



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:12 PM UTC

PDB ID : 3P0J / pdb_00003p0j
Title : Leishmania major Tyrosyl-tRNA synthetase in complex with tyrosinol, triclinic crystal form 1
Authors : Larson, E.T.; Merritt, E.A.; Medical Structural Genomics of Pathogenic Protozoa (MSGPP)
Deposited on : 2010-09-28
Resolution : 2.89 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

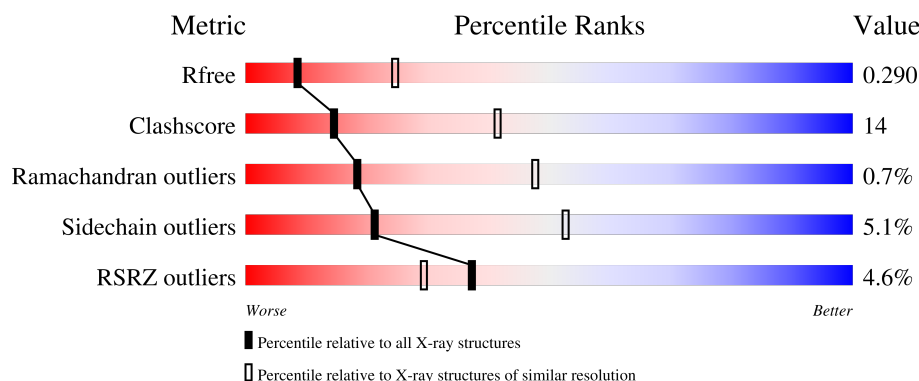
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.89 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	690	 8% 70% 23% • 5%
1	B	690	 4% 60% 22% • 16%
1	C	690	 3% 68% 26% • •
1	D	690	 2% 70% 24% • •

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 19700 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 Tyrosyl-tRNA synthetase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	657	Total	C	N	O	S	0	0	0
			5067	3188	881	961	37			
1	B	581	Total	C	N	O	S	0	0	0
			4381	2763	746	837	35			
1	C	663	Total	C	N	O	S	0	1	0
			5106	3216	884	969	37			
1	D	665	Total	C	N	O	S	0	0	0
			5096	3206	885	967	38			

There are 32 discrepancies between the modelled and reference sequences:

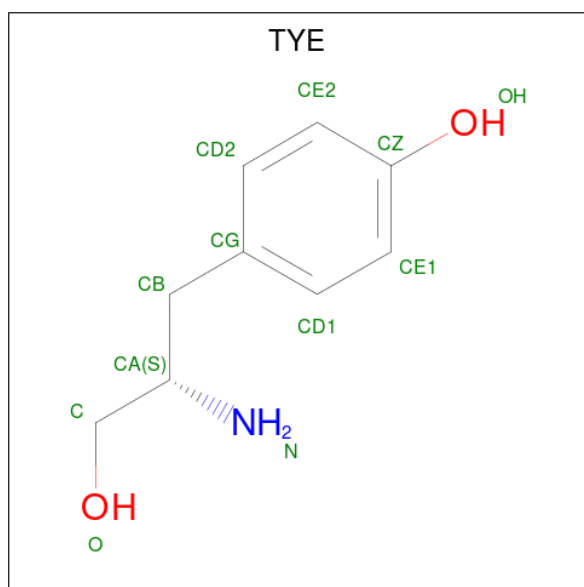
Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	MET	-	expression tag	UNP Q4QFJ7
A	-6	ALA	-	expression tag	UNP Q4QFJ7
A	-5	HIS	-	expression tag	UNP Q4QFJ7
A	-4	HIS	-	expression tag	UNP Q4QFJ7
A	-3	HIS	-	expression tag	UNP Q4QFJ7
A	-2	HIS	-	expression tag	UNP Q4QFJ7
A	-1	HIS	-	expression tag	UNP Q4QFJ7
A	0	HIS	-	expression tag	UNP Q4QFJ7
B	-7	MET	-	expression tag	UNP Q4QFJ7
B	-6	ALA	-	expression tag	UNP Q4QFJ7
B	-5	HIS	-	expression tag	UNP Q4QFJ7
B	-4	HIS	-	expression tag	UNP Q4QFJ7
B	-3	HIS	-	expression tag	UNP Q4QFJ7
B	-2	HIS	-	expression tag	UNP Q4QFJ7
B	-1	HIS	-	expression tag	UNP Q4QFJ7
B	0	HIS	-	expression tag	UNP Q4QFJ7
C	-7	MET	-	expression tag	UNP Q4QFJ7
C	-6	ALA	-	expression tag	UNP Q4QFJ7
C	-5	HIS	-	expression tag	UNP Q4QFJ7
C	-4	HIS	-	expression tag	UNP Q4QFJ7
C	-3	HIS	-	expression tag	UNP Q4QFJ7

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	HIS	-	expression tag	UNP Q4QFJ7
C	-1	HIS	-	expression tag	UNP Q4QFJ7
C	0	HIS	-	expression tag	UNP Q4QFJ7
D	-7	MET	-	expression tag	UNP Q4QFJ7
D	-6	ALA	-	expression tag	UNP Q4QFJ7
D	-5	HIS	-	expression tag	UNP Q4QFJ7
D	-4	HIS	-	expression tag	UNP Q4QFJ7
D	-3	HIS	-	expression tag	UNP Q4QFJ7
D	-2	HIS	-	expression tag	UNP Q4QFJ7
D	-1	HIS	-	expression tag	UNP Q4QFJ7
D	0	HIS	-	expression tag	UNP Q4QFJ7

- Molecule 2 is 4-[(2S)-2-amino-3-hydroxypropyl]phenol (CCD ID: TYE) (formula: $C_9H_{13}NO_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			12	9	1	2		
2	B	1	Total	C	N	O	0	0
			12	9	1	2		
2	C	1	Total	C	N	O	0	0
			12	9	1	2		
2	D	1	Total	C	N	O	0	0
			12	9	1	2		

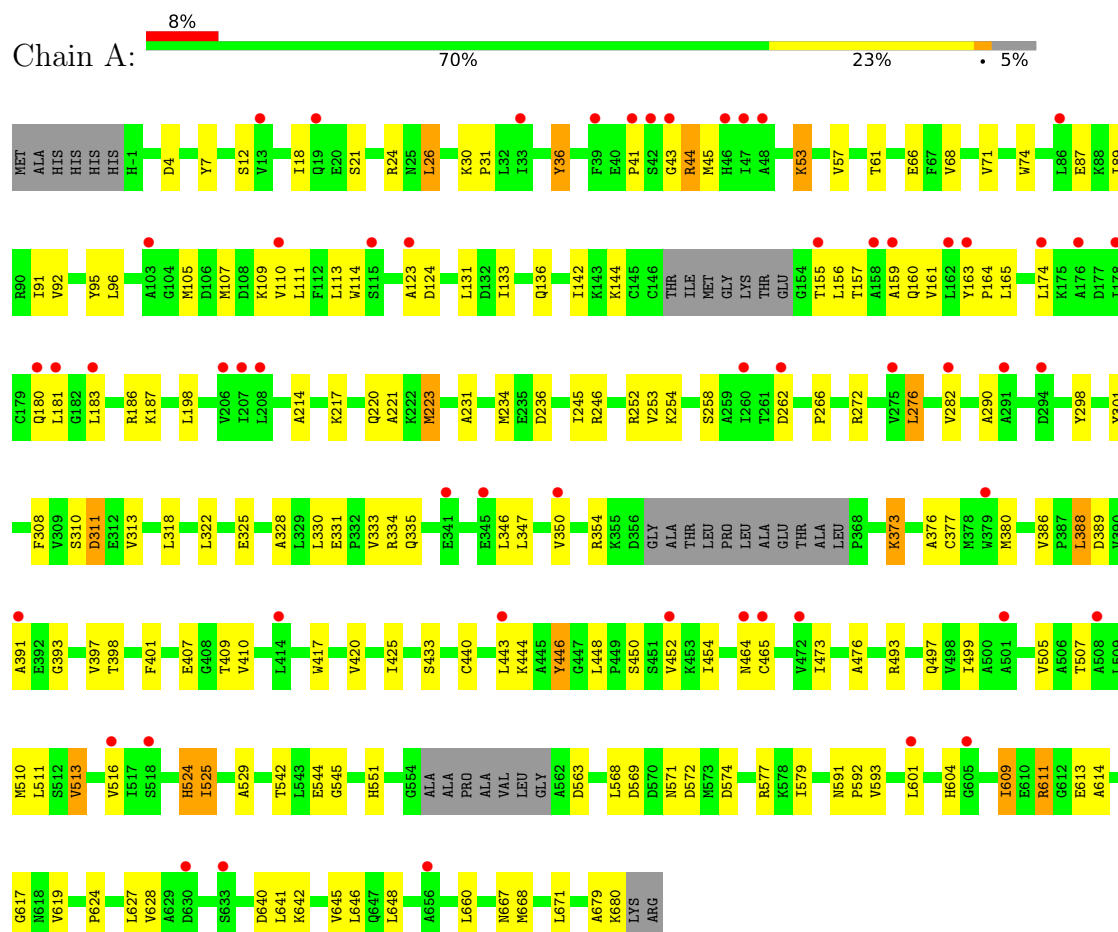
- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total 1	Na 1	0	0
3	C	1	Total 1	Na 1	0	0

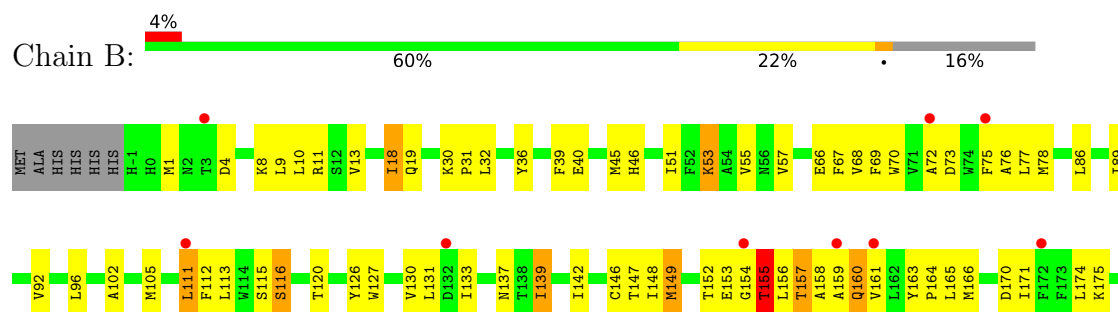
3 Residue-property plots [i](#)

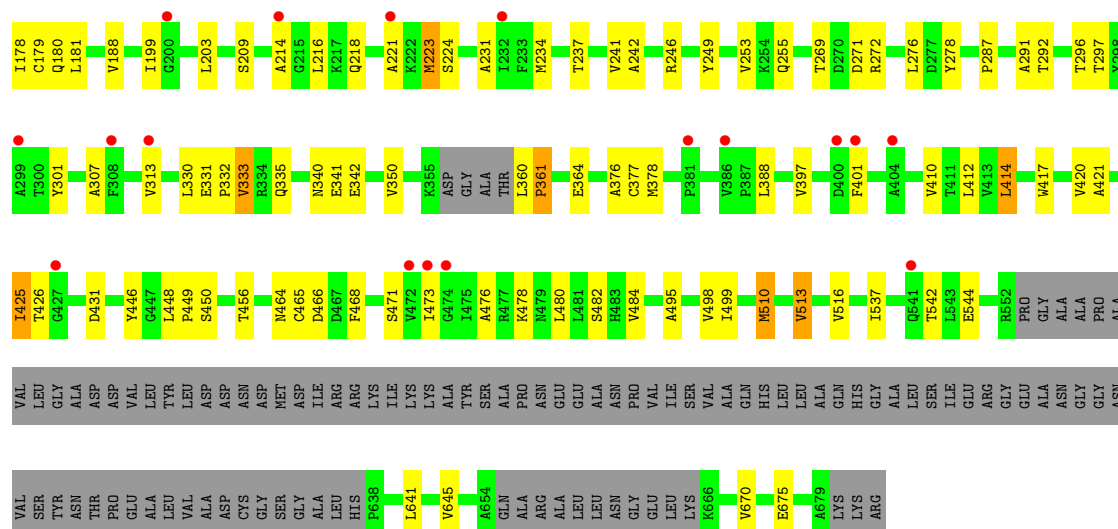
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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



• Molecule 1: Tyrosyl-tRNA synthetase



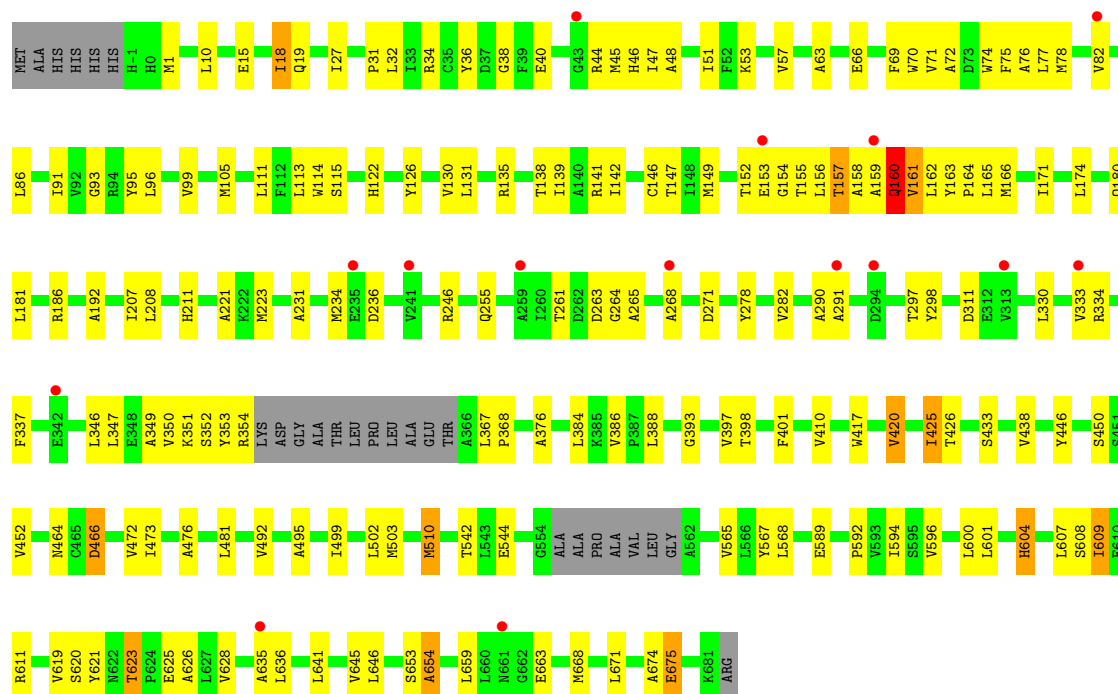


• Molecule 1: Tyrosyl-tRNA synthetase



• Molecule 1: Tyrosyl-tRNA synthetase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	57.52Å 111.93Å 141.04Å 74.47° 80.83° 77.13°	Depositor
Resolution (Å)	50.00 – 2.89 50.00 – 2.89	Depositor EDS
% Data completeness (in resolution range)	98.1 (50.00-2.89) 98.4 (50.00-2.89)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.48 (at 2.91Å)	Xtriage
Refinement program	REFMAC refmac _5.5.0110	Depositor
R, R_{free}	0.244 , 0.293 0.239 , 0.290	Depositor DCC
R_{free} test set	3719 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	73.3	Xtriage
Anisotropy	0.328	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 0.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.029 for h,h-k,h-l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	19700	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.99% 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: NA, TYE

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	0.64	0/5153	0.87	1/6976 (0.0%)
1	B	0.57	1/4456 (0.0%)	0.86	2/6051 (0.0%)
1	C	0.52	0/5194	0.80	0/7033
1	D	0.48	0/5182	0.79	2/7022 (0.0%)
All	All	0.55	1/19985 (0.0%)	0.83	5/27082 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	D	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	157	THR	CA-CB	5.31	1.62	1.53

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	160	GLN	N-CA-C	-8.10	103.48	112.72
1	B	153	GLU	N-CA-C	6.49	118.01	111.07
1	A	160	GLN	N-CA-C	-5.92	105.64	112.92
1	D	186	ARG	N-CA-C	5.15	117.56	111.33
1	D	160	GLN	N-CA-C	-5.07	106.68	112.92

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	635	ALA	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	5067	0	5063	112	0
1	B	4381	0	4268	140	0
1	C	5106	0	5119	157	0
1	D	5096	0	5076	149	0
2	A	12	0	12	0	0
2	B	12	0	12	0	0
2	C	12	0	12	1	0
2	D	12	0	12	0	0
3	A	1	0	0	0	0
3	C	1	0	0	0	0
All	All	19700	0	19574	550	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 14.

All (550) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:417:TRP:O	1:B:420:VAL:HG22	1.47	1.11
1:D:159:ALA:HB3	1:D:476:ALA:HB1	1.11	1.10
1:D:45:MET:HE3	1:D:99:VAL:HG11	1.38	1.06
1:C:45:MET:CE	1:C:99:VAL:HG11	1.87	1.05
1:C:45:MET:HE2	1:C:99:VAL:HG11	1.07	1.04
1:B:148:ILE:HD13	1:B:188:VAL:HG21	1.33	1.04
1:D:159:ALA:CB	1:D:476:ALA:HB1	1.87	1.04
1:C:45:MET:HE2	1:C:99:VAL:CG1	1.91	1.01
1:D:159:ALA:HB3	1:D:476:ALA:CB	1.92	0.99
1:C:417:TRP:O	1:C:420:VAL:HG22	1.64	0.97
1:B:149:MET:HE3	1:B:163:TYR:CG	1.99	0.96
1:A:660:LEU:HD22	1:A:668:MET:HE2	1.47	0.95
1:D:142:ILE:HD13	1:D:165:LEU:HD21	1.49	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:102:ALA:HB3	1:B:333:VAL:HG23	1.50	0.93
1:B:148:ILE:HD13	1:B:188:VAL:CG2	2.00	0.90
1:A:330:LEU:O	1:A:333:VAL:HG12	1.72	0.90
1:A:159:ALA:HB3	1:A:476:ALA:HB1	1.54	0.89
1:A:440:CYS:SG	1:A:454:ILE:HD13	2.12	0.88
1:D:77:LEU:HD22	1:D:86:LEU:HD12	1.54	0.88
1:B:68:VAL:HG22	1:B:111:LEU:HD13	1.55	0.87
1:C:391:ALA:HB2	1:C:568:LEU:HD11	1.58	0.85
1:B:155:THR:OG1	1:B:482:SER:N	2.09	0.85
1:C:91:ILE:HG23	1:C:346:LEU:HD23	1.56	0.85
1:B:1:MET:HE1	1:B:9:LEU:HD22	1.58	0.84
1:C:45:MET:CE	1:C:99:VAL:CG1	2.52	0.83
1:D:57:VAL:HG21	1:D:105:MET:CE	2.09	0.82
1:B:78:MET:SD	1:B:166:MET:CE	2.68	0.81
1:A:660:LEU:CD2	1:A:668:MET:HE2	2.11	0.81
1:B:78:MET:SD	1:B:166:MET:HE1	2.21	0.81
1:B:148:ILE:CD1	1:B:188:VAL:HG21	2.10	0.80
1:D:142:ILE:CD1	1:D:165:LEU:HD21	2.10	0.80
1:B:149:MET:CE	1:B:163:TYR:HB3	2.12	0.79
1:B:417:TRP:O	1:B:420:VAL:CG2	2.29	0.79
1:A:45:MET:HE3	1:A:96:LEU:HD22	1.64	0.79
1:D:40:GLU:HB2	1:D:76:ALA:HB2	1.65	0.78
1:D:162:LEU:HG	1:D:166:MET:HE2	1.65	0.77
1:D:57:VAL:HG21	1:D:105:MET:HE2	1.65	0.77
1:A:386:VAL:HB	1:A:568:LEU:HD13	1.67	0.77
1:B:155:THR:HG1	1:B:482:SER:H	1.33	0.76
1:D:596:VAL:HG12	1:D:600:LEU:HD11	1.67	0.76
1:C:95:TYR:CE1	1:C:347:LEU:HD13	2.21	0.76
1:C:420:VAL:HG12	1:C:432:ILE:HD12	1.66	0.76
1:B:102:ALA:CB	1:B:333:VAL:HG23	2.15	0.76
1:C:223:MET:HG3	1:C:231:ALA:HB2	1.66	0.76
1:B:45:MET:HE2	1:B:96:LEU:HD22	1.69	0.75
1:A:105:MET:HB3	1:A:107:MET:HE2	1.70	0.74
1:C:51:ILE:HG12	1:C:234:MET:HE1	1.69	0.74
1:D:45:MET:CE	1:D:99:VAL:HG11	2.16	0.74
1:D:438:VAL:HG11	1:D:668:MET:CE	2.18	0.74
1:A:391:ALA:HB2	1:A:568:LEU:HD11	1.70	0.73
1:C:159:ALA:HB3	1:C:476:ALA:HB1	1.70	0.73
1:A:159:ALA:CB	1:A:476:ALA:HB1	2.18	0.73
1:C:213:LEU:HD23	1:C:223:MET:HE1	1.70	0.73
1:C:585:ALA:HB3	1:C:588:GLU:CD	2.13	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:596:VAL:O	1:D:600:LEU:HD12	1.89	0.72
1:C:51:ILE:HG21	1:C:329:LEU:HD13	1.71	0.72
1:A:18:ILE:HG22	1:A:18:ILE:O	1.90	0.71
1:B:75:PHE:HA	1:B:78:MET:HE2	1.72	0.71
1:D:18:ILE:HD13	1:D:207:ILE:HD12	1.71	0.71
1:A:66:GLU:OE2	1:B:66:GLU:OE2	2.10	0.70
1:A:131:LEU:HD22	1:A:499:ILE:HG21	1.73	0.70
1:C:223:MET:HE3	1:C:231:ALA:HB1	1.73	0.69
1:C:516:VAL:HG13	1:C:537:ILE:HG21	1.74	0.69
1:D:69:PHE:CZ	1:D:105:MET:HE1	2.28	0.69
1:D:163:TYR:HB3	1:D:164:PRO:HD3	1.74	0.69
1:D:51:ILE:HG21	1:D:282:VAL:HG13	1.75	0.69
1:D:78:MET:SD	1:D:166:MET:HE1	2.32	0.69
1:D:149:MET:HE3	1:D:160:GLN:HB2	1.75	0.69
1:B:149:MET:CE	1:B:160:GLN:HB3	2.23	0.68
1:D:46:HIS:NE2	1:D:48:ALA:HB3	2.08	0.68
1:A:133:ILE:HG22	1:A:165:LEU:HD22	1.75	0.68
1:C:611:ARG:NH2	1:C:635:ALA:O	2.25	0.68
1:D:346:LEU:O	1:D:350:VAL:HG13	1.94	0.68
1:A:542:THR:HG21	1:A:544:GLU:OE2	1.92	0.68
1:D:542:THR:HG21	1:D:544:GLU:OE2	1.93	0.68
1:B:149:MET:HE3	1:B:163:TYR:HB3	1.75	0.68
1:A:507:THR:HG22	1:A:511:LEU:HD12	1.76	0.68
1:B:149:MET:HE3	1:B:163:TYR:CB	2.23	0.68
1:B:116:SER:O	1:B:120:THR:HG23	1.94	0.68
1:B:307:ALA:HB1	1:B:313:VAL:HG13	1.75	0.68
1:C:601:LEU:HD23	1:C:601:LEU:C	2.19	0.68
1:A:542:THR:HG23	1:A:544:GLU:HG3	1.75	0.68
1:D:589:GLU:O	1:D:594:ILE:HD13	1.95	0.67
1:A:464:ASN:CB	1:A:510:MET:HE1	2.25	0.67
1:C:113:LEU:HD11	1:D:32:LEU:HD11	1.77	0.67
1:C:569:ASP:HB2	1:C:668:MET:HE1	1.77	0.67
1:B:142:ILE:CD1	1:B:165:LEU:HD21	2.25	0.67
1:A:613:GLU:O	1:A:614:ALA:HB3	1.93	0.67
1:B:142:ILE:HD12	1:B:495:ALA:CB	2.25	0.66
1:D:473:ILE:O	1:D:476:ALA:HB3	1.95	0.66
1:C:32:LEU:HD11	1:D:113:LEU:HD11	1.77	0.66
1:C:91:ILE:HG23	1:C:346:LEU:CD2	2.25	0.66
1:C:386:VAL:HB	1:C:568:LEU:HD13	1.77	0.66
1:D:157:THR:HG22	1:D:160:GLN:HG3	1.77	0.66
1:B:78:MET:SD	1:B:166:MET:HE2	2.36	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:105:MET:HE3	1:A:110:VAL:HG21	1.75	0.66
1:B:78:MET:HG2	1:B:473:ILE:HG23	1.78	0.66
1:C:397:VAL:HG21	1:C:542:THR:HA	1.78	0.66
1:A:113:LEU:HD12	1:A:174:LEU:HD21	1.78	0.65
1:B:542:THR:HG21	1:B:544:GLU:OE2	1.96	0.65
1:A:660:LEU:HD22	1:A:668:MET:CE	2.25	0.65
1:C:440:CYS:SG	1:C:454:ILE:HD13	2.37	0.65
1:D:417:TRP:O	1:D:420:VAL:HG22	1.95	0.65
1:B:154:GLY:O	1:B:155:THR:HG22	1.97	0.65
1:D:347:LEU:HA	1:D:350:VAL:HG22	1.79	0.64
1:D:261:THR:HG23	1:D:264:GLY:H	1.63	0.64
1:A:105:MET:HE1	1:A:110:VAL:HG11	1.79	0.64
1:A:417:TRP:O	1:A:420:VAL:HG22	1.97	0.64
1:D:565:VAL:CG1	1:D:567:TYR:CZ	2.81	0.64
1:B:376:ALA:HB2	1:B:401:PHE:CZ	2.32	0.64
1:B:142:ILE:HD12	1:B:495:ALA:HB2	1.80	0.64
1:A:53:LYS:O	1:A:57:VAL:HG23	1.98	0.63
1:A:464:ASN:HB2	1:A:510:MET:HE1	1.80	0.63
1:A:668:MET:HA	1:A:671:LEU:HD13	1.80	0.63
1:C:542:THR:HG23	1:C:544:GLU:HG3	1.80	0.63
1:A:105:MET:CB	1:A:107:MET:HE2	2.27	0.63
1:B:149:MET:HE1	1:B:163:TYR:HB3	1.79	0.63
1:D:57:VAL:CB	1:D:105:MET:HE2	2.29	0.63
1:D:221:ALA:HB2	1:D:268:ALA:N	2.13	0.62
1:D:347:LEU:HD12	1:D:350:VAL:CG2	2.29	0.62
1:B:269:THR:HG22	1:B:271:ASP:H	1.63	0.62
1:A:611:ARG:NH1	1:A:617:GLY:O	2.32	0.62
1:D:57:VAL:CG2	1:D:105:MET:HE2	2.28	0.62
1:D:255:GLN:NE2	1:D:271:ASP:O	2.30	0.62
1:D:438:VAL:HG11	1:D:668:MET:HE2	1.81	0.62
1:D:95:TYR:CE1	1:D:347:LEU:HD13	2.35	0.62
1:A:113:LEU:HD12	1:A:174:LEU:CD2	2.29	0.62
1:A:245:ILE:HD12	1:A:322:LEU:HD23	1.81	0.62
1:C:613:GLU:O	1:C:614:ALA:HB3	1.99	0.62
1:B:149:MET:CE	1:B:163:TYR:CB	2.78	0.61
1:D:180:GLN:NE2	1:D:208:LEU:O	2.33	0.61
1:D:565:VAL:HG12	1:D:567:TYR:CZ	2.36	0.61
1:C:376:ALA:HB2	1:C:401:PHE:CZ	2.34	0.61
1:D:159:ALA:CB	1:D:476:ALA:CB	2.66	0.61
1:C:223:MET:HG3	1:C:231:ALA:CB	2.30	0.61
1:C:660:LEU:HD22	1:C:668:MET:HE2	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:78:MET:HE1	1:D:163:TYR:N	2.14	0.61
1:A:113:LEU:HD11	1:B:32:LEU:HD11	1.83	0.61
1:C:111:LEU:HD11	1:D:111:LEU:HD21	1.82	0.61
1:B:78:MET:HE1	1:B:166:MET:CE	2.30	0.61
1:A:214:ALA:HB2	1:A:272:ARG:HD3	1.82	0.61
1:B:112:PHE:O	1:B:113:LEU:HD23	2.00	0.60
1:C:131:LEU:HD22	1:C:499:ILE:HG21	1.83	0.60
1:C:236:ASP:O	1:C:334:ARG:NH1	2.33	0.60
1:B:154:GLY:C	1:B:156:LEU:H	2.10	0.60
1:C:516:VAL:HG23	1:C:516:VAL:O	2.02	0.60
1:B:127:TRP:O	1:B:130:VAL:HG22	2.02	0.60
1:C:142:ILE:HD13	1:C:165:LEU:HD21	1.83	0.60
1:B:155:THR:C	1:B:157:THR:N	2.58	0.60
1:D:69:PHE:CE2	1:D:105:MET:HE1	2.36	0.60
1:A:57:VAL:HG11	1:A:105:MET:HE3	1.84	0.59
1:C:142:ILE:CD1	1:C:165:LEU:HD21	2.32	0.59
1:A:71:VAL:HG11	1:A:114:TRP:CZ3	2.37	0.59
1:A:308:PHE:HA	1:A:313:VAL:HG23	1.83	0.59
1:D:82:VAL:HG12	1:D:82:VAL:O	2.03	0.59
1:B:388:LEU:HD11	1:B:446:TYR:CZ	2.38	0.59
1:C:131:LEU:CD2	1:C:499:ILE:HG21	2.33	0.59
1:D:72:ALA:HB1	1:D:75:PHE:HB2	1.85	0.59
1:B:78:MET:CE	1:B:166:MET:CE	2.80	0.59
1:C:279:PHE:O	1:C:283:VAL:HG23	2.03	0.59
1:C:261:THR:HG23	1:C:264:GLY:H	1.68	0.58
1:D:57:VAL:HG11	1:D:105:MET:HE2	1.85	0.58
1:B:360:LEU:CB	1:B:361:PRO:CD	2.81	0.58
1:A:613:GLU:O	1:A:614:ALA:CB	2.51	0.58
1:D:10:LEU:HD23	1:D:27:ILE:HD11	1.86	0.58
1:D:171:ILE:HD12	1:D:192:ALA:HB2	1.85	0.58
1:B:214:ALA:HA	1:B:221:ALA:O	2.03	0.58
1:C:660:LEU:HD22	1:C:668:MET:CE	2.33	0.58
1:D:159:ALA:HB3	1:D:476:ALA:CA	2.33	0.58
1:C:142:ILE:HD12	1:C:495:ALA:CB	2.34	0.58
1:D:466:ASP:N	1:D:466:ASP:OD1	2.35	0.58
1:D:142:ILE:HD12	1:D:495:ALA:CB	2.33	0.58
1:A:524:HIS:O	1:A:525:ILE:C	2.46	0.58
1:A:398:THR:HG21	1:A:452:VAL:HG21	1.86	0.58
1:D:45:MET:HE3	1:D:99:VAL:CG1	2.24	0.58
1:D:386:VAL:HB	1:D:568:LEU:HD12	1.86	0.57
1:B:51:ILE:O	1:B:55:VAL:HG23	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:154:GLY:O	1:D:155:THR:HG22	2.05	0.57
1:B:92:VAL:HG22	1:B:350:VAL:HG22	1.86	0.57
1:D:91:ILE:HG23	1:D:346:LEU:CD2	2.34	0.57
1:B:77:LEU:HD22	1:B:86:LEU:HD13	1.85	0.57
1:C:391:ALA:CB	1:C:568:LEU:HD11	2.33	0.57
1:D:393:GLY:O	1:D:397:VAL:HG23	2.04	0.57
1:C:18:ILE:CG2	1:C:18:ILE:O	2.52	0.57
1:D:18:ILE:O	1:D:18:ILE:HG22	2.03	0.57
1:A:667:ASN:O	1:A:671:LEU:CD1	2.53	0.57
1:C:378:MET:HE3	1:C:394:MET:HE3	1.86	0.57
1:A:440:CYS:SG	1:A:454:ILE:CD1	2.89	0.56
1:C:351:LYS:HG2	1:C:354:ARG:CZ	2.35	0.56
1:C:95:TYR:CZ	1:C:347:LEU:HD13	2.40	0.56
1:A:144:LYS:HD3	1:A:187:LYS:HZ3	1.70	0.56
1:B:142:ILE:HD13	1:B:165:LEU:HD21	1.86	0.56
1:B:155:THR:OG1	1:B:482:SER:CB	2.54	0.56
1:C:509:LEU:HD11	1:C:532:TYR:HE2	1.70	0.56
1:D:601:LEU:C	1:D:601:LEU:HD23	2.31	0.56
1:A:91:ILE:HG23	1:A:346:LEU:HD23	1.87	0.56
1:B:78:MET:HE3	1:B:163:TYR:CD2	2.41	0.56
1:C:45:MET:HE3	1:C:50:GLY:HA2	1.86	0.56
1:B:149:MET:CE	1:B:163:TYR:CG	2.84	0.56
1:C:242:ALA:HA	1:C:323:ILE:HD13	1.88	0.56
1:A:393:GLY:O	1:A:397:VAL:HG23	2.06	0.56
1:B:159:ALA:HB3	1:B:476:ALA:HB1	1.88	0.56
1:C:417:TRP:O	1:C:420:VAL:CG2	2.49	0.56
1:B:76:ALA:HB3	1:B:89:ILE:HD13	1.88	0.56
1:B:142:ILE:HD11	1:B:165:LEU:HD21	1.87	0.56
1:C:51:ILE:O	1:C:55:VAL:HG23	2.07	0.56
1:B:68:VAL:CG2	1:B:111:LEU:HD13	2.33	0.55
1:B:149:MET:HE2	1:B:160:GLN:HB3	1.89	0.55
1:C:245:ILE:HD12	1:C:322:LEU:HD23	1.88	0.55
1:C:275:VAL:HG13	1:C:322:LEU:HD22	1.89	0.55
1:D:353:TYR:O	1:D:354:ARG:C	2.50	0.55
1:B:18:ILE:CG2	1:B:18:ILE:O	2.55	0.55
1:D:398:THR:HG21	1:D:452:VAL:CG2	2.37	0.55
1:A:43:GLY:HA2	1:A:92:VAL:HG13	1.89	0.55
1:B:1:MET:HE1	1:B:9:LEU:CD2	2.32	0.54
1:C:542:THR:HG21	1:C:544:GLU:OE2	2.07	0.54
1:D:263:ASP:OD2	1:D:265:ALA:HB2	2.07	0.54
1:C:138:THR:HG23	1:C:141:ARG:H	1.73	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:139:ILE:HD11	1:C:481:LEU:HD11	1.89	0.54
1:C:378:MET:O	1:C:412:LEU:HD12	2.07	0.54
1:D:608:SER:HA	1:D:620:SER:HA	1.89	0.54
1:B:139:ILE:CD1	1:B:498:VAL:HG21	2.37	0.54
1:D:93:GLY:HA2	1:D:96:LEU:HD12	1.89	0.53
1:A:440:CYS:HG	1:A:454:ILE:HD13	1.73	0.53
1:B:420:VAL:HG23	1:B:421:ALA:N	2.23	0.53
1:A:61:THR:HG22	1:A:109:LYS:HB2	1.89	0.53
1:B:13:VAL:HG12	1:B:180:GLN:NE2	2.24	0.53
1:C:18:ILE:O	1:C:18:ILE:HG22	2.08	0.53
1:B:149:MET:HE3	1:B:163:TYR:CD1	2.41	0.53
1:D:18:ILE:O	1:D:18:ILE:CG2	2.57	0.53
1:C:105:MET:HE2	1:C:107:MET:CE	2.38	0.53
1:A:159:ALA:HA	1:A:161:VAL:HG12	1.90	0.53
1:B:149:MET:HE1	1:B:160:GLN:HB3	1.91	0.53
1:A:221:ALA:HB1	1:A:266:PRO:HB2	1.91	0.53
1:B:376:ALA:HB2	1:B:401:PHE:CE1	2.44	0.53
1:C:159:ALA:CB	1:C:476:ALA:HB1	2.38	0.53
1:B:70:TRP:CZ3	1:B:174:LEU:HD21	2.44	0.52
1:D:146:CYS:SG	1:D:147:THR:N	2.82	0.52
1:B:133:ILE:O	1:B:137:ASN:ND2	2.42	0.52
1:C:456:THR:O	1:C:460:VAL:HG23	2.09	0.52
1:D:45:MET:HE1	1:D:99:VAL:HB	1.91	0.52
1:D:464:ASN:HB2	1:D:510:MET:HE1	1.90	0.52
1:C:74:TRP:CZ3	1:C:473:ILE:HD12	2.45	0.52
1:C:613:GLU:O	1:C:614:ALA:CB	2.57	0.52
1:D:135:ARG:HD3	1:D:420:VAL:O	2.09	0.52
1:A:74:TRP:HZ3	1:A:473:ILE:HD12	1.75	0.52
1:B:46:HIS:CD2	1:B:223:MET:HG2	2.45	0.52
1:B:32:LEU:HD22	1:B:66:GLU:HB2	1.92	0.52
1:B:255:GLN:HE22	1:B:272:ARG:HA	1.74	0.52
1:D:32:LEU:HD22	1:D:66:GLU:HB2	1.91	0.52
1:B:11:ARG:O	1:B:11:ARG:HD3	2.09	0.52
1:A:444:LYS:NZ	1:A:454:ILE:HD12	2.25	0.51
1:B:51:ILE:HG12	1:B:234:MET:HE1	1.92	0.51
1:A:667:ASN:O	1:A:671:LEU:HD12	2.10	0.51
1:B:641:LEU:O	1:B:645:VAL:HG23	2.11	0.51
1:C:74:TRP:HZ3	1:C:473:ILE:HD12	1.75	0.51
1:D:157:THR:HG22	1:D:160:GLN:CG	2.40	0.51
1:A:593:VAL:HG11	1:A:642:LYS:HG2	1.93	0.51
1:B:57:VAL:HG22	1:B:67:PHE:CE1	2.45	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:223:MET:C	1:B:223:MET:SD	2.94	0.51
1:C:388:LEU:HD22	1:C:566:LEU:HB3	1.93	0.51
1:D:146:CYS:SG	1:D:153:GLU:HB2	2.50	0.51
1:C:337:PHE:CE1	1:C:343:ALA:HB1	2.45	0.51
1:C:337:PHE:CD1	1:C:343:ALA:HB1	2.46	0.51
1:A:36:TYR:HA	1:A:68:VAL:O	2.11	0.51
1:C:487:LEU:HD13	1:C:528:PHE:HB2	1.92	0.51
1:D:236:ASP:O	1:D:334:ARG:NH1	2.43	0.51
1:A:331:GLU:OE2	1:A:335:GLN:NE2	2.45	0.50
1:B:146:CYS:SG	1:B:147:THR:N	2.85	0.50
1:C:45:MET:HE3	1:C:50:GLY:CA	2.41	0.50
1:D:57:VAL:HG21	1:D:105:MET:HE1	1.91	0.50
1:D:74:TRP:CD1	1:D:86:LEU:HD11	2.46	0.50
1:A:641:LEU:O	1:A:645:VAL:HG23	2.12	0.50
1:C:146:CYS:SG	1:C:147:THR:N	2.84	0.50
1:C:283:VAL:HG11	1:C:325:GLU:HG2	1.94	0.50
1:B:73:ASP:OD2	1:B:116:SER:OG	2.30	0.50
1:C:604:HIS:C	1:C:604:HIS:HD1	2.19	0.50
1:A:105:MET:CE	1:A:110:VAL:HG21	2.42	0.50
1:B:158:ALA:C	1:B:160:GLN:H	2.19	0.50
1:C:159:ALA:HB3	1:C:476:ALA:CB	2.40	0.50
1:B:18:ILE:O	1:B:18:ILE:HG22	2.11	0.50
1:C:43:GLY:HA2	1:C:92:VAL:HG13	1.93	0.50
1:D:69:PHE:HZ	1:D:105:MET:HE1	1.74	0.50
1:D:398:THR:HG21	1:D:452:VAL:HG21	1.93	0.50
1:A:579:ILE:HG21	1:A:646:LEU:HB2	1.94	0.50
1:C:45:MET:HE1	1:C:99:VAL:CG1	2.42	0.50
1:C:601:LEU:HD23	1:C:602:ALA:N	2.26	0.50
1:D:34:ARG:NH1	1:D:174:LEU:O	2.45	0.50
1:D:425:ILE:HG22	1:D:426:THR:H	1.77	0.50
1:A:619:VAL:HG13	1:A:619:VAL:O	2.12	0.50
1:C:18:ILE:HG21	1:C:193:ARG:HD2	1.94	0.50
1:C:187:LYS:HG2	1:C:190:MET:HE2	1.92	0.50
1:C:466:ASP:OD1	1:C:466:ASP:N	2.43	0.50
1:D:1:MET:HE1	1:D:63:ALA:HA	1.93	0.50
1:D:18:ILE:HG23	1:D:19:GLN:HG2	1.94	0.50
1:B:516:VAL:HG13	1:B:537:ILE:HG21	1.94	0.49
1:C:511:LEU:O	1:C:512:SER:HB3	2.12	0.49
1:A:44:ARG:CZ	1:A:354:ARG:HD3	2.42	0.49
1:B:102:ALA:HB1	1:B:333:VAL:HA	1.95	0.49
1:C:105:MET:HE3	1:C:110:VAL:HG11	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:378:MET:HE3	1:C:394:MET:CE	2.43	0.49
1:D:180:GLN:C	1:D:181:LEU:HD12	2.38	0.49
1:D:604:HIS:ND1	1:D:604:HIS:C	2.70	0.49
1:A:290:ALA:HB1	1:A:298:TYR:O	2.13	0.49
1:A:516:VAL:O	1:A:516:VAL:HG23	2.13	0.49
1:C:223:MET:HE3	1:C:231:ALA:CB	2.42	0.49
1:D:142:ILE:CG2	1:D:161:VAL:HG22	2.43	0.49
1:D:347:LEU:HD12	1:D:350:VAL:HG22	1.94	0.49
1:D:384:LEU:HD23	1:D:675:GLU:OE1	2.12	0.49
1:D:438:VAL:HG11	1:D:668:MET:HE3	1.93	0.49
1:B:45:MET:CE	1:B:96:LEU:HD22	2.42	0.49
1:B:163:TYR:HB3	1:B:164:PRO:HD3	1.94	0.49
1:B:181:LEU:N	1:B:181:LEU:HD12	2.28	0.49
1:D:57:VAL:CG1	1:D:105:MET:HE2	2.43	0.48
1:C:362:LEU:HD12	1:C:362:LEU:O	2.13	0.48
1:C:392:GLU:OE2	1:C:652:ARG:NH2	2.47	0.48
1:A:350:VAL:HG12	1:A:354:ARG:NH2	2.29	0.48
1:D:609:ILE:HD12	1:D:636:LEU:HD11	1.96	0.48
1:B:13:VAL:CG1	1:B:180:GLN:NE2	2.76	0.48
1:D:388:LEU:HD11	1:D:446:TYR:CZ	2.47	0.48
1:A:563:ASP:OD2	1:A:577:ARG:NH1	2.35	0.48
1:B:414:LEU:HB2	1:B:456:THR:HG22	1.95	0.48
1:C:407:GLU:OE1	1:C:407:GLU:N	2.45	0.48
1:B:464:ASN:CB	1:B:510:MET:HE1	2.44	0.48
1:C:246:ARG:NH2	1:C:320:SER:OG	2.45	0.48
1:C:461:ILE:O	1:C:465:CYS:N	2.47	0.48
1:D:131:LEU:CD2	1:D:499:ILE:HG21	2.43	0.48
1:B:376:ALA:HB3	1:B:410:VAL:HG13	1.95	0.48
1:A:111:LEU:HD21	1:B:66:GLU:HG3	1.95	0.48
1:A:142:ILE:HD12	1:A:165:LEU:HD21	1.95	0.48
1:A:236:ASP:O	1:A:334:ARG:NH1	2.46	0.47
1:A:308:PHE:HA	1:A:313:VAL:CG2	2.44	0.47
1:C:273:ASN:HB3	1:C:276:LEU:HD12	1.95	0.47
1:D:646:LEU:C	1:D:646:LEU:HD23	2.39	0.47
1:D:625:GLU:O	1:D:628:VAL:HG22	2.14	0.47
1:A:376:ALA:HB2	1:A:401:PHE:CZ	2.49	0.47
1:B:388:LEU:HD11	1:B:446:TYR:OH	2.14	0.47
1:C:86:LEU:HG	1:C:90:ARG:HE	1.80	0.47
1:C:346:LEU:O	1:C:350:VAL:HG23	2.15	0.47
1:A:105:MET:HE2	1:A:107:MET:CE	2.44	0.47
1:D:510:MET:HE2	1:D:510:MET:HB3	1.83	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:653:SER:O	1:D:654:ALA:C	2.57	0.47
1:C:249:TYR:CZ	1:C:251:PRO:HG3	2.50	0.47
1:C:392:GLU:OE1	1:C:599:HIS:NE2	2.41	0.47
1:A:133:ILE:HG22	1:A:165:LEU:CD2	2.43	0.47
1:A:569:ASP:HB2	1:A:668:MET:HE1	1.97	0.47
1:B:377:CYS:HB2	1:B:513:VAL:HG21	1.97	0.47
1:D:592:PRO:O	1:D:596:VAL:HG23	2.15	0.47
1:A:388:LEU:HD11	1:A:446:TYR:CZ	2.49	0.47
1:B:171:ILE:HD13	1:B:179:CYS:SG	2.54	0.47
1:B:223:MET:SD	1:B:224:SER:N	2.88	0.47
1:A:217:LYS:O	1:A:220:GLN:HG2	2.15	0.47
1:A:310:SER:O	1:A:311:ASP:CB	2.63	0.47
1:B:468:PHE:O	1:B:471:SER:HB3	2.15	0.46
1:C:45:MET:HG3	1:C:49:GLN:HB3	1.97	0.46
1:C:591:ASN:C	1:C:591:ASN:OD1	2.57	0.46
1:D:46:HIS:ND1	1:D:223:MET:HE2	2.30	0.46
1:D:234:MET:HE3	1:D:330:LEU:HD22	1.97	0.46
1:A:505:VAL:HG22	1:A:529:ALA:HA	1.97	0.46
1:C:133:ILE:HG22	1:C:165:LEU:HD22	1.97	0.46
1:C:18:ILE:HB	1:C:207:ILE:HB	1.97	0.46
1:D:74:TRP:HD1	1:D:86:LEU:HD11	1.80	0.46
1:B:142:ILE:HD12	1:B:495:ALA:HB1	1.97	0.46
1:A:163:TYR:HB2	1:A:164:PRO:HD3	1.97	0.46
1:D:149:MET:HE3	1:D:160:GLN:CB	2.45	0.46
1:B:364:GLU:C	1:B:478:LYS:HZ1	2.23	0.46
1:D:78:MET:SD	1:D:166:MET:CE	3.03	0.46
1:D:611:ARG:NH2	1:D:619:VAL:HG11	2.31	0.46
1:B:53:LYS:HE3	1:B:53:LYS:HA	1.96	0.46
1:C:273:ASN:CB	1:C:276:LEU:HD12	2.46	0.46
1:A:123:ALA:HB1	1:A:465:CYS:SG	2.56	0.46
1:B:181:LEU:O	1:B:209:SER:HA	2.15	0.46
1:C:142:ILE:CG2	1:C:161:VAL:HG23	2.47	0.46
1:D:139:ILE:HD11	1:D:481:LEU:HD11	1.98	0.46
1:D:139:ILE:HD13	1:D:492:VAL:HG21	1.97	0.46
1:D:623:THR:HG23	1:D:626:ALA:HB3	1.98	0.46
1:D:641:LEU:O	1:D:645:VAL:HG23	2.15	0.46
1:C:493:ARG:N	1:C:497:GLN:OE1	2.48	0.45
1:D:351:LYS:HA	1:D:354:ARG:HG2	1.98	0.45
1:A:609:ILE:HD12	1:A:640:ASP:HB3	1.98	0.45
1:B:1:MET:CE	1:B:9:LEU:HD22	2.39	0.45
1:D:71:VAL:HG11	1:D:114:TRP:CZ3	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:7:TYR:CD1	1:A:24:ARG:HG2	2.52	0.45
1:A:276:LEU:CD1	1:A:318:LEU:HD21	2.47	0.45
1:A:591:ASN:HB2	1:A:592:PRO:CD	2.46	0.45
1:B:425:ILE:HG22	1:B:426:THR:H	1.81	0.45
1:C:71:VAL:HG11	1:C:114:TRP:CZ3	2.51	0.45
1:D:45:MET:CE	1:D:99:VAL:CG1	2.90	0.45
1:B:464:ASN:CG	1:B:510:MET:HE1	2.41	0.45
1:B:39:PHE:O	1:B:72:ALA:HB3	2.16	0.45
1:B:130:VAL:HG23	1:B:131:LEU:N	2.31	0.45
1:D:78:MET:HE1	1:D:163:TYR:CA	2.47	0.45
1:D:347:LEU:O	1:D:350:VAL:HG22	2.16	0.45
1:B:253:VAL:HG11	1:B:271:ASP:HB3	1.99	0.45
1:C:516:VAL:HG13	1:C:537:ILE:CG2	2.42	0.45
1:D:163:TYR:HB3	1:D:164:PRO:CD	2.46	0.45
1:A:234:MET:SD	1:A:330:LEU:HD21	2.57	0.45
1:C:163:TYR:HB2	1:C:164:PRO:HD3	1.97	0.45
1:D:15:GLU:OE1	1:D:211:HIS:ND1	2.44	0.45
1:D:45:MET:HE2	1:D:96:LEU:HD23	1.98	0.45
1:C:41:PRO:HB3	1:C:92:VAL:HG12	1.99	0.45
1:C:46:HIS:HD1	1:C:48:ALA:HB3	1.82	0.45
1:D:671:LEU:O	1:D:674:ALA:HB3	2.16	0.45
1:A:253:VAL:HG12	1:A:254:LYS:O	2.17	0.44
1:B:69:PHE:CE2	1:B:105:MET:HE1	2.52	0.44
1:B:127:TRP:HA	1:B:130:VAL:HG22	1.99	0.44
1:C:367:LEU:HD23	1:C:368:PRO:HD2	1.98	0.44
1:C:393:GLY:O	1:C:397:VAL:HG23	2.18	0.44
1:B:126:TYR:OH	1:B:170:ASP:OD1	2.35	0.44
1:C:463:GLY:C	1:C:464:ASN:OD1	2.61	0.44
1:C:464:ASN:OD1	1:C:464:ASN:N	2.50	0.44
1:C:610:GLU:O	1:C:611:ARG:HG2	2.17	0.44
1:C:627:LEU:HD23	1:C:627:LEU:O	2.17	0.44
1:D:10:LEU:CD2	1:D:27:ILE:HD11	2.46	0.44
1:A:183:LEU:O	1:A:186:ARG:HB2	2.18	0.44
1:A:624:PRO:O	1:A:628:VAL:HG23	2.17	0.44
1:B:465:CYS:O	1:B:466:ASP:C	2.60	0.44
1:A:181:LEU:HD12	1:A:181:LEU:N	2.32	0.44
1:B:223:MET:HB2	1:B:231:ALA:HB2	1.99	0.44
1:C:46:HIS:CE1	1:C:48:ALA:HB3	2.52	0.44
1:C:95:TYR:O	1:C:99:VAL:HG23	2.17	0.44
1:D:126:TYR:O	1:D:130:VAL:HG23	2.18	0.44
1:C:159:ALA:HB3	1:C:476:ALA:CA	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:232:ILE:HD13	1:C:244:LYS:HD2	2.00	0.44
1:C:245:ILE:HD12	1:C:322:LEU:CD2	2.48	0.44
1:C:307:ALA:HB1	1:C:313:VAL:HG13	2.00	0.44
1:D:46:HIS:CD2	1:D:48:ALA:HB3	2.52	0.44
1:C:279:PHE:HA	1:C:283:VAL:CG2	2.48	0.43
1:D:91:ILE:HG23	1:D:346:LEU:HD23	1.99	0.43
1:D:159:ALA:CB	1:D:476:ALA:CA	2.94	0.43
1:B:301:TYR:CD2	1:B:301:TYR:C	2.96	0.43
1:C:185:GLN:O	1:C:189:ASN:ND2	2.49	0.43
1:C:604:HIS:C	1:C:604:HIS:ND1	2.76	0.43
1:B:142:ILE:CD1	1:B:495:ALA:CB	2.96	0.43
1:D:142:ILE:HG21	1:D:161:VAL:HG22	2.00	0.43
1:B:39:PHE:N	1:B:39:PHE:CD2	2.86	0.43
1:D:417:TRP:HB3	1:D:503:MET:HE1	2.00	0.43
1:B:78:MET:HE1	1:B:166:MET:HE3	2.01	0.43
1:B:397:VAL:HG21	1:B:542:THR:HA	2.00	0.43
1:C:185:GLN:NE2	2:C:701:TYE:H	2.33	0.43
1:D:138:THR:HG23	1:D:141:ARG:H	1.83	0.43
1:D:623:THR:HG23	1:D:626:ALA:CB	2.48	0.43
1:B:340:ASN:HD21	1:B:342:GLU:HG2	1.83	0.43
1:C:376:ALA:HB2	1:C:401:PHE:CE1	2.54	0.43
1:C:398:THR:HG21	1:C:452:VAL:HG21	2.01	0.43
1:D:156:LEU:O	1:D:156:LEU:HD23	2.19	0.43
1:C:195:TYR:HA	1:C:198:LEU:HD12	2.00	0.43
1:A:571:ASN:OD1	1:A:574:ASP:N	2.41	0.43
1:B:360:LEU:CB	1:B:361:PRO:HD2	2.48	0.43
1:A:604:HIS:NE2	1:A:648:LEU:HD22	2.34	0.43
1:C:78:MET:HE1	1:C:474:GLY:CA	2.49	0.43
1:C:165:LEU:CD1	1:C:499:ILE:HD11	2.48	0.43
1:C:344:HIS:CE1	1:C:348:GLU:HG3	2.54	0.43
1:C:434:ALA:O	1:C:438:VAL:HG23	2.19	0.43
1:D:346:LEU:O	1:D:349:ALA:HB3	2.19	0.43
1:D:607:LEU:O	1:D:621:TYR:N	2.40	0.43
1:B:154:GLY:C	1:B:156:LEU:N	2.73	0.43
1:B:155:THR:O	1:B:156:LEU:C	2.61	0.43
1:B:165:LEU:HD12	1:B:499:ILE:HD11	2.01	0.43
1:C:18:ILE:HG22	1:C:207:ILE:H	1.84	0.43
1:C:46:HIS:ND1	1:C:48:ALA:HB3	2.34	0.43
1:C:601:LEU:HD12	1:C:607:LEU:HB3	2.01	0.42
1:D:376:ALA:HB2	1:D:401:PHE:CZ	2.53	0.42
1:A:223:MET:HB2	1:A:231:ALA:HB2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:216:LEU:HD21	1:B:278:TYR:CZ	2.54	0.42
1:C:43:GLY:CA	1:C:92:VAL:HG13	2.49	0.42
1:A:389:ASP:OD2	1:A:545:GLY:N	2.49	0.42
1:A:493:ARG:N	1:A:497:GLN:OE1	2.41	0.42
1:A:551:HIS:C	1:A:551:HIS:HD1	2.28	0.42
1:C:53:LYS:HE3	1:C:53:LYS:HA	2.02	0.42
1:C:111:LEU:HD12	1:C:111:LEU:N	2.34	0.42
1:C:234:MET:HE3	1:C:330:LEU:CD2	2.49	0.42
1:C:242:ALA:HA	1:C:323:ILE:CD1	2.48	0.42
1:D:223:MET:HB2	1:D:231:ALA:HB2	2.01	0.42
1:B:154:GLY:O	1:B:156:LEU:N	2.46	0.42
1:B:241:VAL:O	1:B:242:ALA:C	2.62	0.42
1:B:431:ASP:OD1	1:B:670:VAL:HG12	2.19	0.42
1:C:409:THR:CG2	1:C:410:VAL:N	2.81	0.42
1:D:38:GLY:HA2	1:D:70:TRP:O	2.19	0.42
1:D:44:ARG:HG2	1:D:354:ARG:NH1	2.34	0.42
1:A:325:GLU:O	1:A:328:ALA:HB3	2.20	0.42
1:A:601:LEU:HD23	1:A:601:LEU:O	2.19	0.42
1:C:187:LYS:HE2	1:C:263:ASP:HA	2.02	0.42
1:D:291:ALA:O	1:D:297:THR:HG23	2.18	0.42
1:A:136:GLN:HB3	1:A:198:LEU:HD13	2.02	0.42
1:C:611:ARG:NH1	1:C:616:GLY:HA3	2.34	0.42
1:A:24:ARG:HG3	1:A:24:ARG:HH11	1.85	0.42
1:A:30:LYS:HA	1:A:31:PRO:HD3	1.90	0.42
1:D:44:ARG:HG2	1:D:354:ARG:CZ	2.50	0.42
1:B:378:MET:O	1:B:412:LEU:HD12	2.20	0.42
1:B:448:LEU:HG	1:B:449:PRO:HD2	2.01	0.42
1:A:331:GLU:O	1:A:335:GLN:HG2	2.20	0.41
1:C:105:MET:CE	1:C:110:VAL:HG11	2.50	0.41
1:A:409:THR:HG22	1:A:410:VAL:N	2.35	0.41
1:C:579:ILE:O	1:C:582:ALA:HB3	2.20	0.41
1:D:44:ARG:HH21	1:D:354:ARG:HD3	1.84	0.41
1:D:659:LEU:HD22	1:D:663:GLU:HB3	2.02	0.41
1:A:111:LEU:HD11	1:B:111:LEU:HD21	2.03	0.41
1:A:444:LYS:HA	1:A:448:LEU:HB3	2.02	0.41
1:B:10:LEU:HD21	1:B:178:ILE:CD1	2.51	0.41
1:B:330:LEU:O	1:B:333:VAL:HG12	2.19	0.41
1:C:370:ALA:HB2	1:C:536:ARG:HD2	2.01	0.41
1:A:679:ALA:O	1:A:680:LYS:C	2.63	0.41
1:B:139:ILE:HD11	1:B:161:VAL:HG21	2.02	0.41
1:B:473:ILE:O	1:B:476:ALA:HB3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:165:LEU:HD12	1:C:499:ILE:HD11	2.02	0.41
1:A:377:CYS:HB2	1:A:513:VAL:HG21	2.02	0.41
1:D:472:VAL:HG13	1:D:502:LEU:HB3	2.03	0.41
1:A:214:ALA:CB	1:A:272:ARG:HD3	2.48	0.41
1:B:292:THR:HG23	1:B:296:THR:O	2.21	0.41
1:C:283:VAL:O	1:C:286:ARG:HB2	2.21	0.41
1:A:95:TYR:CE1	1:A:347:LEU:HD13	2.55	0.41
1:B:218:GLN:HA	1:B:249:TYR:CD2	2.55	0.41
1:C:156:LEU:C	1:C:156:LEU:HD12	2.46	0.41
1:C:659:LEU:O	1:C:660:LEU:C	2.64	0.41
1:D:398:THR:HG23	1:D:410:VAL:HG11	2.03	0.41
1:A:443:LEU:HD23	1:A:448:LEU:HD22	2.02	0.41
1:B:291:ALA:O	1:B:297:THR:HG23	2.21	0.41
1:C:31:PRO:HA	1:D:122:HIS:CE1	2.56	0.41
1:C:34:ARG:NH1	1:C:174:LEU:O	2.54	0.41
1:C:161:VAL:O	1:C:161:VAL:HG22	2.20	0.41
1:C:245:ILE:HG22	1:C:319:LYS:HB3	2.02	0.41
1:C:343:ALA:HA	1:C:346:LEU:HD12	2.01	0.41
1:D:349:ALA:O	1:D:352:SER:OG	2.33	0.41
1:A:26:LEU:HD12	1:A:26:LEU:HA	1.96	0.41
1:B:331:GLU:HB3	1:B:332:PRO:HD3	2.02	0.41
1:B:414:LEU:N	1:B:414:LEU:HD23	2.36	0.41
1:B:480:LEU:O	1:B:484:VAL:HG23	2.21	0.41
1:B:516:VAL:HG23	1:B:516:VAL:O	2.21	0.41
1:C:291:ALA:O	1:C:297:THR:HG23	2.21	0.41
1:C:351:LYS:HA	1:C:354:ARG:NE	2.35	0.41
1:C:480:LEU:N	1:C:480:LEU:HD12	2.36	0.41
1:D:47:ILE:HG21	1:D:278:TYR:CE1	2.55	0.41
1:D:91:ILE:HG23	1:D:346:LEU:HD22	2.01	0.41
1:D:159:ALA:HB3	1:D:476:ALA:C	2.46	0.41
1:A:161:VAL:O	1:A:161:VAL:HG22	2.21	0.40
1:A:301:TYR:CD2	1:A:301:TYR:C	2.98	0.40
1:B:30:LYS:HA	1:B:31:PRO:HD3	1.89	0.40
1:C:428:VAL:HG12	1:C:430:LYS:H	1.86	0.40
1:D:337:PHE:CE1	1:D:347:LEU:HB2	2.56	0.40
1:B:237:THR:O	1:B:241:VAL:HG23	2.21	0.40
1:D:156:LEU:O	1:D:156:LEU:CD2	2.70	0.40
1:A:373:LYS:HG2	1:A:407:GLU:HG2	2.04	0.40
1:A:380:MET:HE2	1:A:440:CYS:N	2.36	0.40
1:B:331:GLU:O	1:B:335:GLN:HG3	2.22	0.40
1:D:290:ALA:HB1	1:D:298:TYR:O	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:155:THR:OG1	1:B:482:SER:OG	2.32	0.40
1:C:79:ASN:OD1	1:C:79:ASN:N	2.55	0.40
1:D:95:TYR:O	1:D:99:VAL:HG23	2.22	0.40
1:D:367:LEU:HD12	1:D:368:PRO:HD2	2.04	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	649/690 (94%)	607 (94%)	37 (6%)	5 (1%)	16	44
1	B	573/690 (83%)	529 (92%)	40 (7%)	4 (1%)	18	48
1	C	656/690 (95%)	622 (95%)	31 (5%)	3 (0%)	24	54
1	D	659/690 (96%)	624 (95%)	30 (5%)	5 (1%)	16	44
All	All	2537/2760 (92%)	2382 (94%)	138 (5%)	17 (1%)	18	48

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	311	ASP
1	A	525	ILE
1	D	31	PRO
1	D	425	ILE
1	A	262	ASP
1	A	425	ILE
1	B	425	ILE
1	C	425	ILE
1	C	451	SER
1	C	512	SER
1	B	155	THR

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Mol	Chain	Res	Type
1	D	158	ALA
1	D	311	ASP
1	D	654	ALA
1	A	41	PRO
1	B	287	PRO
1	B	361	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	540/567 (95%)	509 (94%)	31 (6%)	18	49
1	B	453/567 (80%)	426 (94%)	27 (6%)	17	47
1	C	544/567 (96%)	516 (95%)	28 (5%)	21	53
1	D	539/567 (95%)	520 (96%)	19 (4%)	32	66
All	All	2076/2268 (92%)	1971 (95%)	105 (5%)	21	53

All (105) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	ASP
1	A	12	SER
1	A	21	SER
1	A	26	LEU
1	A	36	TYR
1	A	44	ARG
1	A	53	LYS
1	A	87	GLU
1	A	89	ILE
1	A	124	ASP
1	A	155	THR
1	A	156	LEU
1	A	157	THR
1	A	180	GLN
1	A	223	MET

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Mol	Chain	Res	Type
1	A	246	ARG
1	A	252	ARG
1	A	258	SER
1	A	276	LEU
1	A	282	VAL
1	A	373	LYS
1	A	388	LEU
1	A	433	SER
1	A	446	TYR
1	A	450	SER
1	A	513	VAL
1	A	524	HIS
1	A	572	ASP
1	A	609	ILE
1	A	611	ARG
1	A	627	LEU
1	B	4	ASP
1	B	8	LYS
1	B	18	ILE
1	B	19	GLN
1	B	36	TYR
1	B	40	GLU
1	B	53	LYS
1	B	111	LEU
1	B	115	SER
1	B	116	SER
1	B	139	ILE
1	B	149	MET
1	B	152	THR
1	B	155	THR
1	B	175	LYS
1	B	199	ILE
1	B	203	LEU
1	B	223	MET
1	B	246	ARG
1	B	276	LEU
1	B	333	VAL
1	B	341	GLU
1	B	414	LEU
1	B	450	SER
1	B	510	MET
1	B	513	VAL

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Mol	Chain	Res	Type
1	B	675	GLU
1	C	18	ILE
1	C	36	TYR
1	C	53	LYS
1	C	73	ASP
1	C	79	ASN
1	C	89	ILE
1	C	157	THR
1	C	160	GLN
1	C	203	LEU
1	C	222	LYS
1	C	223	MET
1	C	226	SER
1	C	246	ARG
1	C	282	VAL
1	C	283	VAL
1	C	342	GLU
1	C	364	GLU
1	C	388	LEU
1	C	451	SER
1	C	464	ASN
1	C	466	ASP
1	C	492	VAL
1	C	510	MET
1	C	518	SER
1	C	551	HIS
1	C	604	HIS
1	C	611	ARG
1	C	646	LEU
1	D	18	ILE
1	D	36	TYR
1	D	53	LYS
1	D	115	SER
1	D	152	THR
1	D	157	THR
1	D	160	GLN
1	D	161	VAL
1	D	246	ARG
1	D	333	VAL
1	D	420	VAL
1	D	433	SER
1	D	450	SER

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Mol	Chain	Res	Type
1	D	466	ASP
1	D	510	MET
1	D	604	HIS
1	D	609	ILE
1	D	623	THR
1	D	675	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	405	HIS
1	A	587	ASN
1	C	25	ASN
1	C	46	HIS
1	C	122	HIS
1	C	137	ASN
1	C	167	GLN
1	C	185	GLN
1	C	405	HIS
1	D	79	ASN
1	D	137	ASN
1	D	189	ASN
1	D	526	ASN
1	D	598	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 for Mogul analysis.

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	TYE	A	701	-	12,12,12	0.75	0	13,15,15	1.05	1 (7%)
2	TYE	C	701	-	12,12,12	0.98	1 (8%)	13,15,15	0.64	0
2	TYE	D	701	-	12,12,12	0.98	1 (8%)	13,15,15	0.63	0
2	TYE	B	701	-	12,12,12	0.86	0	13,15,15	0.54	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	TYE	A	701	-	-	0/6/6/6	0/1/1/1
2	TYE	C	701	-	-	3/6/6/6	0/1/1/1
2	TYE	D	701	-	-	4/6/6/6	0/1/1/1
2	TYE	B	701	-	-	2/6/6/6	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	701	TYE	C-CA	3.05	1.57	1.52
2	D	701	TYE	C-CA	2.86	1.57	1.52

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	701	TYE	CB-CA-C	-2.20	108.22	112.21

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	701	TYE	N-CA-CB-CG
2	D	701	TYE	N-CA-CB-CG
2	C	701	TYE	C-CA-CB-CG
2	D	701	TYE	C-CA-CB-CG
2	B	701	TYE	O-C-CA-CB
2	C	701	TYE	O-C-CA-CB
2	B	701	TYE	O-C-CA-N
2	D	701	TYE	CA-CB-CG-CD2
2	D	701	TYE	CA-CB-CG-CD1

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	701	TYE	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

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	657/690 (95%)	0.72	55 (8%)	17	14	60, 60, 60, 60	0
1	B	581/690 (84%)	0.60	26 (4%)	38	30	60, 60, 60, 60	0
1	C	663/690 (96%)	0.28	21 (3%)	50	41	36, 60, 60, 60	1 (0%)
1	D	665/690 (96%)	0.29	15 (2%)	61	52	60, 60, 60, 60	0
All	All	2566/2760 (92%)	0.47	117 (4%)	37	29	36, 60, 60, 60	1 (0%)

All (117) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	350	VAL	4.3
1	A	159	ALA	3.8
1	B	474	GLY	3.6
1	A	260	ILE	3.6
1	A	465	CYS	3.5
1	B	159	ALA	3.5
1	A	341	GLU	3.4
1	B	386	VAL	3.3
1	B	221	ALA	3.3
1	C	91	ILE	3.2
1	D	43	GLY	3.1
1	D	635	ALA	3.1
1	A	379	TRP	3.0
1	D	241	VAL	3.0
1	A	41	PRO	3.0
1	A	294	ASP	3.0
1	C	42	SER	3.0
1	C	82	VAL	3.0
1	B	75	PHE	2.9
1	B	161	VAL	2.9
1	B	404	ALA	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	176	ALA	2.9
1	A	605	GLY	2.8
1	B	473	ILE	2.8
1	C	346	LEU	2.8
1	B	541	GLN	2.8
1	A	86	LEU	2.8
1	A	633	SER	2.8
1	A	208	LEU	2.8
1	A	42	SER	2.8
1	C	229	ASP	2.7
1	A	47	ILE	2.7
1	A	262	ASP	2.7
1	B	232	ILE	2.7
1	B	299	ALA	2.7
1	C	176	ALA	2.6
1	C	223	MET	2.6
1	B	154	GLY	2.6
1	A	123	ALA	2.6
1	A	501	ALA	2.6
1	C	75	PHE	2.6
1	A	13	VAL	2.6
1	A	43	GLY	2.6
1	C	275	VAL	2.6
1	B	132	ASP	2.5
1	D	313	VAL	2.5
1	A	630	ASP	2.5
1	A	178	ILE	2.5
1	D	291	ALA	2.5
1	A	391	ALA	2.5
1	A	516	VAL	2.4
1	A	464	ASN	2.4
1	B	400	ASP	2.4
1	A	206	VAL	2.4
1	D	333	VAL	2.4
1	D	268	ALA	2.4
1	C	174	LEU	2.4
1	A	518	SER	2.4
1	C	226	SER	2.4
1	D	153	GLU	2.3
1	D	342	GLU	2.3
1	A	291	ALA	2.3
1	A	345	GLU	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	72	ALA	2.3
1	A	656	ALA	2.3
1	D	159	ALA	2.3
1	A	46	HIS	2.3
1	D	294	ASP	2.3
1	C	38	GLY	2.3
1	C	531[A]	GLU	2.2
1	A	181	LEU	2.2
1	B	3	THR	2.2
1	B	214	ALA	2.2
1	D	235	GLU	2.2
1	C	-1	HIS	2.2
1	B	381	PRO	2.2
1	A	48	ALA	2.2
1	A	174	LEU	2.2
1	B	172	PHE	2.2
1	A	162	LEU	2.2
1	A	452	VAL	2.2
1	D	82	VAL	2.2
1	A	601	LEU	2.2
1	A	103	ALA	2.2
1	A	508	ALA	2.2
1	A	207	ILE	2.1
1	D	661	ASN	2.1
1	A	472	VAL	2.1
1	C	267	VAL	2.1
1	A	414	LEU	2.1
1	A	163	TYR	2.1
1	A	33	ILE	2.1
1	A	39	PHE	2.1
1	A	443	LEU	2.1
1	B	111	LEU	2.1
1	A	19	GLN	2.1
1	A	180	GLN	2.1
1	B	313	VAL	2.1
1	C	78	MET	2.1
1	A	155	THR	2.1
1	A	110	VAL	2.1
1	A	275	VAL	2.1
1	B	472	VAL	2.1
1	A	183	LEU	2.1
1	C	77	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	308	PHE	2.1
1	B	427	GLY	2.1
1	C	227	ASP	2.1
1	A	158	ALA	2.1
1	A	115	SER	2.0
1	A	282	VAL	2.0
1	B	401	PHE	2.0
1	C	43	GLY	2.0
1	C	660	LEU	2.0
1	B	200	GLY	2.0
1	D	259	ALA	2.0
1	C	455	VAL	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	NA	C	711	1/1	0.63	0.22	119,119,119,119	0
3	NA	A	711	1/1	0.71	0.27	68,68,68,68	0
2	TYE	D	701	12/12	0.75	0.12	60,60,60,60	0
2	TYE	C	701	12/12	0.85	0.09	60,60,60,60	0
2	TYE	B	701	12/12	0.88	0.15	60,60,60,60	0
2	TYE	A	701	12/12	0.93	0.16	60,60,60,60	0

6.5 Other polymers [i](#)

There are no such residues in this entry.