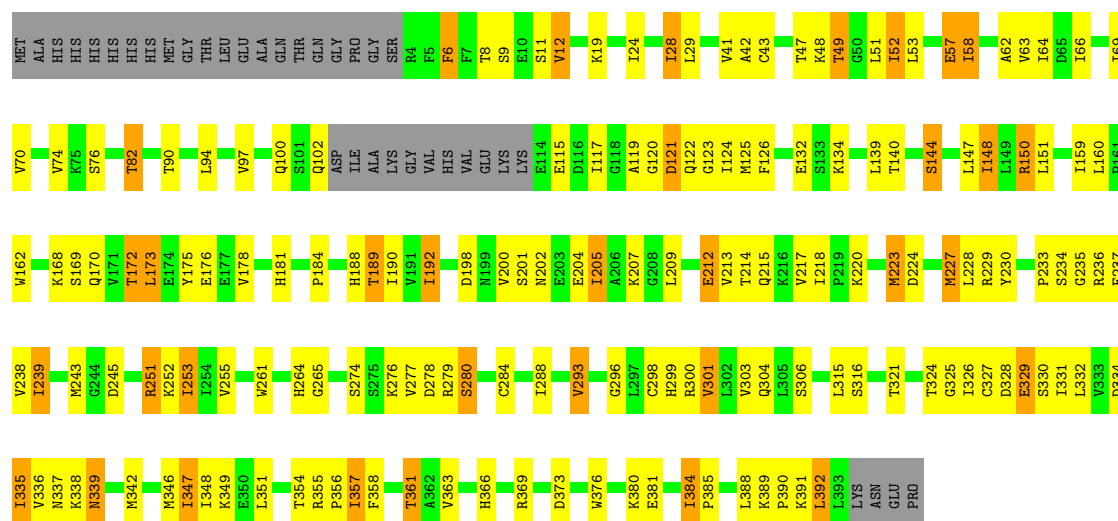
• Molecule 1: Methionine-adenosyltransferase

Chain C: 



• Molecule 1: Methionine-adenosyltransferase

Chain D: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	58.49Å 113.16Å 220.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 3.18 50.00 – 3.18	Depositor EDS
% Data completeness (in resolution range)	91.1 (50.00-3.18) 91.3 (50.00-3.18)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.50 (at 3.18Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.228 , 0.284 0.230 , 0.281	Depositor DCC
R_{free} test set	1296 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	76.0	Xtriage
Anisotropy	0.453	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 55.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	11498	wwPDB-VP
Average B, all atoms (Å ²)	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.43% 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: ACT

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	0.62	0/2914	0.96	5/3952 (0.1%)
1	B	0.64	0/2937	0.99	5/3978 (0.1%)
1	C	0.58	0/2921	0.97	5/3959 (0.1%)
1	D	0.58	0/2930	0.95	4/3969 (0.1%)
All	All	0.61	0/11702	0.96	19/15858 (0.1%)

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

There are no bond length outliers.

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	329	GLU	N-CA-C	8.33	120.44	111.36
1	C	225	ASP	N-CA-C	-8.09	103.19	113.23
1	A	213	VAL	CB-CA-C	-6.42	100.76	111.29
1	B	82	THR	N-CA-C	-6.40	104.57	112.38
1	A	119	ALA	N-CA-C	-6.28	101.25	110.42
1	B	239	ILE	N-CA-C	6.28	117.76	108.65
1	C	391	LYS	N-CA-C	5.80	120.15	112.72
1	B	239	ILE	CA-C-O	5.76	127.28	121.64
1	A	335	ILE	CB-CA-C	-5.75	104.51	112.04
1	D	392	LEU	N-CA-C	5.35	117.20	109.07
1	D	335	ILE	CB-CA-C	-5.29	105.11	111.88
1	D	123	GLY	N-CA-C	5.21	118.36	110.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	123	GLY	N-CA-C	5.17	118.30	110.75
1	C	326	ILE	N-CA-C	5.13	115.91	110.72
1	A	97	VAL	CB-CA-C	-5.08	103.82	110.98
1	B	328	ASP	N-CA-C	-5.06	103.91	110.19
1	A	123	GLY	N-CA-C	5.06	118.13	110.75
1	C	347	ILE	CB-CA-C	-5.04	105.44	112.04
1	B	101	SER	N-CA-C	-5.01	104.24	110.41

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	239	ILE	Peptide

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	2858	0	2785	155	0
1	B	2881	0	2830	190	0
1	C	2865	0	2799	185	0
1	D	2874	0	2824	166	0
2	A	4	0	3	1	0
2	B	4	0	3	0	0
2	C	4	0	3	3	0
2	D	4	0	3	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
All	All	11498	0	11250	652	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 29.

All (652) 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:A:209:LEU:O	1:A:213:VAL:HG22	1.24	1.29
1:A:213:VAL:CG2	1:A:214:THR:H	1.37	1.28
1:D:47:THR:CG2	1:D:52:ILE:HD13	1.63	1.27
1:A:213:VAL:HG23	1:A:214:THR:N	1.26	1.20
1:C:326:ILE:HD12	1:C:327:CYS:N	1.58	1.16
1:A:327:CYS:SG	1:A:388:LEU:CD2	2.36	1.13
1:D:47:THR:HG21	1:D:52:ILE:HD13	1.24	1.10
1:B:209:LEU:HD23	1:B:209:LEU:H	1.02	1.10
1:B:385:PRO:CG	1:B:388:LEU:HD12	1.81	1.09
1:B:385:PRO:HG2	1:B:388:LEU:CD1	1.81	1.08
1:B:205:ILE:O	1:B:209:LEU:CD2	2.00	1.08
1:C:187:VAL:HB	1:C:227:MET:HE1	1.36	1.08
1:A:331:ILE:O	1:A:335:ILE:HD13	1.51	1.07
1:D:181:HIS:ND1	1:D:299:HIS:ND1	2.01	1.06
1:B:205:ILE:O	1:B:209:LEU:HD23	1.52	1.04
1:A:198:ASP:OD2	1:A:236:ARG:NH2	1.91	1.04
1:B:385:PRO:HG2	1:B:388:LEU:HD12	1.07	1.03
1:C:224:ASP:OD2	1:C:226:LYS:HB2	1.58	1.03
1:D:173:LEU:HD12	1:D:173:LEU:C	1.85	1.02
1:C:392:LEU:HD23	1:C:393:LEU:H	1.25	1.01
1:B:308:ALA:HB2	1:B:315:LEU:HD23	1.43	1.00
1:D:173:LEU:HD12	1:D:173:LEU:O	1.61	1.00
1:A:248:LEU:HD23	1:B:248:LEU:HD23	1.44	1.00
1:D:147:LEU:HD12	1:D:217:VAL:HG21	1.45	0.98
1:C:288:ILE:HD11	1:C:342:MET:HE2	1.44	0.98
1:B:209:LEU:HD23	1:B:209:LEU:N	1.73	0.97
1:A:327:CYS:SG	1:A:388:LEU:HD23	2.04	0.97
1:A:209:LEU:O	1:A:213:VAL:CG2	2.11	0.97
1:C:117:ILE:HG23	1:C:117:ILE:O	1.64	0.96
1:D:173:LEU:HD11	1:D:175:TYR:CE1	1.99	0.96
1:D:299:HIS:CD2	1:D:392:LEU:HD11	2.01	0.95
1:C:173:LEU:HD11	1:C:175:TYR:CE1	2.01	0.94
1:D:391:LYS:HG3	1:D:392:LEU:H	1.29	0.94
1:A:385:PRO:O	1:A:388:LEU:HD12	1.66	0.94
1:B:63:VAL:HG11	1:C:63:VAL:HG11	1.49	0.93
1:C:327:CYS:SG	1:C:328:ASP:N	2.32	0.93
1:C:37:PRO:HB3	1:C:354:THR:HG23	1.51	0.93
1:C:213:VAL:O	1:C:217:VAL:HG12	1.66	0.92
1:C:174:GLU:HG3	1:C:175:TYR:N	1.84	0.92
1:C:326:ILE:HD12	1:C:327:CYS:H	1.21	0.91
1:B:209:LEU:O	1:B:213:VAL:HG22	1.69	0.91
1:B:298:CYS:SG	1:B:301:VAL:HG22	2.10	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:335:ILE:HD13	1:A:335:ILE:H	1.35	0.91
1:A:388:LEU:HD12	1:A:388:LEU:H	1.33	0.91
1:B:188:HIS:ND1	1:B:189:THR:HG22	1.87	0.90
1:D:173:LEU:C	1:D:173:LEU:CD1	2.44	0.90
1:A:327:CYS:SG	1:A:388:LEU:HD21	2.08	0.90
1:D:298:CYS:SG	1:D:301:VAL:HG22	2.12	0.90
1:C:187:VAL:HB	1:C:227:MET:CE	2.01	0.89
1:B:209:LEU:H	1:B:209:LEU:CD2	1.86	0.88
1:C:336:VAL:CG1	1:C:342:MET:HE1	2.04	0.88
1:D:299:HIS:HD2	1:D:392:LEU:HD11	1.32	0.88
1:C:173:LEU:C	1:C:173:LEU:HD12	1.99	0.88
1:D:48:LYS:HG3	1:D:49:THR:N	1.88	0.88
1:B:213:VAL:O	1:B:217:VAL:HG13	1.71	0.88
1:B:90:THR:HB	1:D:243:MET:HE1	1.54	0.87
1:D:47:THR:HG21	1:D:52:ILE:CD1	2.02	0.87
1:C:55:LEU:HD11	1:D:48:LYS:CB	2.04	0.87
1:C:117:ILE:HG21	1:C:345:GLY:CA	2.04	0.87
1:D:173:LEU:HD11	1:D:175:TYR:CD1	2.11	0.86
1:B:209:LEU:CD2	1:B:209:LEU:N	2.39	0.86
1:A:150:ARG:NH1	1:A:212:GLU:O	2.09	0.85
1:B:150:ARG:NH1	1:B:212:GLU:O	2.10	0.85
1:B:49:THR:HG22	1:B:50:GLY:N	1.87	0.85
1:D:150:ARG:NH1	1:D:212:GLU:O	2.11	0.84
1:C:150:ARG:NH1	1:C:212:GLU:O	2.10	0.84
1:A:213:VAL:HG23	1:A:214:THR:H	0.72	0.84
1:C:187:VAL:CB	1:C:227:MET:HE1	2.07	0.84
1:A:213:VAL:CG2	1:A:214:THR:N	2.01	0.84
1:C:117:ILE:HG21	1:C:345:GLY:HA3	1.60	0.84
1:A:385:PRO:O	1:A:388:LEU:CD1	2.26	0.83
1:D:336:VAL:HG12	1:D:342:MET:HE1	1.60	0.82
1:C:384:ILE:HG23	1:C:385:PRO:HD2	1.61	0.82
1:D:47:THR:CG2	1:D:52:ILE:CD1	2.55	0.82
1:A:124:ILE:HD11	1:B:8:THR:O	1.78	0.82
1:C:37:PRO:HB3	1:C:354:THR:CG2	2.08	0.82
1:C:28:ILE:HD11	1:C:66:ILE:HG23	1.63	0.81
1:B:205:ILE:O	1:B:209:LEU:HD21	1.78	0.81
1:B:41:VAL:HG13	1:B:58:ILE:HG22	1.63	0.80
1:A:335:ILE:N	1:A:335:ILE:CD1	2.45	0.79
1:A:200:VAL:HG11	1:A:205:ILE:HG12	1.63	0.79
1:D:47:THR:HG22	1:D:52:ILE:HD13	1.64	0.79
1:D:48:LYS:CG	1:D:49:THR:H	1.95	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:151:LEU:HD21	1:B:192:ILE:HD11	1.63	0.79
1:A:308:ALA:HB2	1:A:315:LEU:HD13	1.66	0.78
1:D:120:GLY:O	1:D:121:ASP:HB2	1.83	0.77
1:A:335:ILE:H	1:A:335:ILE:CD1	1.97	0.77
1:A:388:LEU:HD12	1:A:388:LEU:N	2.00	0.77
1:C:151:LEU:HD21	1:C:192:ILE:HD11	1.65	0.77
1:C:173:LEU:HD12	1:C:173:LEU:O	1.82	0.77
1:C:252:LYS:NZ	2:C:400:ACT:H3	2.00	0.77
1:D:209:LEU:O	1:D:213:VAL:HG22	1.85	0.77
1:D:151:LEU:HD21	1:D:192:ILE:HD11	1.67	0.77
1:C:147:LEU:HD12	1:C:217:VAL:HG11	1.66	0.76
1:B:384:ILE:HG13	1:B:385:PRO:HD2	1.67	0.76
1:A:41:VAL:HG13	1:A:58:ILE:HG22	1.68	0.76
1:D:48:LYS:CG	1:D:49:THR:N	2.48	0.76
1:D:188:HIS:ND1	1:D:189:THR:HG22	2.00	0.76
1:B:298:CYS:SG	1:B:301:VAL:CG2	2.73	0.76
1:B:385:PRO:CD	1:B:388:LEU:HD12	2.16	0.75
1:B:139:LEU:HA	1:B:142:VAL:HG23	1.68	0.75
1:D:181:HIS:HD1	1:D:299:HIS:HD1	0.75	0.75
1:C:61:ASN:OD1	1:C:102:GLN:NE2	2.19	0.75
1:A:172:THR:HG21	1:B:315:LEU:HD13	1.69	0.74
1:C:173:LEU:C	1:C:173:LEU:CD1	2.61	0.74
1:C:298:CYS:SG	1:C:301:VAL:CG2	2.76	0.74
1:B:178:VAL:O	1:B:179:GLU:CG	2.36	0.73
1:A:159:ILE:O	1:A:160:LEU:HD12	1.89	0.73
1:A:53:LEU:HD22	1:B:53:LEU:HD22	1.69	0.73
1:C:41:VAL:HG13	1:C:58:ILE:HG22	1.69	0.73
1:C:336:VAL:HG12	1:C:342:MET:HE1	1.69	0.73
1:D:332:LEU:HA	1:D:335:ILE:HD13	1.70	0.73
1:A:124:ILE:CD1	1:B:8:THR:O	2.36	0.72
1:C:188:HIS:ND1	1:C:189:THR:HG22	2.03	0.72
1:C:336:VAL:CG1	1:C:342:MET:CE	2.68	0.72
1:C:53:LEU:HD22	1:D:53:LEU:HD22	1.71	0.72
1:C:209:LEU:O	1:C:213:VAL:HG13	1.90	0.72
1:B:134:LYS:O	1:B:134:LYS:HD3	1.88	0.71
1:B:201:SER:O	1:B:205:ILE:HG23	1.90	0.71
1:B:322:TYR:O	1:B:324:THR:HG22	1.91	0.71
1:D:299:HIS:CD2	1:D:392:LEU:CD1	2.74	0.71
1:D:336:VAL:HG12	1:D:342:MET:CE	2.20	0.71
1:A:331:ILE:O	1:A:335:ILE:CD1	2.33	0.71
1:B:173:LEU:HD11	1:B:175:TYR:CE1	2.26	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:200:VAL:HG11	1:D:205:ILE:HD11	1.72	0.71
1:B:247:GLY:C	1:B:248:LEU:HD12	2.15	0.71
1:D:298:CYS:SG	1:D:301:VAL:CG2	2.78	0.70
1:C:28:ILE:CD1	1:C:66:ILE:HG23	2.21	0.70
1:B:82:THR:HB	1:D:82:THR:HG21	1.73	0.70
1:A:284:CYS:HB2	1:A:347:ILE:HD12	1.72	0.70
1:C:28:ILE:HD12	1:C:66:ILE:HD13	1.72	0.70
1:C:173:LEU:HD11	1:C:175:TYR:CD1	2.26	0.70
1:D:384:ILE:CG2	1:D:388:LEU:HD12	2.22	0.70
1:C:117:ILE:HG21	1:C:345:GLY:N	2.07	0.70
1:B:49:THR:CG2	1:B:50:GLY:N	2.54	0.70
1:B:140:THR:O	1:B:144:SER:OG	2.10	0.70
1:C:174:GLU:HG2	1:C:185:ILE:HG23	1.72	0.70
1:B:49:THR:HG22	1:B:50:GLY:H	1.56	0.69
1:D:355:ARG:O	1:D:357:ILE:HD13	1.92	0.69
1:D:326:ILE:HD12	1:D:327:CYS:N	2.07	0.69
1:C:140:THR:O	1:C:144:SER:OG	2.11	0.69
1:A:213:VAL:HG23	1:A:214:THR:CA	2.21	0.69
1:C:326:ILE:CD1	1:C:327:CYS:N	2.48	0.69
1:D:47:THR:HG22	1:D:52:ILE:HA	1.74	0.69
1:A:188:HIS:ND1	1:A:189:THR:HG22	2.08	0.69
1:C:355:ARG:NH2	1:C:373:ASP:OD1	2.26	0.69
1:A:123:GLY:C	1:A:125:MET:HE3	2.18	0.68
1:C:160:LEU:HD22	1:C:162:TRP:CZ2	2.27	0.68
1:B:66:ILE:O	1:B:69:ILE:HG23	1.92	0.68
1:D:200:VAL:HG21	1:D:205:ILE:HG12	1.76	0.68
1:A:284:CYS:SG	1:A:347:ILE:HD11	2.33	0.68
1:D:140:THR:O	1:D:144:SER:OG	2.10	0.68
1:B:81:ASP:HB2	1:B:84:LYS:HG3	1.75	0.68
1:B:82:THR:OG1	1:D:90:THR:HG21	1.94	0.68
1:A:140:THR:O	1:A:144:SER:OG	2.11	0.67
1:A:48:LYS:O	1:A:242:PRO:O	2.12	0.67
1:A:284:CYS:SG	1:A:347:ILE:CD1	2.82	0.67
1:C:317:ILE:HD12	1:C:318:ASN:N	2.07	0.67
1:A:57:GLU:OE2	1:B:244:GLY:O	2.11	0.67
1:B:63:VAL:HG11	1:C:63:VAL:CG1	2.23	0.67
1:B:139:LEU:HA	1:B:142:VAL:CG2	2.25	0.67
1:B:48:LYS:O	1:B:242:PRO:O	2.13	0.66
1:C:392:LEU:HD23	1:C:393:LEU:N	2.05	0.66
1:A:151:LEU:HD23	1:A:163:LEU:HD21	1.78	0.66
1:C:55:LEU:HD11	1:D:48:LYS:HB2	1.76	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:385:PRO:HD2	1:D:388:LEU:HD12	1.77	0.66
1:C:28:ILE:CD1	1:C:66:ILE:HD13	2.26	0.66
1:C:317:ILE:HD11	1:C:333:VAL:HG13	1.76	0.66
1:B:349:LYS:HG3	1:B:350:GLU:N	2.10	0.66
1:B:27:ALA:C	1:B:69:ILE:HD11	2.21	0.66
1:B:213:VAL:HG23	1:B:214:THR:N	2.11	0.66
1:A:12:VAL:C	1:A:148:ILE:CD1	2.69	0.65
1:B:160:LEU:HD22	1:B:162:TRP:CZ2	2.30	0.65
1:C:336:VAL:HG12	1:C:342:MET:CE	2.26	0.65
1:C:298:CYS:SG	1:C:301:VAL:HG22	2.36	0.65
1:A:94:LEU:HD21	1:C:51:LEU:HD13	1.79	0.65
1:D:148:ILE:CD1	1:D:169:SER:OG	2.44	0.65
1:C:288:ILE:CD1	1:C:342:MET:HE2	2.24	0.65
1:B:28:ILE:HD11	1:B:66:ILE:HG23	1.79	0.64
1:C:187:VAL:CG1	1:C:227:MET:HE1	2.27	0.64
1:C:187:VAL:O	1:C:227:MET:HE2	1.97	0.64
1:C:85:GLY:HA2	1:C:243:MET:HE3	1.80	0.64
1:C:87:ASP:CG	1:C:90:THR:HG23	2.23	0.64
1:C:384:ILE:CG2	1:C:385:PRO:HD2	2.28	0.64
1:A:308:ALA:CB	1:A:315:LEU:HD13	2.27	0.64
1:D:215:GLN:O	1:D:220:LYS:NZ	2.27	0.64
1:B:297:LEU:HD11	1:B:335:ILE:CD1	2.28	0.63
1:B:6:PHE:CE2	1:B:174:GLU:HG3	2.32	0.63
1:C:363:VAL:O	1:C:369:ARG:NH1	2.31	0.63
1:B:289:ALA:O	1:B:293:VAL:HG12	1.99	0.63
1:C:174:GLU:HG2	1:C:185:ILE:HG12	1.81	0.63
1:A:326:ILE:HD12	1:A:327:CYS:N	2.14	0.62
1:B:190:ILE:HD12	1:B:227:MET:CE	2.29	0.62
1:C:289:ALA:O	1:C:293:VAL:HG12	1.98	0.62
1:B:148:ILE:HA	1:B:151:LEU:HD12	1.81	0.62
1:B:9:SER:HB2	1:B:140:THR:HG23	1.79	0.62
1:A:289:ALA:O	1:A:293:VAL:HG12	1.98	0.62
1:B:177:GLU:O	1:B:177:GLU:HG3	1.97	0.62
1:C:253:ILE:HD12	1:C:264:HIS:CE1	2.34	0.62
1:C:328:ASP:O	1:C:331:ILE:HG22	1.99	0.62
1:D:329:GLU:O	1:D:332:LEU:HD23	2.00	0.62
1:A:162:TRP:HB2	1:A:200:VAL:HG21	1.80	0.61
1:B:139:LEU:CA	1:B:142:VAL:HG23	2.31	0.61
1:D:384:ILE:HG23	1:D:388:LEU:HD12	1.82	0.61
1:A:162:TRP:HB2	1:A:200:VAL:CG2	2.31	0.61
1:C:252:LYS:HZ1	2:C:400:ACT:H3	1.63	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:304:GLN:HE22	1:D:8:THR:H	1.48	0.61
1:D:389:LYS:CB	1:D:390:PRO:HD3	2.30	0.61
1:D:253:ILE:HD12	1:D:264:HIS:CE1	2.35	0.61
1:B:160:LEU:HD22	1:B:162:TRP:CE2	2.36	0.61
1:D:159:ILE:O	1:D:160:LEU:HD12	2.01	0.61
1:B:41:VAL:HG13	1:B:58:ILE:CG2	2.29	0.61
1:D:42:ALA:HB2	1:D:276:LYS:HE2	1.83	0.61
1:A:41:VAL:HG13	1:A:58:ILE:CG2	2.31	0.60
1:B:80:ASP:OD1	1:B:84:LYS:NZ	2.34	0.60
1:B:332:LEU:HA	1:B:335:ILE:HG23	1.82	0.60
1:C:41:VAL:HG13	1:C:58:ILE:CG2	2.31	0.60
1:C:140:THR:HG21	1:C:261:TRP:HE1	1.67	0.60
1:C:321:THR:HG21	1:C:325:GLY:N	2.17	0.60
1:C:301:VAL:HG12	1:C:302:LEU:N	2.15	0.60
1:B:321:THR:HG21	1:B:325:GLY:N	2.17	0.60
1:B:257:THR:HG21	1:B:264:HIS:HE2	1.67	0.60
1:D:119:ALA:HB1	1:D:277:VAL:HG11	1.83	0.60
1:C:55:LEU:HD11	1:D:48:LYS:HB3	1.82	0.59
1:C:12:VAL:HG12	1:C:168:LYS:HG3	1.83	0.59
1:C:42:ALA:HB2	1:C:276:LYS:HE2	1.84	0.59
1:A:140:THR:HG21	1:A:261:TRP:HE1	1.67	0.59
1:C:148:ILE:HA	1:C:151:LEU:HD12	1.84	0.59
1:A:361:THR:HG23	1:A:366:HIS:CE1	2.36	0.59
1:D:9:SER:HB2	1:D:140:THR:HG23	1.83	0.59
1:D:293:VAL:O	1:D:293:VAL:CG2	2.50	0.59
1:D:147:LEU:HD12	1:D:217:VAL:CG2	2.24	0.59
1:B:42:ALA:HB2	1:B:276:LYS:HE2	1.85	0.59
1:B:178:VAL:O	1:B:179:GLU:HG2	2.03	0.59
1:C:252:LYS:HZ2	2:C:400:ACT:H3	1.68	0.59
1:D:140:THR:HG21	1:D:261:TRP:HE1	1.68	0.59
1:A:253:ILE:HD12	1:A:264:HIS:CE1	2.37	0.59
1:A:13:SER:N	1:A:148:ILE:HD13	2.17	0.59
1:B:83:ASN:O	1:B:239:ILE:HG21	2.03	0.59
1:B:213:VAL:HG23	1:B:214:THR:H	1.68	0.59
1:C:279:ARG:HH21	1:C:361:THR:HG21	1.68	0.59
1:A:279:ARG:HH21	1:A:361:THR:HG21	1.67	0.58
1:B:173:LEU:C	1:B:173:LEU:HD12	2.28	0.58
1:C:147:LEU:HD12	1:C:217:VAL:CG1	2.32	0.58
1:C:361:THR:HG23	1:C:366:HIS:CE1	2.38	0.58
1:A:44:GLU:OE2	1:B:248:LEU:HD13	2.03	0.58
1:B:65:ASP:O	1:B:69:ILE:HG22	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:297:LEU:O	1:B:326:ILE:HG22	2.04	0.58
1:A:207:LYS:O	1:A:211:GLU:HG3	2.04	0.58
1:C:174:GLU:HG2	1:C:185:ILE:CG2	2.33	0.58
1:C:189:THR:OG1	1:D:315:LEU:HD11	2.03	0.58
1:D:293:VAL:O	1:D:293:VAL:HG22	2.03	0.58
1:A:48:LYS:CG	1:A:49:THR:N	2.62	0.58
1:A:55:LEU:C	1:A:55:LEU:HD12	2.29	0.58
1:C:382:LEU:N	1:C:382:LEU:CD1	2.67	0.58
1:B:190:ILE:HD12	1:B:227:MET:HE1	1.85	0.58
1:C:384:ILE:HG22	1:C:385:PRO:O	2.04	0.58
1:D:11:SER:HB3	1:D:148:ILE:HD13	1.85	0.57
1:C:9:SER:HB2	1:C:140:THR:HG23	1.85	0.57
1:D:279:ARG:HH21	1:D:361:THR:HG21	1.67	0.57
1:B:253:ILE:HD12	1:B:264:HIS:CE1	2.40	0.57
1:D:363:VAL:O	1:D:369:ARG:NH2	2.37	0.57
1:A:9:SER:HB2	1:A:140:THR:HG23	1.85	0.57
1:C:391:LYS:O	1:C:392:LEU:HB2	2.03	0.57
1:C:392:LEU:CD2	1:C:393:LEU:H	2.11	0.57
1:D:160:LEU:HD23	1:D:162:TRP:CZ2	2.39	0.57
1:B:213:VAL:O	1:B:217:VAL:CG1	2.49	0.57
1:D:361:THR:HG23	1:D:366:HIS:CE1	2.39	0.57
1:A:94:LEU:HD21	1:C:51:LEU:CD1	2.34	0.56
1:A:124:ILE:N	1:A:125:MET:HE3	2.19	0.56
1:B:140:THR:HG21	1:B:261:TRP:HE1	1.70	0.56
1:C:116:ASP:C	1:C:117:ILE:HG22	2.30	0.56
1:D:198:ASP:OD1	1:D:236:ARG:NH2	2.38	0.56
1:B:339:ASN:ND2	1:B:382:LEU:HB3	2.20	0.56
1:C:37:PRO:CB	1:C:354:THR:HG23	2.30	0.56
1:C:139:LEU:HD22	1:C:184:PRO:HG3	1.87	0.56
1:C:87:ASP:HB3	1:C:90:THR:HG23	1.88	0.56
1:C:336:VAL:HG11	1:C:342:MET:HE1	1.87	0.56
1:B:198:ASP:OD1	1:B:236:ARG:NH2	2.38	0.56
1:D:57:GLU:C	1:D:58:ILE:CG2	2.79	0.56
1:B:243:MET:HE1	1:D:90:THR:HB	1.88	0.56
1:D:358:PHE:O	1:D:361:THR:HG22	2.05	0.56
1:A:331:ILE:HG22	1:A:335:ILE:HD11	1.87	0.56
1:D:339:ASN:O	1:D:380:LYS:NZ	2.37	0.56
1:B:329:GLU:O	1:B:333:VAL:HG23	2.06	0.56
1:B:134:LYS:HD3	1:B:134:LYS:C	2.30	0.55
1:A:42:ALA:HB2	1:A:276:LYS:HE2	1.88	0.55
1:D:181:HIS:CG	1:D:299:HIS:HD1	2.14	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:12:VAL:HG12	1:A:168:LYS:HG3	1.88	0.55
1:B:178:VAL:O	1:B:179:GLU:HG3	2.05	0.55
1:A:199:ASN:C	1:A:199:ASN:OD1	2.49	0.55
1:A:48:LYS:HG2	1:A:49:THR:N	2.20	0.55
1:D:70:VAL:O	1:D:74:VAL:HG23	2.07	0.55
1:A:357:ILE:HG21	1:A:373:ASP:HB3	1.88	0.55
1:A:53:LEU:HD12	1:A:94:LEU:HB2	1.88	0.55
1:A:116:ASP:OD1	1:A:116:ASP:C	2.50	0.55
1:A:223:MET:HG3	1:A:227:MET:HE3	1.88	0.55
1:B:202:ASN:O	1:B:205:ILE:HG13	2.07	0.55
1:B:139:LEU:HD22	1:B:184:PRO:HG3	1.87	0.55
1:C:117:ILE:HG21	1:C:345:GLY:H	1.70	0.55
1:A:124:ILE:O	1:A:265:GLY:HA3	2.06	0.55
1:A:12:VAL:C	1:A:148:ILE:HD13	2.32	0.54
1:A:247:GLY:C	1:A:248:LEU:HD12	2.32	0.54
1:B:213:VAL:CG2	1:B:214:THR:H	2.20	0.54
1:B:82:THR:HB	1:D:82:THR:CG2	2.37	0.54
1:B:326:ILE:O	1:B:326:ILE:HG13	2.08	0.54
1:C:354:THR:CG2	1:C:354:THR:O	2.53	0.54
1:D:148:ILE:HA	1:D:151:LEU:HD12	1.89	0.54
1:A:321:THR:HG21	1:A:325:GLY:N	2.23	0.54
1:B:385:PRO:HD2	1:B:388:LEU:HD12	1.89	0.54
1:A:378:PHE:CD1	1:A:378:PHE:N	2.75	0.54
1:A:382:LEU:N	1:A:382:LEU:CD1	2.71	0.54
1:A:339:ASN:ND2	1:A:382:LEU:HB3	2.23	0.54
1:C:127:GLY:O	1:C:302:LEU:HD12	2.07	0.54
1:B:124:ILE:HG23	1:B:265:GLY:CA	2.38	0.54
1:D:391:LYS:HG3	1:D:392:LEU:N	2.10	0.54
1:C:300:ARG:O	1:C:321:THR:HG23	2.08	0.54
1:D:53:LEU:HD12	1:D:94:LEU:HB2	1.89	0.54
1:D:181:HIS:ND1	1:D:299:HIS:CE1	2.74	0.54
1:A:82:THR:HG21	1:C:82:THR:CG2	2.38	0.53
1:A:213:VAL:HG22	1:A:214:THR:H	1.57	0.53
1:C:116:ASP:O	1:C:117:ILE:HG22	2.08	0.53
1:C:117:ILE:O	1:C:117:ILE:CG2	2.41	0.53
1:D:300:ARG:O	1:D:321:THR:HG23	2.08	0.53
1:D:336:VAL:CG1	1:D:342:MET:HE1	2.36	0.53
1:A:139:LEU:O	1:A:143:LEU:HD12	2.08	0.53
1:A:139:LEU:HD22	1:A:184:PRO:HG3	1.89	0.53
1:A:82:THR:CG2	1:C:82:THR:HG21	2.39	0.53
1:A:358:PHE:O	1:A:361:THR:HG22	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:335:ILE:N	1:A:335:ILE:HD12	2.21	0.53
1:C:358:PHE:O	1:C:361:THR:HG22	2.08	0.53
1:D:213:VAL:HG23	1:D:214:THR:H	1.73	0.53
1:B:162:TRP:CB	1:B:200:VAL:HG11	2.39	0.53
1:B:304:GLN:O	1:B:304:GLN:HG2	2.08	0.53
1:C:205:ILE:CG2	1:C:209:LEU:HD22	2.38	0.53
1:A:162:TRP:CB	1:A:200:VAL:CG2	2.86	0.53
1:C:308:ALA:HB2	1:C:315:LEU:HD22	1.90	0.53
1:D:351:LEU:HD13	1:D:376:TRP:HB3	1.91	0.53
1:A:284:CYS:CB	1:A:347:ILE:HD12	2.38	0.53
1:B:53:LEU:HD12	1:B:94:LEU:HB2	1.91	0.53
1:A:162:TRP:CB	1:A:200:VAL:HG21	2.39	0.52
1:B:321:THR:HG21	1:B:325:GLY:CA	2.39	0.52
1:C:124:ILE:HG23	1:C:265:GLY:CA	2.39	0.52
1:C:268:ALA:HB3	1:D:251:ARG:CZ	2.39	0.52
1:D:24:ILE:O	1:D:28:ILE:HD12	2.09	0.52
1:D:321:THR:HG21	1:D:325:GLY:N	2.25	0.52
1:D:332:LEU:HA	1:D:335:ILE:CD1	2.37	0.52
1:B:382:LEU:N	1:B:382:LEU:CD1	2.73	0.52
1:D:173:LEU:CD1	1:D:175:TYR:CD1	2.88	0.52
1:D:296:GLY:O	1:D:392:LEU:HB2	2.09	0.52
1:D:384:ILE:CG2	1:D:385:PRO:HD2	2.39	0.52
1:D:57:GLU:O	1:D:58:ILE:HG22	2.10	0.52
1:A:172:THR:HG21	1:B:315:LEU:CD1	2.38	0.52
1:A:284:CYS:SG	1:A:347:ILE:HD12	2.49	0.52
1:A:335:ILE:HG22	1:A:339:ASN:ND2	2.25	0.52
1:B:213:VAL:CG2	1:B:214:THR:N	2.72	0.52
1:C:317:ILE:CG1	1:C:333:VAL:HG13	2.40	0.52
1:D:124:ILE:HG23	1:D:265:GLY:CA	2.40	0.52
1:D:209:LEU:O	1:D:213:VAL:CG2	2.56	0.52
1:B:296:GLY:O	1:B:392:LEU:HB2	2.10	0.52
1:D:213:VAL:HG23	1:D:214:THR:N	2.25	0.52
1:B:102:GLN:O	1:B:103:ASP:C	2.53	0.51
1:A:300:ARG:O	1:A:321:THR:HG23	2.09	0.51
1:B:209:LEU:HB2	1:B:231:TYR:CE1	2.46	0.51
1:C:173:LEU:CD1	1:C:175:TYR:CD1	2.93	0.51
1:D:299:HIS:HD2	1:D:392:LEU:CD1	2.12	0.51
1:D:391:LYS:CG	1:D:392:LEU:H	2.07	0.51
1:B:124:ILE:HG23	1:B:265:GLY:HA2	1.93	0.51
1:D:200:VAL:CG2	1:D:205:ILE:HG12	2.40	0.51
1:A:29:LEU:HD13	1:A:41:VAL:HB	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:66:ILE:HA	1:B:69:ILE:CG2	2.41	0.51
1:A:288:ILE:HD13	1:A:336:VAL:HG13	1.92	0.51
1:C:187:VAL:CB	1:C:227:MET:CE	2.77	0.51
1:C:331:ILE:CG2	1:C:332:LEU:N	2.74	0.51
1:B:81:ASP:C	1:B:83:ASN:N	2.63	0.50
1:C:205:ILE:HG22	1:C:209:LEU:HD22	1.93	0.50
1:C:315:LEU:HD11	1:D:230:TYR:CE2	2.46	0.50
1:C:357:ILE:HG21	1:C:373:ASP:HB3	1.93	0.50
1:C:187:VAL:CG1	1:C:227:MET:CE	2.89	0.50
1:D:148:ILE:HD12	1:D:169:SER:OG	2.11	0.50
1:C:317:ILE:HD12	1:C:317:ILE:C	2.36	0.50
1:D:192:ILE:HD13	1:D:209:LEU:HD22	1.93	0.50
1:A:124:ILE:C	1:A:125:MET:HG3	2.37	0.50
1:B:209:LEU:O	1:B:213:VAL:CG2	2.53	0.50
1:C:117:ILE:CG2	1:C:345:GLY:HA3	2.37	0.50
1:A:162:TRP:CG	1:A:200:VAL:HG21	2.47	0.50
1:B:210:GLU:OE2	1:B:215:GLN:HG2	2.11	0.50
1:D:8:THR:HG23	1:D:172:THR:HG22	1.93	0.49
1:D:218:ILE:HG22	1:D:223:MET:HG3	1.94	0.49
1:D:384:ILE:HG23	1:D:385:PRO:HD2	1.94	0.49
1:C:321:THR:HG21	1:C:325:GLY:CA	2.42	0.49
1:D:139:LEU:HD22	1:D:184:PRO:HG3	1.93	0.49
1:B:11:SER:HA	1:B:255:VAL:HG22	1.94	0.49
1:D:227:MET:HE2	1:D:229:ARG:HH21	1.77	0.49
1:A:20:MET:HE1	1:A:47:THR:HG21	1.94	0.49
1:B:332:LEU:HA	1:B:335:ILE:CG2	2.42	0.49
1:C:124:ILE:HG23	1:C:265:GLY:HA2	1.95	0.49
1:A:316:SER:O	1:A:317:ILE:HD13	2.13	0.49
1:B:146:LYS:HA	1:B:149:LEU:HD12	1.94	0.49
1:B:59:THR:HG22	1:B:102:GLN:HG3	1.95	0.49
1:B:61:ASN:OD1	1:B:102:GLN:HB2	2.13	0.49
1:B:331:ILE:O	1:B:332:LEU:C	2.56	0.49
1:C:224:ASP:OD2	1:C:226:LYS:CB	2.45	0.49
1:C:317:ILE:CD1	1:C:333:VAL:HG13	2.42	0.49
1:D:57:GLU:C	1:D:58:ILE:HG22	2.38	0.49
1:D:218:ILE:CG2	1:D:223:MET:HG3	2.43	0.49
1:B:300:ARG:O	1:B:321:THR:HG23	2.13	0.48
1:D:119:ALA:HB1	1:D:277:VAL:CG1	2.43	0.48
1:A:62:ALA:HB1	1:A:64:ILE:CD1	2.43	0.48
1:B:297:LEU:HD11	1:B:335:ILE:HD11	1.95	0.48
1:B:391:LYS:O	1:B:392:LEU:HD12	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:148:ILE:HD11	1:D:169:SER:CB	2.42	0.48
1:A:118:GLY:O	1:A:119:ALA:C	2.57	0.48
1:A:159:ILE:O	1:A:160:LEU:CD1	2.59	0.48
1:A:173:LEU:HD23	1:A:175:TYR:CE1	2.49	0.48
1:B:332:LEU:O	1:B:335:ILE:HG23	2.13	0.48
1:C:302:LEU:O	1:C:302:LEU:HG	2.11	0.48
1:D:189:THR:HB	1:D:228:LEU:HB2	1.94	0.48
1:C:69:ILE:O	1:C:73:VAL:HG23	2.13	0.48
1:C:354:THR:HG23	1:C:354:THR:O	2.12	0.48
1:A:213:VAL:O	1:A:214:THR:C	2.56	0.48
1:B:147:LEU:HD12	1:B:217:VAL:HG21	1.94	0.48
1:B:297:LEU:HD11	1:B:335:ILE:HD13	1.94	0.48
1:C:87:ASP:CG	1:C:90:THR:CG2	2.86	0.48
1:A:339:ASN:O	1:A:380:LYS:NZ	2.42	0.48
1:A:388:LEU:CD1	1:A:388:LEU:N	2.73	0.48
1:C:160:LEU:HD22	1:C:162:TRP:CE2	2.48	0.48
1:D:357:ILE:HG21	1:D:373:ASP:HB3	1.94	0.48
1:D:385:PRO:HD2	1:D:388:LEU:CD1	2.43	0.48
1:C:321:THR:HG21	1:C:325:GLY:H	1.79	0.48
1:A:12:VAL:HA	1:A:148:ILE:HD11	1.95	0.48
1:B:299:HIS:CD2	1:B:392:LEU:HD23	2.49	0.48
1:B:315:LEU:HD12	1:B:315:LEU:C	2.38	0.48
1:B:381:GLU:C	1:B:382:LEU:HD12	2.39	0.48
1:D:178:VAL:HG12	1:D:178:VAL:O	2.13	0.48
1:B:384:ILE:CG1	1:B:385:PRO:HD2	2.40	0.48
1:D:296:GLY:O	1:D:391:LYS:HB2	2.14	0.48
1:D:58:ILE:HG12	1:D:58:ILE:O	2.12	0.47
1:A:62:ALA:HB1	1:A:64:ILE:HD13	1.96	0.47
1:A:248:LEU:HD13	1:B:44:GLU:OE1	2.14	0.47
1:C:224:ASP:HB3	1:C:226:LYS:H	1.79	0.47
1:B:391:LYS:C	1:B:392:LEU:HD12	2.39	0.47
1:C:8:THR:HG23	1:C:172:THR:HG22	1.97	0.47
1:C:87:ASP:CB	1:C:90:THR:HG23	2.43	0.47
1:C:253:ILE:HD12	1:C:264:HIS:NE2	2.29	0.47
1:D:237:PHE:CD1	1:D:237:PHE:C	2.92	0.47
1:B:384:ILE:HG12	1:B:385:PRO:O	2.15	0.47
1:A:151:LEU:HD11	1:A:192:ILE:CD1	2.44	0.47
1:C:391:LYS:O	1:C:392:LEU:CB	2.63	0.47
1:D:233:PRO:C	1:D:235:GLY:H	2.23	0.47
1:D:253:ILE:HD12	1:D:264:HIS:NE2	2.30	0.47
1:B:382:LEU:N	1:B:382:LEU:HD12	2.29	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:115:GLU:O	1:A:345:GLY:HA3	2.15	0.47
1:A:140:THR:CG2	1:A:261:TRP:HE1	2.28	0.47
1:B:186:ARG:HA	1:B:222:LEU:HD23	1.97	0.47
1:D:124:ILE:HG23	1:D:265:GLY:HA2	1.96	0.47
1:D:148:ILE:CD1	1:D:169:SER:CB	2.93	0.47
1:D:237:PHE:C	1:D:237:PHE:HD1	2.23	0.47
1:D:332:LEU:HD23	1:D:332:LEU:H	1.80	0.47
1:A:279:ARG:NH2	1:A:361:THR:HG21	2.30	0.47
1:B:9:SER:CB	1:B:140:THR:HG23	2.42	0.47
1:C:121:ASP:OD1	1:C:122:GLN:N	2.48	0.47
1:D:392:LEU:HA	1:D:392:LEU:HD23	1.69	0.47
1:A:233:PRO:C	1:A:235:GLY:H	2.23	0.47
1:B:357:ILE:HG21	1:B:373:ASP:HB3	1.97	0.47
1:D:62:ALA:HB1	1:D:64:ILE:CD1	2.45	0.47
1:B:81:ASP:OD1	1:B:83:ASN:OD1	2.33	0.46
1:D:279:ARG:NH2	1:D:361:THR:HG21	2.30	0.46
1:A:315:LEU:HD11	1:B:230:TYR:CE2	2.50	0.46
1:C:49:THR:HG22	1:C:50:GLY:N	2.30	0.46
1:C:182:LEU:CD1	1:C:300:ARG:HD2	2.44	0.46
1:D:125:MET:HE2	1:D:278:ASP:HA	1.97	0.46
1:D:162:TRP:CZ3	1:D:205:ILE:HD13	2.50	0.46
1:D:391:LYS:CG	1:D:392:LEU:N	2.75	0.46
1:D:148:ILE:HD11	1:D:169:SER:OG	2.15	0.46
1:A:12:VAL:CA	1:A:148:ILE:CD1	2.92	0.46
1:A:132:GLU:OE2	1:A:175:TYR:OH	2.24	0.46
1:B:48:LYS:CG	1:B:49:THR:N	2.75	0.46
1:B:79:TYR:HB3	1:B:86:PHE:O	2.15	0.46
1:C:321:THR:HG21	1:C:325:GLY:HA3	1.96	0.46
1:D:8:THR:OG1	1:D:172:THR:HB	2.16	0.46
1:C:335:ILE:HD11	1:C:385:PRO:HD3	1.97	0.46
1:D:9:SER:CB	1:D:140:THR:HG23	2.46	0.46
1:A:311:VAL:HG21	1:B:230:TYR:CE1	2.50	0.46
1:A:372:PRO:HA	1:A:378:PHE:HZ	1.81	0.46
1:B:139:LEU:C	1:B:142:VAL:HG23	2.41	0.46
1:B:288:ILE:HD13	1:B:336:VAL:HG13	1.98	0.46
1:A:121:ASP:OD1	1:A:122:GLN:N	2.48	0.46
1:B:343:ARG:HH21	1:B:346:MET:HE3	1.80	0.46
1:C:329:GLU:O	1:C:333:VAL:HG23	2.15	0.46
1:D:48:LYS:HG3	1:D:49:THR:H	1.57	0.46
1:A:82:THR:HG21	1:C:82:THR:HG21	1.97	0.46
1:C:233:PRO:C	1:C:235:GLY:H	2.24	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:284:CYS:SG	1:D:347:ILE:HD13	2.56	0.46
1:B:81:ASP:CB	1:B:84:LYS:HG3	2.42	0.46
1:B:201:SER:HG	1:B:204:GLU:H	1.59	0.46
1:C:125:MET:HE2	1:C:278:ASP:HA	1.97	0.46
1:C:326:ILE:HD12	1:C:327:CYS:CA	2.42	0.46
1:C:339:ASN:HD21	1:C:383:GLU:HG2	1.81	0.45
1:B:358:PHE:HA	1:B:361:THR:OG1	2.16	0.45
1:C:186:ARG:NH1	1:C:226:LYS:O	2.49	0.45
1:B:81:ASP:C	1:B:83:ASN:H	2.24	0.45
1:B:147:LEU:O	1:B:151:LEU:HD12	2.16	0.45
1:B:233:PRO:C	1:B:235:GLY:H	2.23	0.45
1:C:227:MET:HE2	1:C:227:MET:HB2	1.68	0.45
1:C:9:SER:CB	1:C:140:THR:HG23	2.47	0.45
1:C:218:ILE:CG2	1:C:223:MET:HG3	2.46	0.45
1:D:326:ILE:HD12	1:D:326:ILE:C	2.41	0.45
1:A:219:PRO:HG2	1:A:222:LEU:HD22	1.99	0.45
1:B:8:THR:OG1	1:B:172:THR:HB	2.17	0.45
1:C:8:THR:HA	1:C:172:THR:HA	1.99	0.45
1:A:243:MET:HE1	1:C:89:GLN:O	2.17	0.45
1:B:274:SER:O	1:B:280:SER:OG	2.33	0.45
1:B:12:VAL:HG12	1:B:168:LYS:HG3	1.99	0.45
1:B:247:GLY:O	1:B:248:LEU:HD12	2.16	0.45
1:A:151:LEU:HD11	1:A:192:ILE:HD12	1.99	0.45
1:A:298:CYS:SG	1:A:301:VAL:HG12	2.57	0.45
1:B:322:TYR:C	1:B:324:THR:CG2	2.89	0.45
1:B:237:PHE:HE1	1:B:240:GLY:HA3	1.82	0.44
1:C:8:THR:OG1	1:C:172:THR:HB	2.17	0.44
1:D:29:LEU:HD13	1:D:41:VAL:HB	1.98	0.44
1:D:284:CYS:HB2	1:D:347:ILE:HD13	1.98	0.44
1:D:120:GLY:O	1:D:121:ASP:CB	2.56	0.44
1:D:140:THR:CG2	1:D:261:TRP:HE1	2.30	0.44
1:D:201:SER:O	1:D:204:GLU:N	2.50	0.44
1:A:9:SER:CB	1:A:140:THR:HG23	2.47	0.44
1:C:174:GLU:CG	1:C:185:ILE:HG23	2.42	0.44
1:C:304:GLN:O	1:C:304:GLN:HG2	2.14	0.44
1:C:321:THR:CG2	1:C:325:GLY:H	2.30	0.44
1:A:327:CYS:HG	1:A:388:LEU:CD2	2.23	0.44
1:B:63:VAL:CG1	1:C:63:VAL:CG1	2.94	0.44
1:B:85:GLY:HA2	1:B:243:MET:HE3	2.00	0.44
1:C:140:THR:CG2	1:C:261:TRP:HE1	2.28	0.44
1:C:279:ARG:NH2	1:C:361:THR:HG21	2.31	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:209:LEU:C	1:B:213:VAL:HG22	2.36	0.44
1:B:210:GLU:OE2	1:B:215:GLN:CG	2.65	0.44
1:C:218:ILE:HG22	1:C:223:MET:HG3	1.98	0.44
1:A:253:ILE:HD12	1:A:264:HIS:NE2	2.33	0.44
1:A:321:THR:HG21	1:A:325:GLY:CA	2.48	0.44
1:D:326:ILE:HD12	1:D:327:CYS:CA	2.47	0.44
1:A:8:THR:OG1	1:A:172:THR:HB	2.18	0.44
1:C:139:LEU:HD21	1:C:222:LEU:HD21	2.00	0.44
1:D:62:ALA:HB1	1:D:64:ILE:HD13	2.00	0.44
1:A:124:ILE:HD12	1:B:8:THR:O	2.17	0.43
1:B:90:THR:CB	1:D:243:MET:HE1	2.38	0.43
1:B:140:THR:CG2	1:B:261:TRP:HE1	2.30	0.43
1:B:257:THR:HG21	1:B:264:HIS:NE2	2.31	0.43
1:A:55:LEU:HD11	1:B:246:ALA:HB1	2.00	0.43
1:A:200:VAL:CG1	1:A:205:ILE:HG12	2.41	0.43
1:B:132:GLU:OE2	1:B:175:TYR:OH	2.24	0.43
1:B:178:VAL:O	1:B:178:VAL:HG23	2.17	0.43
1:D:48:LYS:O	1:D:49:THR:C	2.59	0.43
1:A:19:LYS:NZ	1:A:256:ASP:OD1	2.51	0.43
1:A:124:ILE:HD11	1:A:304:GLN:NE2	2.32	0.43
1:A:214:THR:HG22	1:A:223:MET:HE1	2.00	0.43
1:B:125:MET:HE2	1:B:278:ASP:HA	1.98	0.43
1:D:8:THR:HA	1:D:172:THR:HA	1.99	0.43
1:D:132:GLU:OE2	1:D:175:TYR:OH	2.24	0.43
1:A:389:LYS:N	1:A:390:PRO:CD	2.81	0.43
1:B:8:THR:HA	1:B:172:THR:HA	2.00	0.43
1:C:55:LEU:HD13	1:C:96:CYS:HB3	2.01	0.43
1:C:382:LEU:N	1:C:382:LEU:HD12	2.32	0.43
1:D:200:VAL:O	1:D:200:VAL:HG13	2.17	0.43
1:A:246:ALA:HB1	1:B:55:LEU:CD1	2.49	0.43
1:A:335:ILE:HD13	1:A:335:ILE:N	2.08	0.43
1:C:62:ALA:HB1	1:C:64:ILE:CD1	2.49	0.43
1:D:331:ILE:HD12	1:D:331:ILE:N	2.34	0.43
1:A:346:MET:N	1:A:346:MET:HE2	2.34	0.43
1:C:308:ALA:HB2	1:C:315:LEU:HD13	2.01	0.43
1:D:126:PHE:CE1	1:D:304:GLN:HB2	2.53	0.43
1:D:215:GLN:CD	1:D:223:MET:HE1	2.43	0.43
1:B:218:ILE:CG2	1:B:223:MET:HE3	2.49	0.43
1:B:383:GLU:O	1:B:383:GLU:HG2	2.18	0.43
1:A:8:THR:HA	1:A:172:THR:HA	2.01	0.43
1:A:381:GLU:C	1:A:382:LEU:HD12	2.44	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:139:LEU:O	1:B:142:VAL:HG23	2.19	0.43
1:C:274:SER:O	1:C:280:SER:OG	2.33	0.43
1:A:82:THR:HG23	1:C:82:THR:HG21	2.01	0.42
1:D:6:PHE:N	1:D:6:PHE:CD2	2.86	0.42
1:A:248:LEU:HD23	1:B:248:LEU:CD2	2.31	0.42
1:C:381:GLU:C	1:C:382:LEU:HD12	2.44	0.42
1:D:12:VAL:HG12	1:D:168:LYS:HG3	2.00	0.42
1:D:47:THR:HG22	1:D:52:ILE:CD1	2.40	0.42
1:A:189:THR:HB	1:A:228:LEU:HB2	2.00	0.42
1:B:162:TRP:HB3	1:B:200:VAL:HG11	2.01	0.42
1:B:218:ILE:HG21	1:B:223:MET:HE3	2.00	0.42
1:A:82:THR:HG21	1:C:82:THR:HG23	2.00	0.42
1:A:384:ILE:HG22	1:A:388:LEU:HD13	2.00	0.42
1:B:322:TYR:C	1:B:324:THR:HG23	2.45	0.42
1:C:62:ALA:HB1	1:C:64:ILE:HD13	2.02	0.42
1:D:274:SER:O	1:D:280:SER:OG	2.33	0.42
1:A:24:ILE:O	1:A:28:ILE:HD12	2.20	0.42
1:B:27:ALA:CB	1:B:69:ILE:HD11	2.49	0.42
1:B:62:ALA:HB1	1:B:64:ILE:CD1	2.49	0.42
1:B:315:LEU:HD12	1:B:316:SER:N	2.35	0.42
1:C:147:LEU:O	1:C:151:LEU:HD12	2.20	0.42
1:C:296:GLY:O	1:C:392:LEU:HB2	2.19	0.42
1:A:252:LYS:HB3	1:A:255:VAL:HG22	2.00	0.42
1:B:116:ASP:OD2	1:B:116:ASP:C	2.62	0.42
1:B:121:ASP:CG	1:B:122:GLN:N	2.78	0.42
1:C:29:LEU:HD13	1:C:41:VAL:HB	2.02	0.42
1:A:51:LEU:HD13	1:C:94:LEU:HD21	2.01	0.42
1:B:62:ALA:HB1	1:B:64:ILE:HD13	2.00	0.42
1:B:253:ILE:HD12	1:B:264:HIS:NE2	2.34	0.42
1:B:291:SER:HB3	1:B:382:LEU:HD11	2.02	0.42
1:B:322:TYR:O	1:B:324:THR:CG2	2.65	0.42
1:C:315:LEU:HD11	1:D:230:TYR:CZ	2.54	0.42
1:A:200:VAL:HG12	1:A:201:SER:O	2.20	0.42
1:D:148:ILE:CD1	1:D:169:SER:HB3	2.50	0.42
1:A:207:LYS:O	1:A:211:GLU:CG	2.66	0.41
1:A:326:ILE:HD12	1:A:326:ILE:C	2.45	0.41
1:C:326:ILE:C	1:C:327:CYS:O	2.59	0.41
1:D:238:VAL:HG13	1:D:239:ILE:HG12	2.01	0.41
1:D:288:ILE:CG2	1:D:303:VAL:HG21	2.50	0.41
1:D:336:VAL:CG1	1:D:342:MET:CE	2.93	0.41
1:C:209:LEU:O	1:C:213:VAL:CG1	2.64	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:125:MET:HE2	1:D:277:VAL:C	2.46	0.41
1:A:163:LEU:HD11	1:A:194:THR:HG21	2.03	0.41
1:C:154:CYS:HB3	1:C:160:LEU:HB2	2.02	0.41
1:C:219:PRO:HG2	1:C:222:LEU:HD22	2.01	0.41
1:D:201:SER:O	1:D:202:ASN:C	2.63	0.41
1:B:94:LEU:HD22	1:C:96:CYS:HB2	2.02	0.41
1:B:321:THR:CG2	1:B:325:GLY:N	2.83	0.41
1:A:94:LEU:HD23	1:A:94:LEU:N	2.36	0.41
1:B:201:SER:OG	1:B:204:GLU:CB	2.69	0.41
1:B:297:LEU:CD1	1:B:335:ILE:HD13	2.51	0.41
1:B:332:LEU:CA	1:B:335:ILE:HG23	2.50	0.41
1:C:136:MET:HE1	1:C:293:VAL:HG22	2.02	0.41
1:C:301:VAL:CG1	1:C:302:LEU:N	2.82	0.41
1:C:328:ASP:O	1:C:331:ILE:N	2.50	0.41
1:D:284:CYS:CB	1:D:347:ILE:HD13	2.51	0.41
1:D:391:LYS:O	1:D:392:LEU:HD23	2.21	0.41
1:A:252:LYS:NZ	2:A:400:ACT:H3	2.36	0.41
1:B:209:LEU:CA	1:B:213:VAL:HG22	2.51	0.41
1:B:231:TYR:N	1:B:231:TYR:CD2	2.89	0.41
1:A:127:GLY:HA3	1:A:285:ALA:HB1	2.03	0.40
1:A:319:VAL:O	1:A:319:VAL:CG2	2.69	0.40
1:B:28:ILE:O	1:B:31:ALA:HB3	2.21	0.40
1:B:154:CYS:HB3	1:B:160:LEU:HB2	2.02	0.40
1:C:173:LEU:HD13	1:C:174:GLU:O	2.20	0.40
1:D:63:VAL:C	1:D:64:ILE:HD12	2.46	0.40
1:D:339:ASN:N	1:D:339:ASN:ND2	2.68	0.40
1:C:389:LYS:CB	1:C:390:PRO:HD3	2.51	0.40
1:B:28:ILE:N	1:B:69:ILE:HD11	2.36	0.40
1:B:209:LEU:HG	1:B:231:TYR:CD1	2.57	0.40
1:C:308:ALA:CB	1:C:315:LEU:HD13	2.51	0.40
1:D:252:LYS:HB3	1:D:255:VAL:HG22	2.02	0.40
1:A:92:SER:OG	1:C:51:LEU:HD12	2.21	0.40
1:C:20:MET:HE2	1:C:79:TYR:CE2	2.57	0.40
1:C:318:ASN:OD1	1:C:319:VAL:N	2.55	0.40
1:D:192:ILE:CD1	1:D:209:LEU:HD22	2.51	0.40
1:D:355:ARG:HA	1:D:356:PRO:HD3	1.98	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	375/415 (90%)	352 (94%)	22 (6%)	1 (0%)	36	66
1	B	375/415 (90%)	351 (94%)	24 (6%)	0	100	100
1	C	375/415 (90%)	349 (93%)	22 (6%)	4 (1%)	11	41
1	D	375/415 (90%)	349 (93%)	25 (7%)	1 (0%)	36	66
All	All	1500/1660 (90%)	1401 (93%)	93 (6%)	6 (0%)	30	60

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	121	ASP
1	A	102	GLN
1	C	327	CYS
1	C	392	LEU
1	C	117	ILE
1	C	356	PRO

5.3.2 Protein sidechains [i](#)

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	306/351 (87%)	239 (78%)	67 (22%)	1	5
1	B	312/351 (89%)	237 (76%)	75 (24%)	1	4
1	C	308/351 (88%)	234 (76%)	74 (24%)	1	4
1	D	310/351 (88%)	247 (80%)	63 (20%)	1	6

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1236/1404 (88%)	957 (77%)	279 (23%)	1 4

All (279) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	VAL
1	A	28	ILE
1	A	43	CYS
1	A	49	THR
1	A	52	ILE
1	A	57	GLU
1	A	58	ILE
1	A	68	LYS
1	A	69	ILE
1	A	76	SER
1	A	82	THR
1	A	89	GLN
1	A	94	LEU
1	A	97	VAL
1	A	98	GLU
1	A	100	GLN
1	A	102	GLN
1	A	103	ASP
1	A	122	GLN
1	A	124	ILE
1	A	125	MET
1	A	134	LYS
1	A	144	SER
1	A	148	ILE
1	A	150	ARG
1	A	152	GLN
1	A	170	GLN
1	A	172	THR
1	A	176	GLU
1	A	189	THR
1	A	190	ILE
1	A	192	ILE
1	A	205	ILE
1	A	211	GLU
1	A	222	LEU
1	A	223	MET
1	A	224	ASP

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Mol	Chain	Res	Type
1	A	227	MET
1	A	228	LEU
1	A	245	ASP
1	A	253	ILE
1	A	277	VAL
1	A	280	SER
1	A	293	VAL
1	A	301	VAL
1	A	306	SER
1	A	316	SER
1	A	319	VAL
1	A	324	THR
1	A	327	CYS
1	A	329	GLU
1	A	330	SER
1	A	333	VAL
1	A	334	ASP
1	A	335	ILE
1	A	346	MET
1	A	354	THR
1	A	357	ILE
1	A	361	THR
1	A	378	PHE
1	A	381	GLU
1	A	382	LEU
1	A	383	GLU
1	A	384	ILE
1	A	387	GLU
1	A	388	LEU
1	A	392	LEU
1	B	12	VAL
1	B	28	ILE
1	B	43	CYS
1	B	44	GLU
1	B	49	THR
1	B	51	LEU
1	B	52	ILE
1	B	57	GLU
1	B	58	ILE
1	B	63	VAL
1	B	69	ILE
1	B	76	SER

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Mol	Chain	Res	Type
1	B	80	ASP
1	B	81	ASP
1	B	86	PHE
1	B	94	LEU
1	B	100	GLN
1	B	102	GLN
1	B	117	ILE
1	B	122	GLN
1	B	134	LYS
1	B	136	MET
1	B	142	VAL
1	B	144	SER
1	B	145	THR
1	B	150	ARG
1	B	152	GLN
1	B	170	GLN
1	B	172	THR
1	B	173	LEU
1	B	176	GLU
1	B	189	THR
1	B	190	ILE
1	B	192	ILE
1	B	205	ILE
1	B	209	LEU
1	B	215	GLN
1	B	217	VAL
1	B	221	GLU
1	B	223	MET
1	B	227	MET
1	B	231	TYR
1	B	234	SER
1	B	238	VAL
1	B	243	MET
1	B	245	ASP
1	B	253	ILE
1	B	277	VAL
1	B	280	SER
1	B	293	VAL
1	B	301	VAL
1	B	306	SER
1	B	315	LEU
1	B	316	SER

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Mol	Chain	Res	Type
1	B	324	THR
1	B	326	ILE
1	B	328	ASP
1	B	330	SER
1	B	331	ILE
1	B	334	ASP
1	B	335	ILE
1	B	337	ASN
1	B	346	MET
1	B	348	ILE
1	B	349	LYS
1	B	354	THR
1	B	355	ARG
1	B	357	ILE
1	B	361	THR
1	B	381	GLU
1	B	382	LEU
1	B	383	GLU
1	B	384	ILE
1	B	387	GLU
1	B	391	LYS
1	C	12	VAL
1	C	19	LYS
1	C	28	ILE
1	C	43	CYS
1	C	44	GLU
1	C	49	THR
1	C	51	LEU
1	C	52	ILE
1	C	53	LEU
1	C	57	GLU
1	C	58	ILE
1	C	63	VAL
1	C	66	ILE
1	C	68	LYS
1	C	69	ILE
1	C	76	SER
1	C	82	THR
1	C	90	THR
1	C	102	GLN
1	C	103	ASP
1	C	122	GLN

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Mol	Chain	Res	Type
1	C	134	LYS
1	C	135	GLU
1	C	136	MET
1	C	144	SER
1	C	148	ILE
1	C	150	ARG
1	C	151	LEU
1	C	170	GLN
1	C	172	THR
1	C	173	LEU
1	C	174	GLU
1	C	176	GLU
1	C	185	ILE
1	C	189	THR
1	C	190	ILE
1	C	192	ILE
1	C	207	LYS
1	C	209	LEU
1	C	213	VAL
1	C	217	VAL
1	C	222	LEU
1	C	224	ASP
1	C	226	LYS
1	C	227	MET
1	C	228	LEU
1	C	234	SER
1	C	239	ILE
1	C	243	MET
1	C	245	ASP
1	C	253	ILE
1	C	280	SER
1	C	293	VAL
1	C	306	SER
1	C	316	SER
1	C	317	ILE
1	C	324	THR
1	C	327	CYS
1	C	328	ASP
1	C	329	GLU
1	C	330	SER
1	C	331	ILE
1	C	334	ASP

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Mol	Chain	Res	Type
1	C	337	ASN
1	C	349	LYS
1	C	354	THR
1	C	355	ARG
1	C	357	ILE
1	C	361	THR
1	C	369	ARG
1	C	381	GLU
1	C	382	LEU
1	C	387	GLU
1	C	392	LEU
1	D	6	PHE
1	D	12	VAL
1	D	19	LYS
1	D	28	ILE
1	D	43	CYS
1	D	49	THR
1	D	51	LEU
1	D	52	ILE
1	D	57	GLU
1	D	58	ILE
1	D	66	ILE
1	D	69	ILE
1	D	76	SER
1	D	82	THR
1	D	97	VAL
1	D	100	GLN
1	D	102	GLN
1	D	115	GLU
1	D	117	ILE
1	D	122	GLN
1	D	134	LYS
1	D	144	SER
1	D	148	ILE
1	D	150	ARG
1	D	170	GLN
1	D	172	THR
1	D	173	LEU
1	D	176	GLU
1	D	189	THR
1	D	190	ILE
1	D	192	ILE

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Mol	Chain	Res	Type
1	D	205	ILE
1	D	207	LYS
1	D	212	GLU
1	D	223	MET
1	D	224	ASP
1	D	227	MET
1	D	234	SER
1	D	239	ILE
1	D	245	ASP
1	D	251	ARG
1	D	253	ILE
1	D	280	SER
1	D	293	VAL
1	D	301	VAL
1	D	306	SER
1	D	316	SER
1	D	324	THR
1	D	328	ASP
1	D	330	SER
1	D	334	ASP
1	D	337	ASN
1	D	338	LYS
1	D	339	ASN
1	D	346	MET
1	D	347	ILE
1	D	348	ILE
1	D	349	LYS
1	D	354	THR
1	D	357	ILE
1	D	361	THR
1	D	381	GLU
1	D	384	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	GLN
1	A	195	GLN
1	A	313	HIS
1	A	339	ASN
1	B	23	GLN
1	B	83	ASN

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Mol	Chain	Res	Type
1	B	313	HIS
1	B	339	ASN
1	C	23	GLN
1	C	40	HIS
1	C	304	GLN
1	D	23	GLN
1	D	100	GLN
1	D	294	HIS
1	D	337	ASN

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](#)

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	ACT	C	400	-	3,3,3	0.84	0	3,3,3	1.51	0
2	ACT	B	400	-	3,3,3	0.90	0	3,3,3	1.42	0
2	ACT	A	400	-	3,3,3	0.81	0	3,3,3	1.60	0
2	ACT	D	400	-	3,3,3	0.85	0	3,3,3	1.37	0

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.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	400	ACT	3	0
2	A	400	ACT	1	0

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	379/415 (91%)	-0.35	0 100 100	47, 68, 99, 129	0
1	B	379/415 (91%)	-0.30	0 100 100	44, 73, 108, 124	0
1	C	379/415 (91%)	-0.22	0 100 100	49, 80, 119, 139	0
1	D	379/415 (91%)	-0.15	0 100 100	54, 89, 122, 142	0
All	All	1516/1660 (91%)	-0.26	0 100 100	44, 76, 114, 142	0

There are no RSRZ outliers to report.

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	ACT	C	400	4/4	0.76	0.12	43,47,47,48	0
2	ACT	A	400	4/4	0.85	0.12	39,40,41,43	0
2	ACT	B	400	4/4	0.86	0.13	31,32,32,35	0
2	ACT	D	400	4/4	0.95	0.06	35,37,38,40	0

6.5 Other polymers [i](#)

There are no such residues in this entry.