



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:31 AM UTC

PDB ID : 3IE2 / pdb_00003ie2
Title : Crystal Structure of H400V mutant TTHA0252 from *Thermus thermophilus* HB8
Authors : Ishikawa, H.; Nakagawa, N.; Kuramitsu, S.; Yokoyama, S.; Masui, R.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2009-07-22
Resolution : 2.80 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

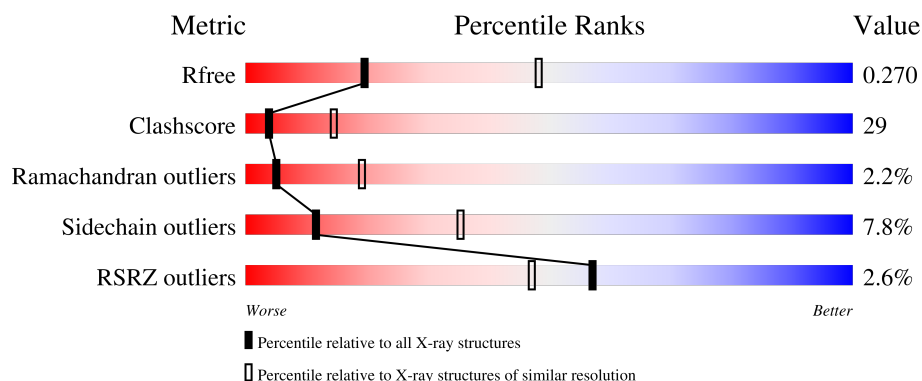
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	431	
1	B	431	
1	C	431	
1	D	431	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 13713 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 Ribonuclease TTHA0252.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	431	Total	C	N	O	S	0	0	0
			3323	2126	595	594	8			
1	B	431	Total	C	N	O	S	0	0	0
			3323	2126	595	594	8			
1	C	431	Total	C	N	O	S	0	0	0
			3323	2126	595	594	8			
1	D	431	Total	C	N	O	S	0	0	0
			3323	2126	595	594	8			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	400	VAL	HIS	engineered mutation	UNP Q5SLP1
B	400	VAL	HIS	engineered mutation	UNP Q5SLP1
C	400	VAL	HIS	engineered mutation	UNP Q5SLP1
D	400	VAL	HIS	engineered mutation	UNP Q5SLP1

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

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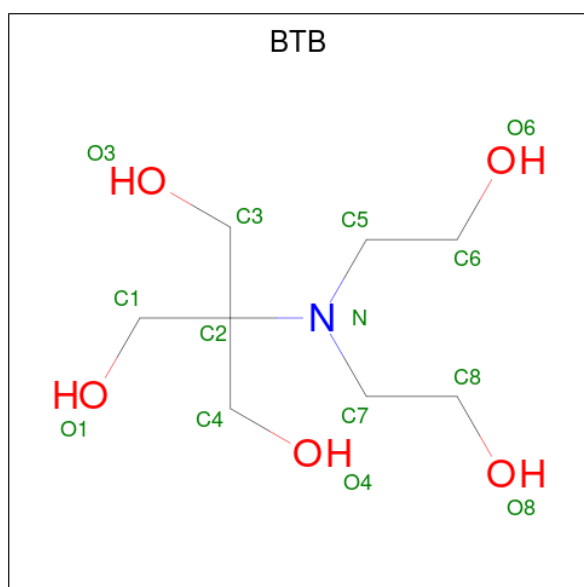
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is 2-[BIS-(2-HYDROXY-ETHYL)-AMINO]-2-HYDROXYMETHYL-PROPAN E-1,3-DIOL (CCD ID: BTB) (formula: $C_8H_{19}NO_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	N	O	
			14	8	1	5	

- Molecule 4 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total 1	Zn 1	0	0
4	B	1	Total 1	Zn 1	0	0
4	C	1	Total 1	Zn 1	0	0
4	D	1	Total 1	Zn 1	0	0

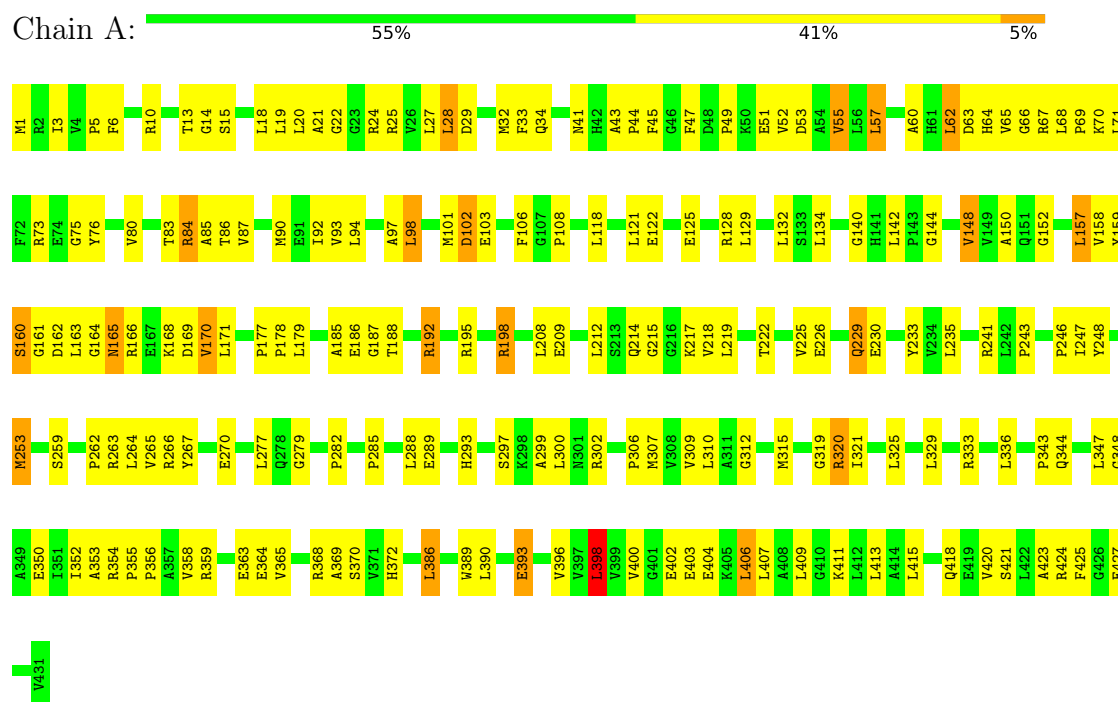
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	27	Total 27	O 27	0	0
5	B	14	Total 14	O 14	0	0
5	C	18	Total 18	O 18	0	0
5	D	19	Total 19	O 19	0	0

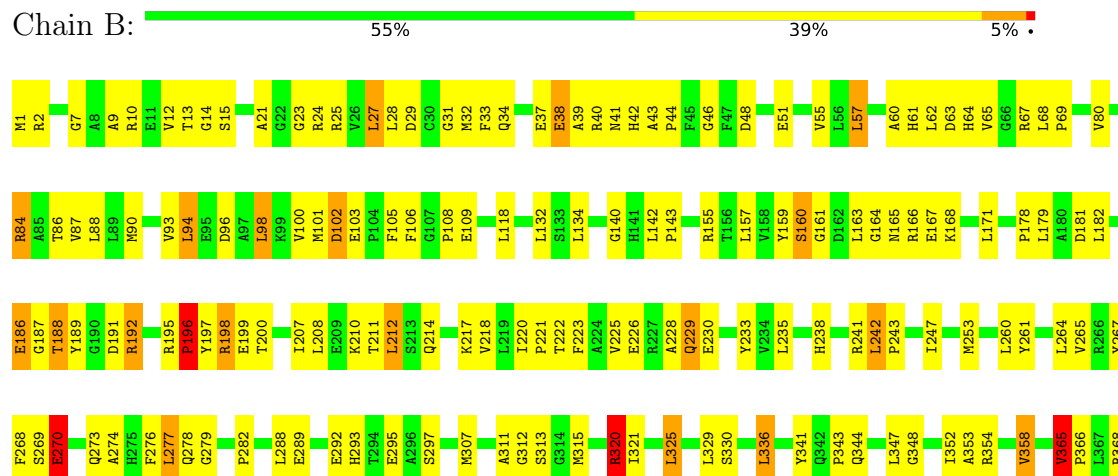
3 Residue-property plots

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

• Molecule 1: Ribonuclease TTHA0252

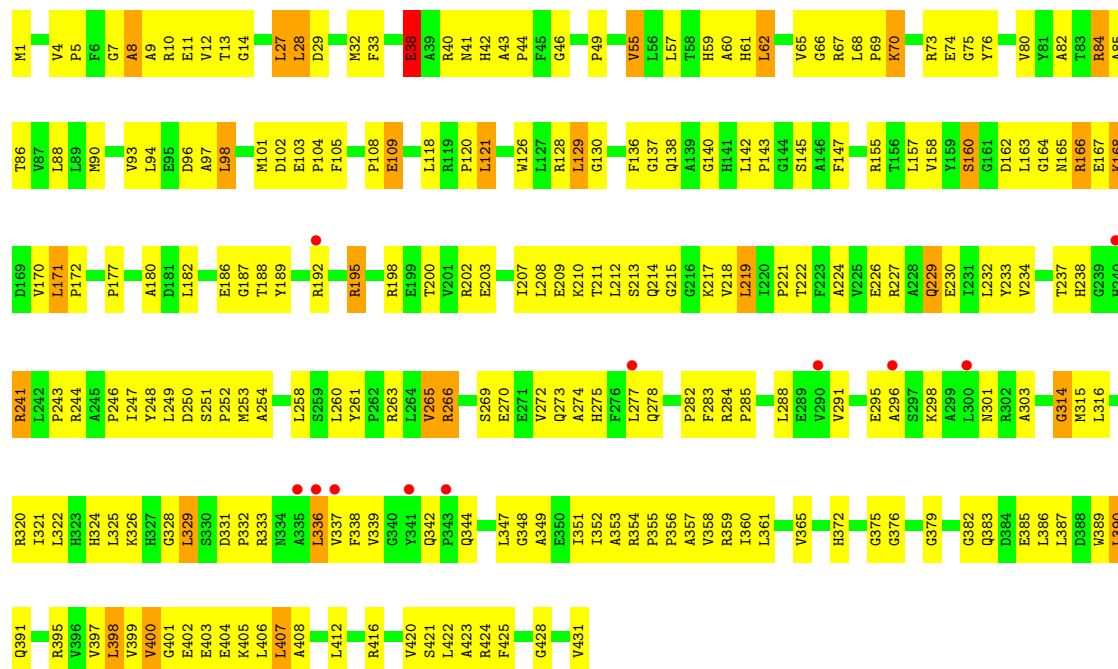


• Molecule 1: Ribonuclease TTHA0252

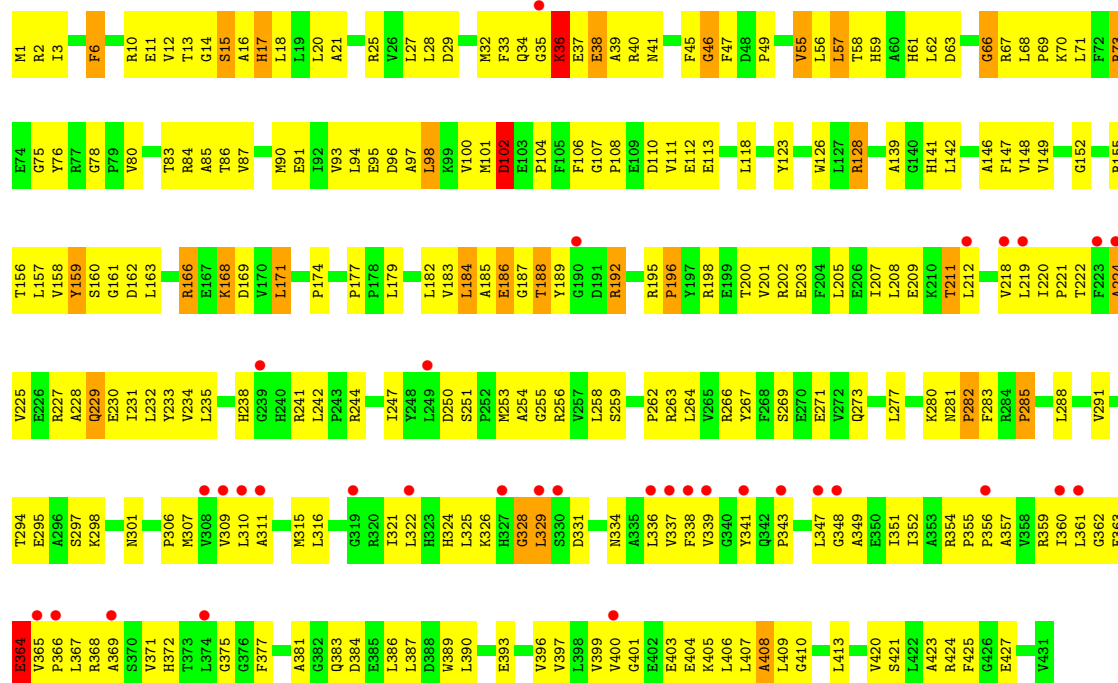




• Molecule 1: Ribonuclease TTHA0252



• Molecule 1: Ribonuclease TTHA0252



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	143.44Å 149.74Å 122.12Å 90.00° 110.61° 90.00°	Depositor
Resolution (Å)	50.00 – 2.80 50.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	89.0 (50.00-2.80) 89.3 (50.00-2.80)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.01 (at 2.81Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.219 , 0.275 0.214 , 0.270	Depositor DCC
R_{free} test set	5608 reflections (10.15%)	wwPDB-VP
Wilson B-factor (Å ²)	54.0	Xtriage
Anisotropy	0.084	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 54.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	13713	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 3.85% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, BTB, SO4

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	0.51	0/3403	1.04	18/4616 (0.4%)
1	B	0.50	0/3403	1.01	17/4616 (0.4%)
1	C	0.41	0/3403	0.96	10/4616 (0.2%)
1	D	0.37	0/3403	0.92	7/4616 (0.2%)
All	All	0.45	0/13612	0.99	52/18464 (0.3%)

There are no bond length outliers.

All (52) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	225	VAL	N-CA-C	9.92	119.86	110.53
1	A	170	VAL	N-CA-C	8.88	118.88	110.53
1	C	60	ALA	N-CA-C	7.66	122.02	113.21
1	D	46	GLY	N-CA-C	-7.58	105.47	114.48
1	B	160	SER	N-CA-C	7.16	119.09	111.28
1	B	293	HIS	N-CA-C	7.03	119.81	109.69
1	A	192	ARG	N-CA-C	6.92	119.44	108.23
1	C	314	GLY	N-CA-C	-6.79	106.42	115.32
1	C	398	LEU	N-CA-C	6.45	119.60	110.50
1	C	103	GLU	CA-C-N	6.42	126.39	119.78
1	C	103	GLU	C-N-CA	6.42	126.39	119.78
1	A	160	SER	N-CA-C	6.40	119.56	111.69
1	C	46	GLY	N-CA-C	-6.34	106.19	115.63
1	D	75	GLY	N-CA-C	6.31	122.20	114.69
1	A	185	ALA	N-CA-C	6.19	119.74	110.14
1	A	60	ALA	N-CA-C	6.11	120.59	113.20
1	A	393	GLU	CA-C-N	6.08	125.80	119.05
1	A	393	GLU	C-N-CA	6.08	125.80	119.05
1	B	46	GLY	N-CA-C	-6.04	106.30	115.00
1	A	161	GLY	N-CA-C	-5.94	101.25	111.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	320	ARG	N-CA-C	5.90	118.47	111.33
1	A	398	LEU	N-CA-C	5.82	118.72	109.24
1	D	159	TYR	N-CA-C	-5.74	102.81	110.55
1	A	225	VAL	N-CA-C	5.71	116.35	110.82
1	C	265	VAL	N-CA-C	5.70	116.34	110.36
1	D	169	ASP	N-CA-C	5.56	121.49	114.31
1	B	181	ASP	N-CA-C	-5.53	105.66	112.90
1	B	398	LEU	N-CA-C	5.52	118.45	110.28
1	B	292	GLU	N-CA-C	5.50	118.10	111.40
1	A	293	HIS	N-CA-C	5.47	118.62	110.14
1	B	365	VAL	CA-C-N	5.45	125.38	119.76
1	B	365	VAL	C-N-CA	5.45	125.38	119.76
1	D	393	GLU	CA-C-N	5.42	124.60	118.97
1	D	393	GLU	C-N-CA	5.42	124.60	118.97
1	B	270	GLU	N-CA-C	-5.41	105.46	111.36
1	D	156	THR	N-CA-C	5.34	117.11	108.41
1	A	195	ARG	N-CA-C	-5.32	103.07	109.83
1	C	160	SER	N-CA-C	5.30	117.48	111.11
1	B	242	LEU	N-CA-C	5.29	116.32	109.65
1	A	222	THR	N-CA-C	5.20	116.98	108.76
1	A	66	GLY	N-CA-C	5.16	120.13	113.27
1	C	296	ALA	N-CA-C	-5.14	106.85	113.23
1	B	167	GLU	N-CA-C	5.14	119.18	113.01
1	C	70	LYS	N-CA-C	-5.12	105.59	111.07
1	B	223	PHE	N-CA-C	-5.12	102.11	110.20
1	A	336	LEU	N-CA-C	-5.09	100.94	109.24
1	A	169	ASP	N-CA-C	5.08	120.47	113.97
1	B	336	LEU	N-CA-C	-5.07	101.04	109.46
1	B	31	GLY	N-CA-C	5.04	119.42	112.37
1	A	148	VAL	N-CA-C	5.03	115.83	108.58
1	A	166	ARG	N-CA-C	5.03	118.88	112.34
1	B	353	ALA	N-CA-C	-5.02	106.42	112.54

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3323	0	3353	141	0
1	B	3323	0	3353	168	0
1	C	3323	0	3353	221	0
1	D	3323	0	3353	255	0
2	A	120	0	0	0	0
2	B	95	0	0	3	0
2	C	70	0	0	2	0
2	D	40	0	0	1	0
3	A	14	0	19	6	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	27	0	0	0	0
5	B	14	0	0	0	0
5	C	18	0	0	3	0
5	D	19	0	0	5	0
All	All	13713	0	13431	772	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 (772) 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:C:33:PHE:N	1:C:41:ASN:HD21	1.60	0.99
1:C:33:PHE:H	1:C:41:ASN:ND2	1.61	0.98
1:C:298:LYS:HA	1:C:301:ASN:ND2	1.77	0.97
1:C:298:LYS:HA	1:C:301:ASN:HD22	1.28	0.96
1:D:220:ILE:HG12	1:D:337:VAL:HB	1.49	0.95
1:D:198:ARG:NH2	1:D:202:ARG:HH21	1.65	0.94
1:C:10:ARG:HH22	1:C:424:ARG:HE	0.96	0.91
1:D:211:THR:HG23	1:D:212:LEU:H	1.34	0.91
1:C:224:ALA:HB3	1:C:253:MET:HE2	1.53	0.91
1:A:32:MET:HE1	1:A:101:MET:HE1	1.51	0.90
1:B:160:SER:HB2	1:B:163:LEU:HD21	1.52	0.89
1:B:208:LEU:HD23	1:B:218:VAL:HG21	1.52	0.89
1:B:315:MET:HE2	1:B:343:PRO:HD3	1.53	0.89
1:B:33:PHE:H	1:B:41:ASN:HD21	1.23	0.87
1:C:207:ILE:O	1:C:211:THR:HG22	1.75	0.87
1:D:221:PRO:HB3	1:D:321:ILE:HG12	1.57	0.86
1:C:168:LYS:HE3	1:C:230:GLU:OE1	1.75	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:224:ALA:HB3	1:D:253:MET:HE2	1.56	0.86
1:C:10:ARG:NH2	1:C:424:ARG:HE	1.74	0.85
1:C:232:LEU:HD11	1:C:249:LEU:HD22	1.56	0.85
1:B:208:LEU:O	1:B:212:LEU:HB2	1.79	0.83
1:D:228:ALA:HB3	1:D:229:GLN:NE2	1.95	0.81
1:D:13:THR:HG21	1:D:34:GLN:HB2	1.64	0.80
1:B:37:GLU:HB3	1:B:40:ARG:HG3	1.64	0.79
1:B:182:LEU:HD11	1:B:397:VAL:HG23	1.63	0.79
1:C:222:THR:HG22	1:C:339:VAL:HG21	1.63	0.79
1:C:202:ARG:HH21	1:C:202:ARG:HG2	1.50	0.77
1:D:207:ILE:O	1:D:211:THR:HG22	1.84	0.77
1:A:320:ARG:HH11	3:A:456:BTB:H31	1.51	0.75
1:D:57:LEU:HD21	1:D:68:LEU:HD22	1.69	0.75
1:D:17:HIS:HD2	1:D:184:LEU:HD21	1.52	0.75
1:D:359:ARG:NH1	1:D:362:GLY:HA2	2.00	0.75
1:D:227:ARG:HG2	1:D:227:ARG:HH21	1.51	0.74
1:A:309:VAL:C	1:A:310:LEU:HD12	2.11	0.74
1:D:20:LEU:CD2	1:D:25:ARG:HG2	2.17	0.74
1:A:160:SER:HB2	1:A:163:LEU:HD21	1.69	0.74
1:A:1:MET:HG3	1:A:21:ALA:HB2	1.70	0.73
1:B:192:ARG:HG3	1:B:192:ARG:HH11	1.54	0.73
1:D:195:ARG:NH2	1:D:375:GLY:H	1.87	0.73
1:C:101:MET:SD	1:C:104:PRO:HA	2.29	0.73
1:D:331:ASP:HB3	1:D:334:ASN:ND2	2.02	0.73
1:B:33:PHE:H	1:B:41:ASN:ND2	1.86	0.72
1:A:33:PHE:H	1:A:41:ASN:HD21	1.37	0.72
1:B:33:PHE:N	1:B:41:ASN:HD21	1.86	0.72
1:D:315:MET:SD	1:D:343:PRO:HD3	2.29	0.72
1:A:45:PHE:HB3	1:A:47:PHE:CE1	2.24	0.72
1:A:102:ASP:CG	1:A:103:GLU:H	1.98	0.71
1:D:97:ALA:O	1:D:101:MET:HB2	1.89	0.71
1:A:208:LEU:HD23	1:A:218:VAL:HG21	1.71	0.71
1:C:12:VAL:HG23	1:C:13:THR:HG23	1.72	0.71
1:D:17:HIS:CD2	1:D:184:LEU:HD21	2.25	0.71
1:D:36:LYS:H	1:D:36:LYS:HD2	1.56	0.71
1:B:179:LEU:HD13	1:C:128:ARG:HH12	1.56	0.71
1:D:250:ASP:HB3	1:D:311:ALA:HB2	1.72	0.70
1:C:10:ARG:HH22	1:C:424:ARG:NE	1.81	0.70
1:B:195:ARG:NH2	1:B:375:GLY:H	1.88	0.70
1:D:195:ARG:HH21	1:D:375:GLY:H	1.39	0.69
1:D:298:LYS:HA	1:D:301:ASN:ND2	2.08	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:62:LEU:HD13	1:D:93:VAL:HG12	1.75	0.69
1:B:1:MET:HG3	1:B:21:ALA:HB2	1.76	0.68
1:A:98:LEU:HD11	1:A:108:PRO:HA	1.75	0.68
1:A:219:LEU:HD23	1:A:325:LEU:HD12	1.74	0.68
1:A:320:ARG:HD3	3:A:456:BTB:H81	1.75	0.68
1:C:360:ILE:HG22	1:C:361:LEU:HD12	1.76	0.68
1:B:179:LEU:HD13	1:C:128:ARG:NH1	2.09	0.68
1:C:160:SER:HB2	1:C:163:LEU:HD11	1.75	0.68
1:B:221:PRO:HB3	1:B:321:ILE:HG12	1.76	0.68
1:B:32:MET:HA	1:B:67:ARG:HG3	1.76	0.67
1:D:263:ARG:O	1:D:263:ARG:HG2	1.93	0.67
1:B:57:LEU:HD23	1:B:90:MET:HE1	1.76	0.67
1:C:348:GLY:O	1:C:352:ILE:HG13	1.94	0.67
1:B:404:GLU:H	1:B:404:GLU:CD	2.02	0.66
1:D:244:ARG:H	1:D:244:ARG:HD2	1.59	0.66
1:C:229:GLN:H	1:C:229:GLN:NE2	1.94	0.66
1:B:86:THR:HG22	1:B:90:MET:HE3	1.78	0.66
1:B:98:LEU:HD11	1:B:108:PRO:HA	1.77	0.66
1:D:187:GLY:HA2	1:D:386:LEU:HD21	1.77	0.66
1:B:37:GLU:O	1:B:39:ALA:N	2.30	0.65
1:D:168:LYS:HE2	1:D:230:GLU:OE1	1.96	0.65
1:D:171:LEU:H	1:D:171:LEU:HD22	1.62	0.65
1:A:3:ILE:HD12	1:A:19:LEU:HA	1.76	0.65
1:C:285:PRO:HD2	1:C:288:LEU:HD22	1.79	0.65
1:D:183:VAL:HG12	1:D:183:VAL:O	1.97	0.65
1:A:128:ARG:HH12	1:D:179:LEU:HD13	1.61	0.65
1:C:59:HIS:CE1	1:C:162:ASP:HB2	2.32	0.65
1:C:355:PRO:HB2	1:C:356:PRO:HD2	1.78	0.65
1:B:168:LYS:HG2	1:B:197:TYR:CD2	2.33	0.64
1:A:413:LEU:HD22	1:A:418:GLN:OE1	1.97	0.64
1:C:230:GLU:O	1:C:234:VAL:HG23	1.97	0.64
1:A:319:GLY:HA2	3:A:456:BTB:H82	1.77	0.64
1:B:179:LEU:HB2	1:C:128:ARG:HH11	1.62	0.64
1:C:420:VAL:HG22	1:C:421:SER:N	2.13	0.64
1:D:355:PRO:O	1:D:367:LEU:HD23	1.98	0.64
1:A:424:ARG:HD3	1:A:427:GLU:OE1	1.97	0.64
1:C:43:ALA:HB1	1:C:44:PRO:HD2	1.79	0.64
1:D:37:GLU:O	1:D:39:ALA:N	2.31	0.64
1:D:33:PHE:H	1:D:41:ASN:HD21	1.45	0.64
1:D:155:ARG:HH11	1:D:155:ARG:HG2	1.63	0.64
1:C:229:GLN:H	1:C:229:GLN:CD	2.06	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:309:VAL:C	1:D:310:LEU:HD12	2.22	0.63
1:C:219:LEU:N	1:C:219:LEU:HD12	2.14	0.63
1:C:252:PRO:HD2	5:C:457:HOH:O	1.99	0.63
1:C:326:LYS:HA	1:C:365:VAL:HG11	1.79	0.63
1:A:32:MET:HA	1:A:67:ARG:HG3	1.81	0.63
1:C:10:ARG:CZ	1:C:424:ARG:HG2	2.29	0.63
1:C:10:ARG:NH1	1:C:424:ARG:HG2	2.13	0.63
1:C:195:ARG:NH2	1:C:375:GLY:H	1.97	0.63
1:D:37:GLU:HB3	1:D:40:ARG:HG3	1.81	0.63
1:D:3:ILE:HD12	1:D:18:LEU:O	1.99	0.62
1:C:163:LEU:HD12	1:C:163:LEU:N	2.15	0.62
1:B:168:LYS:HE2	1:B:230:GLU:OE1	1.99	0.62
1:D:12:VAL:HG23	1:D:13:THR:N	2.15	0.62
1:D:186:GLU:HA	1:D:399:VAL:O	1.99	0.62
1:C:140:GLY:O	1:C:164:GLY:HA3	2.00	0.62
1:D:102:ASP:O	1:D:104:PRO:HD3	1.99	0.62
1:B:189:TYR:OH	1:B:341:TYR:HB2	1.99	0.62
1:C:260:LEU:HD12	1:C:263:ARG:HD3	1.80	0.62
1:D:17:HIS:ND1	1:D:17:HIS:N	2.46	0.62
1:D:231:ILE:HG21	1:D:310:LEU:HD21	1.80	0.62
1:A:97:ALA:O	1:A:101:MET:HB2	2.00	0.62
1:A:217:LYS:HE2	1:A:307:MET:HE3	1.82	0.61
1:B:195:ARG:HG3	1:B:195:ARG:HH11	1.64	0.61
1:C:221:PRO:HB3	1:C:321:ILE:HG12	1.82	0.61
1:C:316:LEU:HD11	1:C:338:PHE:HE1	1.65	0.61
1:B:295:GLU:OE2	1:B:295:GLU:N	2.23	0.61
1:B:315:MET:CE	1:B:343:PRO:HD3	2.26	0.61
1:D:16:ALA:C	1:D:17:HIS:ND1	2.57	0.61
1:D:359:ARG:CA	1:D:364:GLU:HG3	2.29	0.61
1:B:43:ALA:HB1	1:B:44:PRO:HD2	1.80	0.61
1:C:214:GLN:NE2	1:C:333:ARG:HA	2.15	0.61
1:C:359:ARG:HG2	1:C:359:ARG:HH11	1.65	0.61
1:D:250:ASP:HA	1:D:291:VAL:HB	1.82	0.61
1:D:90:MET:HE1	1:D:118:LEU:HD22	1.81	0.61
1:B:68:LEU:N	1:B:69:PRO:HD2	2.16	0.60
1:B:200:THR:HG23	1:B:374:LEU:HB3	1.82	0.60
1:C:202:ARG:HG2	1:C:202:ARG:NH2	2.12	0.60
1:B:192:ARG:HH11	1:B:192:ARG:CG	2.14	0.60
1:D:357:ALA:HB1	1:D:365:VAL:O	2.00	0.60
1:A:85:ALA:HB2	1:A:267:TYR:CD2	2.37	0.60
1:C:269:SER:O	1:C:273:GLN:HG3	2.00	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:168:LYS:HE3	1:C:230:GLU:CD	2.27	0.60
1:D:211:THR:HG21	1:D:218:VAL:HG22	1.83	0.60
1:A:68:LEU:N	1:A:69:PRO:HD2	2.17	0.60
1:B:28:LEU:O	1:B:29:ASP:HB2	2.01	0.60
1:D:36:LYS:HD2	1:D:36:LYS:N	2.15	0.60
1:D:208:LEU:HD23	1:D:218:VAL:HG11	1.84	0.60
1:D:139:ALA:O	1:D:174:PRO:HG3	2.02	0.60
1:C:360:ILE:HD12	1:C:365:VAL:HG21	1.84	0.59
1:C:155:ARG:HD3	1:C:431:VAL:O	2.02	0.59
1:C:322:LEU:HB3	1:C:361:LEU:HD11	1.84	0.59
1:D:163:LEU:HD11	1:D:389:TRP:CE2	2.37	0.59
1:C:32:MET:HA	1:C:67:ARG:HG3	1.84	0.59
1:D:198:ARG:HH22	1:D:202:ARG:HH21	1.47	0.59
1:D:309:VAL:O	1:D:310:LEU:HD12	2.02	0.59
1:B:192:ARG:HG2	2:B:441:SO4:S	2.41	0.59
1:C:316:LEU:HD11	1:C:338:PHE:CE1	2.37	0.59
1:C:62:LEU:HD13	1:C:93:VAL:CG1	2.32	0.59
1:C:157:LEU:HD12	1:C:182:LEU:O	2.03	0.59
1:A:235:LEU:HD13	1:A:247:ILE:HD13	1.85	0.59
1:D:123:TYR:HE1	1:D:146:ALA:HB2	1.67	0.59
1:D:219:LEU:HD23	1:D:325:LEU:HD12	1.84	0.59
1:A:198:ARG:HG2	1:A:198:ARG:NH2	2.16	0.59
1:B:88:LEU:HD13	1:B:260:LEU:HD11	1.85	0.59
1:B:229:GLN:CD	1:B:229:GLN:H	2.08	0.59
1:A:404:GLU:CD	1:A:404:GLU:H	2.10	0.59
1:C:11:GLU:OE1	1:C:33:PHE:HE1	1.87	0.58
1:D:285:PRO:HD2	1:D:288:LEU:HD22	1.84	0.58
1:A:157:LEU:HG	1:A:158:VAL:N	2.17	0.58
1:B:220:ILE:HG22	1:B:222:THR:HG23	1.84	0.58
1:C:97:ALA:O	1:C:101:MET:HB2	2.04	0.58
1:D:91:GLU:O	1:D:95:GLU:HG2	2.03	0.58
1:C:251:SER:HB3	1:C:254:ALA:HB3	1.84	0.58
1:D:61:HIS:HB3	1:D:63:ASP:OD1	2.03	0.58
1:B:23:GLY:O	1:B:24:ARG:HD3	2.04	0.58
1:B:398:LEU:HB3	1:B:406:LEU:HG	1.85	0.58
1:C:195:ARG:CZ	1:C:375:GLY:H	2.16	0.58
1:D:45:PHE:HB3	1:D:47:PHE:CE1	2.38	0.58
1:B:188:THR:HG22	1:B:189:TYR:CD1	2.38	0.58
1:D:171:LEU:HD22	1:D:171:LEU:N	2.19	0.58
1:B:57:LEU:HG	1:B:65:VAL:HG22	1.85	0.58
1:B:277:LEU:C	1:B:279:GLY:H	2.12	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:7:GLY:HA3	1:C:14:GLY:O	2.03	0.58
1:B:182:LEU:HD11	1:B:397:VAL:CG2	2.31	0.58
1:D:363:GLU:O	1:D:365:VAL:HG13	2.04	0.58
1:A:229:GLN:H	1:A:229:GLN:CD	2.12	0.58
1:B:210:LYS:O	1:B:214:GLN:HG2	2.04	0.58
1:B:221:PRO:HA	1:B:311:ALA:O	2.03	0.57
1:D:59:HIS:CD2	1:D:61:HIS:HB2	2.39	0.57
1:D:195:ARG:NH2	1:D:375:GLY:N	2.51	0.57
1:A:128:ARG:NH1	1:D:179:LEU:HB2	2.19	0.57
1:B:165:ASN:HD22	1:B:165:ASN:C	2.12	0.57
1:B:168:LYS:HG2	1:B:197:TYR:CE2	2.38	0.57
1:B:178:PRO:HB3	1:C:126:TRP:CE3	2.39	0.57
1:D:10:ARG:HG3	1:D:10:ARG:HH11	1.69	0.57
1:D:33:PHE:HB3	1:D:37:GLU:HB2	1.86	0.57
1:D:49:PRO:HB3	1:D:71:LEU:HD12	1.87	0.57
1:A:262:PRO:O	1:A:265:VAL:HG23	2.04	0.57
1:A:310:LEU:HD12	1:A:310:LEU:N	2.19	0.57
1:B:191:ASP:OD1	1:B:405:LYS:HE2	2.05	0.57
1:C:171:LEU:HD21	1:C:226:GLU:OE1	2.04	0.57
1:D:258:LEU:HD11	1:D:283:PHE:HB3	1.87	0.57
1:D:396:VAL:HG12	1:D:397:VAL:N	2.19	0.57
1:A:55:VAL:HG13	1:A:80:VAL:HG22	1.85	0.57
1:C:164:GLY:HA2	1:C:379:GLY:O	2.04	0.57
1:D:55:VAL:C	1:D:56:LEU:HD12	2.29	0.57
1:C:211:THR:HG21	1:C:218:VAL:HG22	1.87	0.57
1:A:122:GLU:O	1:A:125:GLU:HB2	2.04	0.56
1:D:266:ARG:HG3	1:D:267:TYR:N	2.20	0.56
1:B:409:LEU:HD22	1:B:413:LEU:HG	1.87	0.56
1:A:263:ARG:HG2	1:D:277:LEU:HD21	1.88	0.56
1:A:266:ARG:HG3	1:A:267:TYR:N	2.21	0.56
1:B:61:HIS:HB3	1:B:63:ASP:OD1	2.06	0.56
1:D:16:ALA:HA	1:D:29:ASP:O	2.04	0.56
1:A:170:VAL:HG21	1:A:230:GLU:HG3	1.88	0.56
1:D:329:LEU:HD13	1:D:369:ALA:HB3	1.88	0.56
1:D:6:PHE:HB2	1:D:16:ALA:O	2.06	0.56
1:D:20:LEU:HD23	1:D:25:ARG:HG2	1.87	0.56
1:D:359:ARG:HA	1:D:364:GLU:HA	1.87	0.56
1:C:230:GLU:O	1:C:233:TYR:HB3	2.05	0.56
1:B:48:ASP:OD2	1:B:51:GLU:HG2	2.05	0.56
1:B:100:VAL:HG12	1:B:100:VAL:O	2.05	0.56
1:D:28:LEU:O	1:D:29:ASP:HB2	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:55:VAL:HG22	1:D:80:VAL:HG13	1.87	0.56
1:B:109:GLU:H	1:B:109:GLU:CD	2.14	0.56
1:C:422:LEU:N	1:C:422:LEU:HD12	2.21	0.56
1:B:297:SER:OG	1:B:320:ARG:HG2	2.06	0.56
1:D:192:ARG:NH1	5:D:456:HOH:O	2.38	0.56
1:A:270:GLU:CD	1:D:266:ARG:HH12	2.13	0.55
1:C:171:LEU:N	1:C:171:LEU:HD22	2.21	0.55
1:D:383:GLN:O	1:D:387:LEU:HG	2.05	0.55
1:C:314:GLY:O	1:C:342:GLN:NE2	2.40	0.55
1:D:183:VAL:O	1:D:185:ALA:N	2.39	0.55
1:B:33:PHE:HB2	1:B:41:ASN:ND2	2.22	0.55
1:B:98:LEU:HD11	1:B:108:PRO:CA	2.36	0.55
1:D:12:VAL:HG23	1:D:13:THR:H	1.71	0.55
1:D:229:GLN:NE2	1:D:229:GLN:H	2.04	0.55
1:B:161:GLY:O	1:B:186:GLU:HG2	2.06	0.55
1:A:348:GLY:O	1:A:352:ILE:HG13	2.05	0.55
1:D:69:PRO:HB3	1:D:111:VAL:HA	1.88	0.55
1:A:198:ARG:HG2	1:A:198:ARG:HH21	1.70	0.55
1:D:87:VAL:HA	1:D:90:MET:CE	2.37	0.55
1:D:160:SER:HB2	1:D:163:LEU:HD21	1.88	0.55
1:B:13:THR:HG21	1:B:34:GLN:HB2	1.88	0.55
1:D:359:ARG:HA	1:D:364:GLU:HG3	1.88	0.55
1:C:214:GLN:HE21	1:C:333:ARG:HA	1.72	0.55
1:D:155:ARG:HG2	1:D:155:ARG:NH1	2.22	0.55
1:D:20:LEU:HD22	1:D:25:ARG:HG2	1.90	0.54
1:D:107:GLY:O	1:D:111:VAL:HG23	2.08	0.54
1:B:10:ARG:HH12	1:B:424:ARG:NH2	2.04	0.54
1:D:3:ILE:HD12	1:D:18:LEU:C	2.31	0.54
1:B:265:VAL:HG23	1:B:276:PHE:CD2	2.42	0.54
1:C:420:VAL:HG22	1:C:421:SER:H	1.72	0.54
1:D:208:LEU:CD2	1:D:218:VAL:HG11	2.38	0.54
1:D:326:LYS:HE3	1:D:363:GLU:HG2	1.90	0.54
1:A:420:VAL:HG22	1:A:421:SER:N	2.22	0.54
1:B:55:VAL:CG1	1:B:80:VAL:HG22	2.38	0.54
1:B:393:GLU:O	1:B:418:GLN:HG2	2.07	0.54
1:C:98:LEU:HD11	1:C:108:PRO:HA	1.90	0.54
1:A:288:LEU:HD12	1:A:289:GLU:H	1.73	0.54
1:D:205:LEU:HD11	1:D:238:HIS:CG	2.43	0.54
1:B:160:SER:HB2	1:B:163:LEU:CD2	2.33	0.54
1:B:195:ARG:HG3	1:B:195:ARG:NH1	2.22	0.54
1:C:211:THR:HG23	1:C:212:LEU:H	1.72	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:141:HIS:O	1:D:142:LEU:HD23	2.07	0.54
1:D:354:ARG:HD2	1:D:367:LEU:HD21	1.90	0.54
1:B:348:GLY:O	1:B:352:ILE:HG13	2.08	0.54
1:D:83:THR:O	1:D:87:VAL:HG23	2.08	0.53
1:B:208:LEU:HD21	1:B:218:VAL:HG11	1.89	0.53
1:B:217:LYS:HE2	1:B:307:MET:HE3	1.90	0.53
1:C:211:THR:HG23	1:C:212:LEU:N	2.24	0.53
1:C:406:LEU:HB3	1:C:422:LEU:HD21	1.91	0.53
1:D:225:VAL:HB	5:D:448:HOH:O	2.08	0.53
1:C:182:LEU:HB2	1:C:395:ARG:NH1	2.24	0.53
1:A:1:MET:HE2	1:A:132:LEU:HD13	1.88	0.53
1:A:57:LEU:HG	1:A:65:VAL:HG22	1.91	0.53
1:B:207:ILE:HD13	1:B:372:HIS:CG	2.44	0.53
1:C:247:ILE:O	1:C:288:LEU:HA	2.08	0.53
1:A:24:ARG:HA	1:A:53:ASP:OD2	2.07	0.53
1:A:63:ASP:CG	1:A:64:HIS:CD2	2.87	0.53
1:A:297:SER:OG	1:A:320:ARG:HG2	2.08	0.53
1:B:208:LEU:CD2	1:B:218:VAL:HG11	2.39	0.53
1:C:27:LEU:HD22	1:C:29:ASP:O	2.09	0.53
1:B:347:LEU:CD1	1:B:358:VAL:HG11	2.39	0.53
1:B:155:ARG:HD3	1:B:431:VAL:O	2.09	0.52
1:C:84:ARG:HD3	2:C:435:SO4:O2	2.09	0.52
1:D:90:MET:HE1	1:D:118:LEU:CD2	2.38	0.52
1:A:142:LEU:CD2	1:A:226:GLU:HB2	2.40	0.52
1:C:382:GLY:O	1:C:386:LEU:HD13	2.09	0.52
1:B:330:SER:O	1:B:368:ARG:HD2	2.10	0.52
1:A:178:PRO:HB3	1:D:126:TRP:CE3	2.44	0.52
1:C:210:LYS:O	1:C:214:GLN:HG2	2.09	0.52
1:D:381:ALA:HB3	1:D:386:LEU:CD1	2.40	0.52
1:D:405:LYS:O	1:D:408:ALA:HB3	2.10	0.52
1:B:80:VAL:HB	1:B:118:LEU:HD23	1.91	0.52
1:C:208:LEU:CD2	1:C:218:VAL:HG11	2.38	0.52
1:B:238:HIS:ND1	1:B:241:ARG:NH1	2.51	0.52
1:C:28:LEU:O	1:C:29:ASP:HB2	2.10	0.52
1:C:138:GLN:NE2	1:C:172:PRO:HG2	2.25	0.52
1:C:241:ARG:HG2	1:C:241:ARG:HH11	1.74	0.52
1:D:266:ARG:HD2	1:D:267:TYR:CE2	2.45	0.52
1:A:87:VAL:HA	1:A:90:MET:HE3	1.92	0.52
1:B:217:LYS:HG2	1:B:307:MET:HG2	1.91	0.52
1:B:420:VAL:HG22	1:B:421:SER:N	2.25	0.52
1:C:33:PHE:CD2	1:C:40:ARG:HB2	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:182:LEU:HD11	1:C:397:VAL:HG23	1.92	0.52
1:D:205:LEU:HD13	1:D:241:ARG:HH21	1.75	0.52
1:A:355:PRO:HB2	1:A:356:PRO:HD2	1.91	0.52
1:B:179:LEU:HB2	1:C:128:ARG:NH1	2.25	0.52
1:A:359:ARG:HD3	1:A:364:GLU:OE1	2.10	0.51
1:A:398:LEU:HB3	1:A:406:LEU:HG	1.92	0.51
1:C:70:LYS:O	1:C:73:ARG:HB3	2.10	0.51
1:D:221:PRO:HA	1:D:311:ALA:O	2.09	0.51
1:B:61:HIS:CD2	1:B:142:LEU:HD11	2.45	0.51
1:C:203:GLU:O	1:C:207:ILE:HG13	2.10	0.51
1:D:221:PRO:HB3	1:D:321:ILE:CG1	2.35	0.51
1:D:32:MET:HG2	1:D:66:GLY:HA3	1.91	0.51
1:C:59:HIS:HE1	1:C:162:ASP:HB2	1.75	0.51
1:D:227:ARG:HG2	1:D:227:ARG:NH2	2.25	0.51
1:D:294:THR:O	1:D:297:SER:HB3	2.10	0.51
1:A:63:ASP:CG	1:A:64:HIS:HD2	2.18	0.51
1:A:85:ALA:HB2	1:A:267:TYR:CE2	2.46	0.51
1:A:259:SER:O	1:A:262:PRO:HD2	2.11	0.51
1:C:88:LEU:HB3	1:C:260:LEU:HD21	1.93	0.51
1:C:195:ARG:HB2	5:C:453:HOH:O	2.09	0.51
1:A:86:THR:HG22	1:A:90:MET:HE2	1.92	0.51
1:B:33:PHE:HB3	1:B:37:GLU:HB2	1.92	0.51
1:D:182:LEU:HD12	1:D:183:VAL:H	1.75	0.51
1:D:269:SER:HB2	2:D:435:SO4:O1	2.10	0.51
1:D:298:LYS:HE2	5:D:451:HOH:O	2.10	0.51
1:A:43:ALA:HB1	1:A:44:PRO:HD2	1.93	0.51
1:C:27:LEU:HB3	1:C:55:VAL:HB	1.93	0.51
1:D:12:VAL:HG23	1:D:13:THR:HG23	1.92	0.51
1:C:217:LYS:HE2	1:C:303:ALA:O	2.11	0.51
1:D:35:GLY:O	1:D:36:LYS:C	2.54	0.51
1:D:57:LEU:HD21	1:D:68:LEU:CD2	2.39	0.51
1:D:90:MET:O	1:D:91:GLU:C	2.53	0.51
1:D:354:ARG:HA	1:D:367:LEU:HD21	1.92	0.51
1:C:170:VAL:HA	1:C:272:VAL:HG21	1.93	0.51
1:D:159:TYR:CD2	1:D:160:SER:N	2.79	0.51
1:D:404:GLU:H	1:D:404:GLU:CD	2.18	0.51
1:A:229:GLN:H	1:A:229:GLN:NE2	2.09	0.50
1:C:213:SER:C	1:C:215:GLY:H	2.19	0.50
1:C:315:MET:HA	1:C:342:GLN:HE22	1.76	0.50
1:D:14:GLY:O	1:D:15:SER:C	2.54	0.50
1:B:274:ALA:O	1:B:278:GLN:HG2	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:321:ILE:HG23	1:C:322:LEU:N	2.26	0.50
1:C:387:LEU:HB3	1:C:416:ARG:NH1	2.26	0.50
1:D:11:GLU:C	1:D:401:GLY:H	2.20	0.50
1:D:232:LEU:HD13	1:D:288:LEU:HD21	1.92	0.50
1:D:336:LEU:C	1:D:336:LEU:HD23	2.36	0.50
1:B:33:PHE:HB2	1:B:41:ASN:HD21	1.77	0.50
1:B:57:LEU:HD22	1:B:80:VAL:HG12	1.93	0.50
1:D:363:GLU:O	1:D:363:GLU:HG3	2.10	0.50
1:C:59:HIS:CD2	1:C:61:HIS:HB2	2.47	0.50
1:C:402:GLU:O	1:C:403:GLU:C	2.54	0.50
1:A:288:LEU:HD12	1:A:289:GLU:N	2.27	0.50
1:C:344:GLN:HG3	1:C:349:ALA:CB	2.41	0.50
1:B:7:GLY:HA3	1:B:14:GLY:O	2.11	0.50
1:D:1:MET:HA	1:D:21:ALA:HB2	1.94	0.50
1:C:80:VAL:HB	1:C:118:LEU:HD23	1.94	0.50
1:C:165:ASN:C	1:C:165:ASN:HD22	2.20	0.50
1:B:9:ALA:O	1:B:10:ARG:HB2	2.10	0.50
1:B:102:ASP:CG	1:B:103:GLU:H	2.20	0.50
1:B:192:ARG:H	1:B:192:ARG:HD2	1.77	0.50
1:D:399:VAL:HG12	1:D:423:ALA:HB3	1.94	0.50
1:A:98:LEU:HG	1:A:106:PHE:HE2	1.76	0.49
1:A:102:ASP:CG	1:A:103:GLU:N	2.68	0.49
1:A:140:GLY:O	1:A:164:GLY:HA3	2.10	0.49
1:C:98:LEU:HD11	1:C:108:PRO:HB3	1.95	0.49
1:C:246:PRO:HB2	1:C:248:TYR:CE1	2.47	0.49
1:D:187:GLY:HA2	1:D:386:LEU:CD2	2.42	0.49
1:D:211:THR:HG23	1:D:212:LEU:N	2.14	0.49
1:D:222:THR:HG22	1:D:339:VAL:CG2	2.42	0.49
1:D:259:SER:O	1:D:262:PRO:HD2	2.12	0.49
1:A:55:VAL:CG1	1:A:80:VAL:HG22	2.42	0.49
1:A:6:PHE:HA	1:A:425:PHE:CE1	2.47	0.49
1:C:229:GLN:HG3	1:C:261:TYR:CZ	2.46	0.49
1:D:229:GLN:H	1:D:229:GLN:CD	2.18	0.49
1:D:363:GLU:O	1:D:364:GLU:C	2.54	0.49
1:A:212:LEU:HB2	1:A:243:PRO:CG	2.42	0.49
1:B:10:ARG:HH12	1:B:424:ARG:HH21	1.59	0.49
1:B:387:LEU:HB3	1:B:416:ARG:NH1	2.27	0.49
1:C:233:TYR:CE1	1:C:282:PRO:HB2	2.48	0.49
1:D:17:HIS:HD2	1:D:184:LEU:CD2	2.23	0.49
1:D:203:GLU:O	1:D:207:ILE:HG13	2.13	0.49
1:D:227:ARG:HH21	1:D:227:ARG:CG	2.24	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1:MET:HG3	1:A:21:ALA:CB	2.41	0.49
1:A:165:ASN:C	1:A:165:ASN:HD22	2.18	0.49
1:C:32:MET:HE2	1:C:105:PHE:HZ	1.78	0.49
1:C:170:VAL:HB	1:C:171:LEU:HD22	1.95	0.49
1:D:200:THR:HG22	1:D:377:PHE:CE1	2.48	0.49
1:D:233:TYR:CE1	1:D:282:PRO:HB2	2.48	0.49
1:C:198:ARG:HG2	1:C:198:ARG:HH21	1.77	0.49
1:D:189:TYR:OH	1:D:341:TYR:HB2	2.12	0.49
1:D:321:ILE:HG23	1:D:322:LEU:N	2.27	0.49
1:A:170:VAL:CG2	1:A:230:GLU:HG3	2.43	0.49
1:C:209:GLU:OE2	1:C:243:PRO:HD3	2.13	0.49
1:C:221:PRO:HD2	1:C:337:VAL:O	2.13	0.49
1:D:253:MET:C	1:D:255:GLY:N	2.71	0.49
1:D:263:ARG:C	1:D:264:LEU:HD12	2.38	0.49
1:D:1:MET:HG3	1:D:21:ALA:HB2	1.93	0.49
1:D:220:ILE:HB	1:D:310:LEU:HG	1.93	0.49
1:D:396:VAL:CG1	1:D:397:VAL:N	2.76	0.49
1:A:63:ASP:OD1	1:A:64:HIS:HD2	1.96	0.49
1:B:142:LEU:O	1:B:143:PRO:C	2.55	0.49
1:D:182:LEU:HD12	1:D:183:VAL:N	2.28	0.48
1:B:198:ARG:HG2	1:B:198:ARG:HH21	1.78	0.48
1:B:312:GLY:HA2	1:B:313:SER:C	2.38	0.48
1:C:325:LEU:O	1:C:329:LEU:HB2	2.13	0.48
1:D:349:ALA:HA	1:D:352:ILE:HD12	1.95	0.48
1:B:178:PRO:HD3	1:C:126:TRP:CD1	2.48	0.48
1:B:196:PRO:O	1:B:199:GLU:HG2	2.14	0.48
1:C:274:ALA:O	1:C:278:GLN:HG2	2.13	0.48
1:D:195:ARG:HH21	1:D:375:GLY:N	2.08	0.48
1:B:86:THR:HG22	1:B:90:MET:CE	2.43	0.48
1:C:163:LEU:HD12	1:C:163:LEU:H	1.78	0.48
1:C:195:ARG:HG3	1:C:195:ARG:HH11	1.78	0.48
1:C:398:LEU:N	1:C:398:LEU:HD22	2.29	0.48
1:B:354:ARG:HG3	1:B:354:ARG:HH11	1.78	0.48
1:D:160:SER:HB2	1:D:163:LEU:CD2	2.42	0.48
1:D:348:GLY:O	1:D:352:ILE:HG13	2.13	0.48
1:B:163:LEU:HD11	1:B:389:TRP:CE2	2.49	0.48
1:D:106:PHE:CD2	1:D:111:VAL:HG22	2.49	0.48
1:D:128:ARG:HD2	5:D:442:HOH:O	2.13	0.48
1:A:177:PRO:HB3	1:A:389:TRP:CZ2	2.49	0.48
1:A:215:GLY:HA2	1:A:306:PRO:HB3	1.96	0.48
1:A:347:LEU:HD11	1:A:358:VAL:HG11	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:396:VAL:O	1:A:420:VAL:HA	2.13	0.48
1:C:298:LYS:CA	1:C:301:ASN:HD22	2.13	0.48
1:D:87:VAL:HA	1:D:90:MET:HE3	1.94	0.48
1:A:350:GLU:O	1:A:353:ALA:HB3	2.14	0.48
1:B:344:GLN:OE1	1:B:344:GLN:HA	2.14	0.48
1:C:207:ILE:HD13	1:C:372:HIS:CD2	2.49	0.48
1:A:98:LEU:HD11	1:A:108:PRO:CA	2.42	0.48
1:B:269:SER:O	1:B:273:GLN:HG3	2.14	0.48
1:A:62:LEU:HD13	1:A:93:VAL:CG1	2.44	0.48
1:B:228:ALA:HB3	1:B:229:GLN:NE2	2.29	0.48
1:C:10:ARG:HH12	1:C:424:ARG:NE	2.12	0.48
1:D:235:LEU:HD13	1:D:247:ILE:HD13	1.95	0.48
1:A:45:PHE:CE2	1:A:70:LYS:HG2	2.49	0.47
1:A:212:LEU:HB2	1:A:243:PRO:HG2	1.96	0.47
1:B:57:LEU:HD22	1:B:80:VAL:CG1	2.44	0.47
1:A:18:LEU:HG	1:A:20:LEU:HD21	1.96	0.47
1:A:22:GLY:C	1:A:24:ARG:H	2.21	0.47
1:C:224:ALA:CB	1:C:253:MET:HE2	2.37	0.47
1:C:387:LEU:HD11	1:C:412:LEU:HD13	1.96	0.47
1:D:208:LEU:HD23	1:D:218:VAL:HG21	1.96	0.47
1:A:163:LEU:N	1:A:163:LEU:HD22	2.28	0.47
1:B:32:MET:HE1	1:B:101:MET:HE1	1.96	0.47
1:B:62:LEU:HD13	1:B:93:VAL:HG12	1.96	0.47
1:C:258:LEU:HD11	1:C:283:PHE:HB3	1.96	0.47
1:D:86:THR:O	1:D:90:MET:HG3	2.15	0.47
1:A:13:THR:HG21	1:A:34:GLN:HB2	1.96	0.47
1:A:90:MET:HE1	1:A:118:LEU:HD22	1.96	0.47
1:D:73:ARG:HG3	1:D:110:ASP:OD2	2.15	0.47
1:A:226:GLU:C	1:A:229:GLN:HE21	2.22	0.47
1:C:208:LEU:HD21	1:C:218:VAL:HG11	1.97	0.47
1:A:277:LEU:C	1:A:279:GLY:H	2.23	0.47
1:C:166:ARG:HG2	1:C:385:GLU:OE1	2.14	0.47
1:B:132:LEU:HD21	1:B:134:LEU:HD21	1.96	0.47
1:C:195:ARG:CB	5:C:453:HOH:O	2.63	0.47
1:C:219:LEU:O	1:C:336:LEU:HD23	2.13	0.47
1:D:280:LYS:HD2	5:D:453:HOH:O	2.15	0.47
1:D:359:ARG:HH12	1:D:362:GLY:HA2	1.77	0.47
1:D:389:TRP:HE3	1:D:390:LEU:CD1	2.28	0.47
1:C:250:ASP:OD1	1:C:320:ARG:NH1	2.43	0.47
1:C:391:GLN:HA	1:C:416:ARG:NH2	2.29	0.47
1:D:316:LEU:HD11	1:D:338:PHE:HE1	1.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:384:ASP:OD2	1:B:384:ASP:N	2.47	0.47
1:D:98:LEU:O	1:D:101:MET:HB3	2.15	0.47
1:D:6:PHE:CZ	1:D:47:PHE:HB3	2.50	0.46
1:A:32:MET:HE1	1:A:101:MET:CE	2.35	0.46
1:A:33:PHE:H	1:A:41:ASN:ND2	2.10	0.46
1:C:389:TRP:HE3	1:C:390:LEU:CD1	2.28	0.46
1:A:320:ARG:CD	3:A:456:BTB:H81	2.44	0.46
1:C:406:LEU:HB3	1:C:422:LEU:CD2	2.45	0.46
1:D:73:ARG:CG	1:D:73:ARG:HH21	2.28	0.46
1:D:87:VAL:HG13	1:D:118:LEU:HD13	1.97	0.46
1:B:2:ARG:NH2	2:B:446:SO4:S	2.89	0.46
1:C:165:ASN:C	1:C:167:GLU:H	2.23	0.46
1:C:224:ALA:HB1	1:C:254:ALA:N	2.30	0.46
1:C:347:LEU:O	1:C:351:ILE:HG13	2.15	0.46
1:A:285:PRO:HD2	1:A:288:LEU:HD22	1.98	0.46
1:B:105:PHE:CD1	1:B:106:PHE:HD1	2.33	0.46
1:B:166:ARG:HG2	1:B:166:ARG:HH11	1.80	0.46
1:B:389:TRP:HE3	1:B:390:LEU:HD13	1.81	0.46
1:C:1:MET:HG2	1:C:431:VAL:CG2	2.44	0.46
1:C:295:GLU:N	1:C:295:GLU:OE2	2.48	0.46
1:C:353:ALA:HB3	1:C:355:PRO:HD3	1.98	0.46
1:C:10:ARG:HB3	1:C:403:GLU:HG3	1.96	0.46
1:C:219:LEU:N	1:C:219:LEU:CD1	2.78	0.46
1:D:409:LEU:CD2	1:D:413:LEU:HG	2.46	0.46
1:A:214:GLN:NE2	1:A:333:ARG:HA	2.31	0.46
1:B:140:GLY:O	1:B:164:GLY:HA3	2.16	0.46
1:B:277:LEU:C	1:B:279:GLY:N	2.74	0.46
1:C:27:LEU:O	1:C:55:VAL:HA	2.16	0.46
1:C:90:MET:HE1	1:C:118:LEU:HD21	1.96	0.46
1:C:266:ARG:HA	1:C:273:GLN:OE1	2.16	0.46
1:D:266:ARG:HD2	1:D:267:TYR:CZ	2.51	0.46
1:A:424:ARG:HD3	1:A:427:GLU:CD	2.41	0.46
1:B:33:PHE:CG	1:B:40:ARG:HB2	2.51	0.46
1:B:165:ASN:C	1:B:165:ASN:ND2	2.72	0.46
1:D:13:THR:CB	1:D:34:GLN:H	2.28	0.46
1:B:157:LEU:HD12	1:B:182:LEU:O	2.16	0.46
1:B:270:GLU:HB2	2:B:438:SO4:O4	2.15	0.46
1:C:98:LEU:HD11	1:C:108:PRO:CA	2.46	0.46
1:C:121:LEU:HD13	1:C:136:PHE:CE2	2.51	0.46
1:D:187:GLY:O	1:D:188:THR:C	2.59	0.46
1:C:96:ASP:C	1:C:98:LEU:N	2.74	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:63:ASP:CG	1:B:64:HIS:HD2	2.24	0.45
1:B:421:SER:C	1:B:422:LEU:HD12	2.41	0.45
1:B:33:PHE:HD2	1:B:41:ASN:HD22	1.64	0.45
1:D:45:PHE:CE1	1:D:67:ARG:HD2	2.51	0.45
1:D:147:PHE:HB2	1:D:159:TYR:O	2.16	0.45
1:A:386:LEU:HD12	1:A:386:LEU:HA	1.82	0.45
1:A:402:GLU:O	1:A:403:GLU:C	2.60	0.45
1:C:49:PRO:HG3	1:C:74:GLU:HB2	1.98	0.45
1:C:359:ARG:HG2	1:C:359:ARG:NH1	2.30	0.45
1:D:157:LEU:HD12	1:D:158:VAL:H	1.82	0.45
1:A:92:ILE:HG22	1:A:253:MET:CE	2.46	0.45
1:A:148:VAL:HB	1:A:159:TYR:HB3	1.97	0.45
1:C:4:VAL:HG22	1:C:428:GLY:CA	2.46	0.45
1:C:61:HIS:NE2	1:C:142:LEU:HD11	2.32	0.45
1:D:62:LEU:HD13	1:D:93:VAL:CG1	2.44	0.45
1:A:212:LEU:CB	1:A:243:PRO:HG2	2.47	0.45
1:B:27:LEU:O	1:B:55:VAL:HA	2.16	0.45
1:D:306:PRO:O	1:D:307:MET:HB3	2.16	0.45
1:D:360:ILE:O	1:D:361:LEU:HB2	2.16	0.45
1:A:393:GLU:HA	1:A:393:GLU:OE2	2.17	0.45
1:B:265:VAL:HA	1:B:268:PHE:CD2	2.52	0.45
1:C:166:ARG:HH11	1:C:166:ARG:HG3	1.81	0.45
1:D:49:PRO:HB3	1:D:71:LEU:CD1	2.46	0.45
1:D:62:LEU:O	1:D:66:GLY:N	2.47	0.45
1:D:351:ILE:CG2	1:D:371:VAL:HG21	2.46	0.45
1:A:320:ARG:HD3	3:A:456:BTB:C8	2.46	0.45
1:B:41:ASN:HD22	1:B:41:ASN:N	2.15	0.45
1:B:398:LEU:HD22	1:B:420:VAL:CG2	2.46	0.45
1:C:57:LEU:HD23	1:C:90:MET:SD	2.57	0.45
1:D:35:GLY:O	1:D:38:GLU:HB2	2.16	0.45
1:D:205:LEU:CD1	1:D:241:ARG:HH21	2.29	0.45
1:D:326:LYS:C	1:D:328:GLY:H	2.24	0.45
1:A:246:PRO:HG2	1:A:248:TYR:HE1	1.80	0.45
1:C:158:VAL:HG23	1:C:180:ALA:HB2	1.98	0.45
1:C:358:VAL:HG12	1:C:359:ARG:N	2.32	0.45
1:C:405:LYS:O	1:C:408:ALA:HB3	2.17	0.45
1:D:241:ARG:O	1:D:242:LEU:HD23	2.17	0.45
1:D:331:ASP:HB3	1:D:334:ASN:HD22	1.78	0.45
1:D:347:LEU:O	1:D:351:ILE:HG13	2.17	0.45
1:A:299:ALA:HA	1:A:302:ARG:HH11	1.82	0.45
1:B:37:GLU:O	1:B:38:GLU:C	2.60	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:200:THR:HG21	1:C:376:GLY:HA3	1.99	0.45
1:D:201:VAL:HG13	1:D:234:VAL:HG11	1.99	0.45
1:A:49:PRO:HB3	1:A:71:LEU:HD12	1.99	0.44
1:A:229:GLN:CD	1:A:229:GLN:N	2.74	0.44
1:A:420:VAL:CG2	1:A:421:SER:N	2.80	0.44
1:B:382:GLY:O	1:B:386:LEU:HB2	2.16	0.44
1:C:241:ARG:HG2	1:C:241:ARG:NH1	2.32	0.44
1:D:15:SER:HA	1:D:17:HIS:HE1	1.82	0.44
1:B:365:VAL:HA	1:B:366:PRO:HD3	1.86	0.44
1:D:12:VAL:HG12	1:D:401:GLY:HA2	1.98	0.44
1:A:187:GLY:O	1:A:188:THR:C	2.59	0.44
3:A:456:BTB:O3	3:A:456:BTB:H71	2.16	0.44
1:C:129:LEU:O	1:C:130:GLY:C	2.59	0.44
1:C:208:LEU:HA	1:C:211:THR:CG2	2.48	0.44
1:D:84:ARG:HB2	1:D:267:TYR:OH	2.18	0.44
1:D:85:ALA:HB2	1:D:267:TYR:CD2	2.53	0.44
1:D:200:THR:HG22	1:D:377:PHE:HE1	1.81	0.44
1:D:420:VAL:HG22	1:D:421:SER:N	2.32	0.44
1:A:52:VAL:HB	1:A:76:TYR:CD2	2.52	0.44
1:B:235:LEU:HD13	1:B:247:ILE:HD13	1.99	0.44
1:B:312:GLY:CA	1:B:313:SER:C	2.91	0.44
1:C:233:TYR:CD1	1:C:282:PRO:HB2	2.53	0.44
1:C:404:GLU:H	1:C:404:GLU:CD	2.24	0.44
1:D:1:MET:HE1	1:D:152:GLY:N	2.32	0.44
1:B:63:ASP:OD1	1:B:64:HIS:HD2	1.99	0.44
1:B:187:GLY:O	1:B:188:THR:C	2.60	0.44
1:C:420:VAL:CG2	1:C:421:SER:N	2.80	0.44
1:B:420:VAL:CG2	1:B:421:SER:N	2.81	0.44
1:C:94:LEU:C	1:C:96:ASP:N	2.75	0.44
1:D:298:LYS:HA	1:D:301:ASN:HD22	1.81	0.44
1:C:9:ALA:O	1:C:10:ARG:HB2	2.17	0.44
1:C:32:MET:HB3	1:C:66:GLY:HA3	2.00	0.44
1:C:90:MET:HE3	1:C:118:LEU:CD1	2.48	0.44
1:A:5:PRO:HG2	1:A:423:ALA:HB1	1.99	0.44
1:B:208:LEU:CD2	1:B:218:VAL:HG21	2.37	0.44
1:B:229:GLN:CD	1:B:229:GLN:N	2.76	0.44
1:B:387:LEU:HB3	1:B:416:ARG:HH12	1.83	0.44
1:C:40:ARG:C	1:C:42:HIS:H	2.25	0.44
1:A:312:GLY:O	1:A:321:ILE:HG22	2.18	0.44
1:B:25:ARG:HH11	1:B:51:GLU:HB3	1.83	0.44
1:B:373:THR:O	1:B:374:LEU:HD23	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:9:ALA:C	1:C:11:GLU:H	2.26	0.44
1:A:178:PRO:HB3	1:D:126:TRP:CD2	2.53	0.43
1:B:12:VAL:HG12	1:B:401:GLY:HA2	1.99	0.43
1:B:94:LEU:HD12	1:B:94:LEU:HA	1.83	0.43
1:B:155:ARG:NH1	1:B:431:VAL:OXT	2.51	0.43
1:D:32:MET:HA	1:D:67:ARG:HG3	1.99	0.43
1:D:62:LEU:HD21	1:D:97:ALA:HB2	2.00	0.43
1:A:28:LEU:O	1:A:29:ASP:HB2	2.18	0.43
1:A:248:TYR:CE2	1:A:300:LEU:HD21	2.53	0.43
1:B:288:LEU:HD12	1:B:289:GLU:H	1.83	0.43
1:C:401:GLY:HA3	1:C:406:LEU:CD1	2.47	0.43
1:A:75:GLY:O	1:A:76:TYR:C	2.62	0.43
1:A:162:ASP:HA	1:A:186:GLU:OE1	2.19	0.43
1:A:179:LEU:HD12	1:A:393:GLU:OE2	2.18	0.43
1:A:315:MET:SD	1:A:343:PRO:HD3	2.58	0.43
1:A:370:SER:OG	1:A:372:HIS:HE1	2.01	0.43
1:B:229:GLN:HG3	1:B:261:TYR:CZ	2.53	0.43
1:B:402:GLU:O	1:B:403:GLU:C	2.61	0.43
1:D:177:PRO:HD3	1:D:389:TRP:NE1	2.33	0.43
1:D:316:LEU:HD11	1:D:338:PHE:CE1	2.53	0.43
1:A:198:ARG:HH21	1:A:198:ARG:CG	2.31	0.43
1:C:167:GLU:O	1:C:168:LYS:O	2.36	0.43
1:A:129:LEU:O	1:A:132:LEU:HB3	2.19	0.43
1:C:59:HIS:NE2	1:C:61:HIS:HB2	2.33	0.43
1:C:284:ARG:HA	1:C:288:LEU:HD23	1.99	0.43
1:D:251:SER:HB3	1:D:254:ALA:HB3	2.01	0.43
1:A:41:ASN:O	1:A:70:LYS:HE3	2.19	0.43
1:D:231:ILE:C	1:D:233:TYR:H	2.26	0.43
1:D:232:LEU:O	1:D:285:PRO:HD3	2.19	0.43
1:A:68:LEU:N	1:A:69:PRO:CD	2.82	0.43
1:A:368:ARG:O	1:A:369:ALA:C	2.61	0.43
1:B:358:VAL:O	1:B:365:VAL:HG13	2.18	0.43
1:C:4:VAL:HA	1:C:5:PRO:HD3	1.90	0.43
1:A:14:GLY:O	1:A:15:SER:C	2.62	0.43
1:B:159:TYR:C	1:B:159:TYR:CD2	2.95	0.43
1:B:229:GLN:H	1:B:229:GLN:NE2	2.16	0.43
1:C:182:LEU:HD11	1:C:397:VAL:CG2	2.49	0.43
1:C:321:ILE:O	1:C:325:LEU:HD13	2.19	0.43
1:D:10:ARG:NH2	1:D:424:ARG:NH2	2.66	0.43
1:D:148:VAL:HG12	1:D:149:VAL:N	2.33	0.43
1:D:187:GLY:O	1:D:189:TYR:N	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:3:ILE:HD12	1:A:19:LEU:CA	2.44	0.43
1:B:84:ARG:HB2	1:B:267:TYR:OH	2.18	0.43
1:C:356:PRO:HG2	1:C:357:ALA:H	1.84	0.43
1:C:382:GLY:O	1:C:383:GLN:C	2.62	0.43
1:D:183:VAL:O	1:D:184:LEU:C	2.62	0.43
1:A:134:LEU:HD23	1:A:150:ALA:HA	2.01	0.43
1:B:55:VAL:HG13	1:B:80:VAL:HG13	2.01	0.43
1:C:325:LEU:O	1:C:329:LEU:N	2.47	0.43
1:C:354:ARG:N	1:C:355:PRO:CD	2.82	0.43
1:C:383:GLN:O	1:C:387:LEU:HG	2.19	0.43
1:D:10:ARG:HB3	1:D:403:GLU:HG3	2.00	0.43
1:D:171:LEU:H	1:D:171:LEU:CD2	2.30	0.43
1:D:207:ILE:HD13	1:D:372:HIS:CG	2.54	0.43
1:C:275:HIS:ND1	1:C:282:PRO:HB3	2.33	0.42
1:C:401:GLY:HA3	1:C:406:LEU:HD11	2.00	0.42
1:D:85:ALA:HB2	1:D:267:TYR:CE2	2.54	0.42
1:B:7:GLY:C	1:B:9:ALA:H	2.27	0.42
1:D:211:THR:HG21	1:D:218:VAL:CG2	2.49	0.42
1:C:399:VAL:HG12	1:C:423:ALA:HB3	2.00	0.42
1:D:253:MET:O	1:D:256:ARG:N	2.53	0.42
1:D:291:VAL:HG11	1:D:297:SER:HB2	2.01	0.42
1:A:165:ASN:C	1:A:165:ASN:ND2	2.77	0.42
1:B:354:ARG:HG3	1:B:354:ARG:NH1	2.34	0.42
1:C:59:HIS:HB3	1:C:145:SER:HA	2.02	0.42
1:D:219:LEU:HD21	1:D:324:HIS:O	2.20	0.42
1:D:262:PRO:C	1:D:264:LEU:H	2.26	0.42
1:A:25:ARG:HD3	1:A:51:GLU:O	2.20	0.42
1:A:163:LEU:HD11	1:A:389:TRP:CE2	2.55	0.42
1:C:70:LYS:NZ	1:C:74:GLU:OE1	2.50	0.42
1:C:270:GLU:HG2	2:C:433:SO4:O3	2.19	0.42
1:D:424:ARG:O	1:D:427:GLU:HB2	2.19	0.42
1:A:57:LEU:HD23	1:A:90:MET:SD	2.59	0.42
1:C:40:ARG:C	1:C:42:HIS:N	2.77	0.42
1:C:57:LEU:HG	1:C:65:VAL:HG22	2.01	0.42
1:D:6:PHE:HE1	1:D:46:GLY:HA3	1.85	0.42
1:D:12:VAL:CG2	1:D:13:THR:N	2.82	0.42
1:D:403:GLU:O	1:D:407:LEU:HD23	2.20	0.42
1:A:233:TYR:CE1	1:A:282:PRO:HB2	2.54	0.42
1:B:86:THR:O	1:B:90:MET:HB2	2.19	0.42
1:C:1:MET:HG2	1:C:431:VAL:HG21	2.02	0.42
1:C:320:ARG:O	1:C:324:HIS:HD2	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:331:ASP:HA	1:C:332:PRO:HD2	1.90	0.42
1:D:91:GLU:O	1:D:95:GLU:CG	2.66	0.42
1:D:112:GLU:HG3	1:D:113:GLU:N	2.34	0.42
1:C:7:GLY:O	1:C:9:ALA:N	2.53	0.42
1:C:138:GLN:NE2	1:C:172:PRO:CG	2.83	0.42
1:C:214:GLN:HE21	1:C:333:ARG:CA	2.31	0.42
1:C:387:LEU:O	1:C:416:ARG:NH2	2.46	0.42
1:B:7:GLY:O	1:B:9:ALA:N	2.53	0.42
1:B:33:PHE:CB	1:B:41:ASN:HD21	2.33	0.42
1:B:192:ARG:CG	1:B:192:ARG:NH1	2.76	0.42
1:C:137:GLY:HA3	1:C:147:PHE:CZ	2.55	0.42
1:C:163:LEU:N	1:C:163:LEU:CD1	2.83	0.42
1:C:187:GLY:O	1:C:189:TYR:N	2.53	0.42
1:D:56:LEU:HD12	1:D:56:LEU:N	2.35	0.42
1:D:76:TYR:CZ	1:D:78:GLY:HA3	2.55	0.42
1:D:367:LEU:HG	1:D:367:LEU:O	2.19	0.42
1:B:7:GLY:C	1:B:9:ALA:N	2.76	0.42
1:B:178:PRO:HB3	1:C:126:TRP:CD2	2.54	0.42
1:C:90:MET:HE1	1:C:118:LEU:CD2	2.50	0.42
1:C:109:GLU:H	1:C:109:GLU:HG3	1.59	0.42
1:C:422:LEU:N	1:C:422:LEU:CD1	2.82	0.42
1:B:40:ARG:C	1:B:42:HIS:H	2.27	0.41
1:B:68:LEU:N	1:B:69:PRO:CD	2.82	0.41
1:D:41:ASN:O	1:D:70:LYS:HE3	2.20	0.41
1:D:68:LEU:N	1:D:69:PRO:HD2	2.35	0.41
1:D:98:LEU:HD11	1:D:108:PRO:HB3	2.02	0.41
1:A:73:ARG:O	1:A:73:ARG:HG3	2.20	0.41
1:C:38:GLU:CD	1:C:38:GLU:C	2.87	0.41
1:C:284:ARG:HA	1:C:288:LEU:CD2	2.50	0.41
1:C:344:GLN:HA	1:C:349:ALA:HB2	2.02	0.41
1:C:403:GLU:O	1:C:407:LEU:CD2	2.68	0.41
1:D:147:PHE:HB2	1:D:160:SER:HA	2.02	0.41
1:D:166:ARG:CD	1:D:166:ARG:C	2.93	0.41
1:D:228:ALA:HB3	1:D:229:GLN:HE22	1.83	0.41
1:D:359:ARG:CB	1:D:364:GLU:HG3	2.50	0.41
1:D:399:VAL:CG1	1:D:423:ALA:HB3	2.49	0.41
1:A:168:LYS:HE2	1:A:230:GLU:OE1	2.20	0.41
1:A:413:LEU:CD2	1:A:418:GLN:OE1	2.67	0.41
1:C:177:PRO:HD3	1:C:389:TRP:CE2	2.56	0.41
1:C:237:THR:O	1:C:238:HIS:CG	2.73	0.41
1:A:85:ALA:HB3	1:A:144:GLY:HA3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:411:LYS:O	1:A:415:LEU:HG	2.20	0.41
1:B:87:VAL:HG13	1:B:118:LEU:CD1	2.51	0.41
1:C:68:LEU:N	1:C:69:PRO:HD2	2.35	0.41
1:C:250:ASP:HA	1:C:291:VAL:HB	2.03	0.41
1:D:11:GLU:OE1	1:D:37:GLU:HG3	2.20	0.41
1:D:220:ILE:HG22	1:D:222:THR:HG23	2.02	0.41
1:D:336:LEU:HD23	1:D:337:VAL:N	2.35	0.41
1:D:425:PHE:C	1:D:427:GLU:H	2.29	0.41
1:B:233:TYR:CE1	1:B:282:PRO:HB2	2.55	0.41
1:D:177:PRO:HB3	1:D:389:TRP:CZ2	2.55	0.41
1:D:351:ILE:HG22	1:D:371:VAL:HG21	2.01	0.41
1:A:226:GLU:CA	1:A:229:GLN:HE21	2.34	0.41
1:A:310:LEU:N	1:A:310:LEU:CD1	2.84	0.41
1:D:141:HIS:CD2	1:D:162:ASP:HB2	2.56	0.41
1:D:269:SER:O	1:D:273:GLN:HG3	2.21	0.41
1:D:355:PRO:HB2	1:D:356:PRO:HD2	2.03	0.41
1:B:65:VAL:HG11	1:B:90:MET:CG	2.50	0.41
1:B:90:MET:O	1:B:94:LEU:HB2	2.21	0.41
1:D:325:LEU:HB3	1:D:329:LEU:HD23	2.01	0.41
1:C:90:MET:HE3	1:C:118:LEU:HD11	2.03	0.41
1:C:166:ARG:HG2	1:C:385:GLU:CD	2.46	0.41
1:C:359:ARG:C	1:C:360:ILE:HG13	2.46	0.41
1:D:58:THR:HG21	1:D:161:GLY:HA3	2.02	0.41
1:D:253:MET:C	1:D:255:GLY:H	2.28	0.41
1:A:1:MET:HE1	1:A:152:GLY:HA3	2.02	0.41
1:A:128:ARG:NH1	1:D:179:LEU:HD13	2.32	0.41
1:A:235:LEU:HD13	1:A:247:ILE:CD1	2.51	0.41
1:C:85:ALA:O	1:C:86:THR:C	2.64	0.41
1:C:390:LEU:O	1:C:391:GLN:C	2.64	0.41
1:D:36:LYS:HE3	1:D:36:LYS:HB3	1.95	0.41
1:D:384:ASP:OD2	1:D:384:ASP:N	2.54	0.41
1:C:424:ARG:O	1:C:425:PHE:C	2.64	0.41
1:D:281:ASN:C	1:D:283:PHE:H	2.28	0.41
1:A:83:THR:HA	1:A:121:LEU:O	2.21	0.40
1:A:84:ARG:HE	1:A:84:ARG:HB2	1.63	0.40
1:B:33:PHE:HD2	1:B:41:ASN:ND2	2.18	0.40
1:B:265:VAL:HA	1:B:268:PHE:HD2	1.84	0.40
1:B:315:MET:HE2	1:B:343:PRO:CD	2.38	0.40
1:B:325:LEU:HA	1:B:325:LEU:HD12	1.78	0.40
1:C:67:ARG:O	1:C:70:LYS:HB3	2.20	0.40
1:D:106:PHE:CE2	1:D:111:VAL:HG22	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:211:THR:O	1:D:212:LEU:C	2.64	0.40
1:D:366:PRO:HG2	1:D:368:ARG:HH12	1.85	0.40
1:D:403:GLU:C	1:D:407:LEU:HD23	2.47	0.40
1:A:98:LEU:HD23	1:A:98:LEU:HA	1.91	0.40
1:B:98:LEU:HD23	1:B:98:LEU:HA	1.91	0.40
1:C:182:LEU:HB3	1:C:431:VAL:HG13	2.03	0.40
1:A:1:MET:CE	1:A:132:LEU:HD13	2.51	0.40
1:B:142:LEU:HD22	1:B:226:GLU:HB2	2.03	0.40
1:C:8:ALA:HB1	1:C:400:VAL:H	1.86	0.40
1:C:75:GLY:O	1:C:76:TYR:C	2.63	0.40
1:D:409:LEU:O	1:D:410:GLY:C	2.64	0.40
1:B:242:LEU:HD22	1:B:243:PRO:HD2	2.04	0.40
1:B:400:VAL:O	1:B:400:VAL:HG12	2.21	0.40
1:C:82:ALA:O	1:C:120:PRO:HA	2.22	0.40
1:C:227:ARG:O	1:C:230:GLU:HB3	2.21	0.40
1:C:386:LEU:HD12	1:C:386:LEU:N	2.36	0.40
1:D:233:TYR:OH	1:D:271:GLU:HG2	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	429/431 (100%)	394 (92%)	32 (8%)	3 (1%)	18	47
1	B	429/431 (100%)	398 (93%)	26 (6%)	5 (1%)	10	34
1	C	429/431 (100%)	367 (86%)	53 (12%)	9 (2%)	5	20
1	D	429/431 (100%)	354 (82%)	55 (13%)	20 (5%)	2	6
All	All	1716/1724 (100%)	1513 (88%)	166 (10%)	37 (2%)	5	19

All (37) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	38	GLU
1	B	102	ASP
1	C	8	ALA
1	C	168	LYS
1	D	38	GLU
1	C	166	ARG
1	D	6	PHE
1	D	15	SER
1	D	36	LYS
1	D	102	ASP
1	D	184	LEU
1	D	211	THR
1	A	102	ASP
1	A	400	VAL
1	B	60	ALA
1	C	188	THR
1	C	244	ARG
1	D	188	THR
1	D	209	GLU
1	D	224	ALA
1	D	328	GLY
1	D	408	ALA
1	B	188	THR
1	C	38	GLU
1	D	100	VAL
1	D	168	LYS
1	D	364	GLU
1	B	196	PRO
1	C	328	GLY
1	D	282	PRO
1	D	400	VAL
1	A	354	ARG
1	C	400	VAL
1	D	66	GLY
1	D	196	PRO
1	D	285	PRO
1	C	143	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	342/342 (100%)	312 (91%)	30 (9%)	9	29
1	B	342/342 (100%)	312 (91%)	30 (9%)	9	29
1	C	342/342 (100%)	317 (93%)	25 (7%)	13	38
1	D	342/342 (100%)	320 (94%)	22 (6%)	16	44
All	All	1368/1368 (100%)	1261 (92%)	107 (8%)	11	35

All (107) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	ARG
1	A	27	LEU
1	A	28	LEU
1	A	55	VAL
1	A	57	LEU
1	A	62	LEU
1	A	84	ARG
1	A	94	LEU
1	A	98	LEU
1	A	157	LEU
1	A	165	ASN
1	A	171	LEU
1	A	192	ARG
1	A	198	ARG
1	A	209	GLU
1	A	229	GLN
1	A	241	ARG
1	A	253	MET
1	A	264	LEU
1	A	320	ARG
1	A	329	LEU
1	A	344	GLN
1	A	363	GLU
1	A	365	VAL
1	A	386	LEU
1	A	390	LEU
1	A	398	LEU
1	A	406	LEU
1	A	407	LEU
1	A	409	LEU

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Mol	Chain	Res	Type
1	B	15	SER
1	B	27	LEU
1	B	57	LEU
1	B	84	ARG
1	B	94	LEU
1	B	96	ASP
1	B	98	LEU
1	B	171	LEU
1	B	186	GLU
1	B	192	ARG
1	B	196	PRO
1	B	198	ARG
1	B	211	THR
1	B	212	LEU
1	B	229	GLN
1	B	253	MET
1	B	264	LEU
1	B	270	GLU
1	B	277	LEU
1	B	320	ARG
1	B	325	LEU
1	B	329	LEU
1	B	336	LEU
1	B	358	VAL
1	B	365	VAL
1	B	390	LEU
1	B	398	LEU
1	B	406	LEU
1	B	407	LEU
1	B	409	LEU
1	C	27	LEU
1	C	28	LEU
1	C	38	GLU
1	C	55	VAL
1	C	62	LEU
1	C	84	ARG
1	C	98	LEU
1	C	102	ASP
1	C	109	GLU
1	C	121	LEU
1	C	129	LEU
1	C	171	LEU

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Mol	Chain	Res	Type
1	C	186	GLU
1	C	192	ARG
1	C	195	ARG
1	C	219	LEU
1	C	229	GLN
1	C	241	ARG
1	C	265	VAL
1	C	266	ARG
1	C	277	LEU
1	C	329	LEU
1	C	336	LEU
1	C	390	LEU
1	C	407	LEU
1	D	2	ARG
1	D	17	HIS
1	D	27	LEU
1	D	36	LYS
1	D	55	VAL
1	D	57	LEU
1	D	73	ARG
1	D	94	LEU
1	D	96	ASP
1	D	98	LEU
1	D	102	ASP
1	D	128	ARG
1	D	166	ARG
1	D	171	LEU
1	D	186	GLU
1	D	192	ARG
1	D	196	PRO
1	D	229	GLN
1	D	295	GLU
1	D	329	LEU
1	D	364	GLU
1	D	406	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (38) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	GLN
1	A	41	ASN
1	A	42	HIS

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Mol	Chain	Res	Type
1	A	59	HIS
1	A	165	ASN
1	A	214	GLN
1	A	229	GLN
1	A	273	GLN
1	A	293	HIS
1	A	301	ASN
1	A	344	GLN
1	A	372	HIS
1	B	34	GLN
1	B	41	ASN
1	B	165	ASN
1	B	229	GLN
1	B	240	HIS
1	B	293	HIS
1	B	380	HIS
1	B	383	GLN
1	C	34	GLN
1	C	41	ASN
1	C	59	HIS
1	C	165	ASN
1	C	214	GLN
1	C	229	GLN
1	C	301	ASN
1	C	323	HIS
1	C	342	GLN
1	D	34	GLN
1	D	41	ASN
1	D	165	ASN
1	D	194	HIS
1	D	229	GLN
1	D	275	HIS
1	D	323	HIS
1	D	324	HIS
1	D	342	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 70 ligands modelled in this entry, 4 are monoatomic - leaving 66 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	SO4	D	432	-	4,4,4	1.06	0	6,6,6	0.58	0
2	SO4	A	445	-	4,4,4	1.07	0	6,6,6	0.57	0
2	SO4	A	449	-	4,4,4	1.07	0	6,6,6	0.58	0
2	SO4	A	437	-	4,4,4	1.07	0	6,6,6	0.56	0
2	SO4	B	432	-	4,4,4	1.07	0	6,6,6	0.59	0
2	SO4	D	434	-	4,4,4	1.06	0	6,6,6	0.57	0
2	SO4	A	439	-	4,4,4	1.10	0	6,6,6	0.57	0
2	SO4	C	438	-	4,4,4	1.06	0	6,6,6	0.59	0
2	SO4	B	436	-	4,4,4	1.07	0	6,6,6	0.57	0
2	SO4	A	444	-	4,4,4	1.15	0	6,6,6	0.56	0
2	SO4	A	442	-	4,4,4	1.06	0	6,6,6	0.58	0
2	SO4	C	444	-	4,4,4	1.05	0	6,6,6	0.58	0
2	SO4	D	437	-	4,4,4	1.06	0	6,6,6	0.57	0
2	SO4	A	450	-	4,4,4	1.09	0	6,6,6	0.56	0
2	SO4	B	446	-	4,4,4	1.09	0	6,6,6	0.57	0
2	SO4	C	436	-	4,4,4	1.06	0	6,6,6	0.57	0
3	BTB	A	456	-	13,13,13	2.55	5 (38%)	7,16,16	0.95	1 (14%)
2	SO4	A	451	-	4,4,4	1.11	0	6,6,6	0.55	0
2	SO4	A	454	-	4,4,4	1.06	0	6,6,6	0.57	0
2	SO4	A	438	-	4,4,4	1.04	0	6,6,6	0.59	0
2	SO4	B	434	-	4,4,4	1.00	0	6,6,6	0.60	0
2	SO4	A	435	-	4,4,4	1.07	0	6,6,6	0.57	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SO4	C	435	-	4,4,4	1.04	0	6,6,6	0.60	0
2	SO4	A	432	-	4,4,4	1.00	0	6,6,6	0.56	0
2	SO4	D	439	-	4,4,4	1.08	0	6,6,6	0.57	0
2	SO4	A	446	-	4,4,4	1.07	0	6,6,6	0.57	0
2	SO4	A	447	-	4,4,4	1.06	0	6,6,6	0.58	0
2	SO4	B	438	-	4,4,4	1.14	0	6,6,6	0.56	0
2	SO4	B	440	-	4,4,4	1.09	0	6,6,6	0.58	0
2	SO4	C	434	-	4,4,4	1.04	0	6,6,6	0.58	0
2	SO4	A	434	-	4,4,4	1.07	0	6,6,6	0.57	0
2	SO4	C	445	-	4,4,4	1.06	0	6,6,6	0.57	0
2	SO4	B	437	-	4,4,4	1.12	0	6,6,6	0.58	0
2	SO4	C	440	-	4,4,4	1.06	0	6,6,6	0.58	0
2	SO4	C	437	-	4,4,4	1.01	0	6,6,6	0.59	0
2	SO4	A	433	-	4,4,4	1.07	0	6,6,6	0.59	0
2	SO4	A	448	-	4,4,4	0.97	0	6,6,6	0.62	0
2	SO4	B	435	-	4,4,4	1.05	0	6,6,6	0.60	0
2	SO4	D	435	-	4,4,4	1.08	0	6,6,6	0.57	0
2	SO4	B	444	-	4,4,4	1.08	0	6,6,6	0.58	0
2	SO4	A	452	-	4,4,4	1.11	0	6,6,6	0.55	0
2	SO4	B	442	-	4,4,4	1.08	0	6,6,6	0.57	0
2	SO4	D	436	-	4,4,4	1.04	0	6,6,6	0.58	0
2	SO4	C	433	-	4,4,4	1.06	0	6,6,6	0.58	0
2	SO4	B	447	-	4,4,4	1.09	0	6,6,6	0.57	0
2	SO4	C	432	-	4,4,4	1.10	0	6,6,6	0.56	0
2	SO4	C	439	-	4,4,4	1.08	0	6,6,6	0.56	0
2	SO4	B	450	-	4,4,4	1.13	0	6,6,6	0.55	0
2	SO4	D	433	-	4,4,4	1.07	0	6,6,6	0.58	0
2	SO4	B	441	-	4,4,4	1.06	0	6,6,6	0.58	0
2	SO4	A	453	-	4,4,4	1.07	0	6,6,6	0.57	0
2	SO4	C	442	-	4,4,4	1.07	0	6,6,6	0.58	0
2	SO4	B	443	-	4,4,4	1.07	0	6,6,6	0.57	0
2	SO4	B	445	-	4,4,4	1.11	0	6,6,6	0.55	0
2	SO4	A	443	-	4,4,4	1.08	0	6,6,6	0.56	0
2	SO4	B	433	-	4,4,4	1.05	0	6,6,6	0.57	0
2	SO4	D	438	-	4,4,4	1.08	0	6,6,6	0.56	0
2	SO4	A	436	-	4,4,4	1.05	0	6,6,6	0.58	0
2	SO4	B	439	-	4,4,4	1.10	0	6,6,6	0.57	0
2	SO4	B	448	-	4,4,4	1.08	0	6,6,6	0.56	0
2	SO4	C	443	-	4,4,4	1.09	0	6,6,6	0.54	0
2	SO4	B	449	-	4,4,4	1.07	0	6,6,6	0.59	0
2	SO4	A	455	-	4,4,4	1.06	0	6,6,6	0.58	0
2	SO4	A	441	-	4,4,4	1.08	0	6,6,6	0.57	0
2	SO4	C	441	-	4,4,4	1.06	0	6,6,6	0.59	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SO4	A	440	-	4,4,4	1.11	0	6,6,6	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
3	BTB	A	456	-	-	2/21/21/21	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	456	BTB	C2-N	6.49	1.61	1.48
3	A	456	BTB	C5-N	4.08	1.53	1.48
3	A	456	BTB	C1-C2	2.96	1.56	1.53
3	A	456	BTB	C4-C2	2.81	1.56	1.53
3	A	456	BTB	C3-C2	2.52	1.56	1.53

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	456	BTB	C6-C5-N	2.01	119.45	111.59

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	456	BTB	N-C5-C6-O6
3	A	456	BTB	C6-C5-N-C2

There are no ring outliers.

7 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	446	SO4	1	0
3	A	456	BTB	6	0
2	C	435	SO4	1	0
2	B	438	SO4	1	0
2	D	435	SO4	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	433	SO4	1	0
2	B	441	SO4	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	431/431 (100%)	-0.35	0 100 100	20, 38, 59, 82	0
1	B	431/431 (100%)	-0.27	0 100 100	22, 40, 65, 83	0
1	C	431/431 (100%)	0.31	11 (2%) 57 47	17, 68, 102, 115	0
1	D	431/431 (100%)	0.77	34 (7%) 18 13	35, 78, 132, 146	0
All	All	1724/1724 (100%)	0.11	45 (2%) 57 47	17, 50, 113, 146	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	366	PRO	3.8
1	D	223	PHE	3.8
1	D	337	VAL	3.6
1	D	329	LEU	3.5
1	D	361	LEU	3.1
1	D	249	LEU	3.1
1	D	356	PRO	3.1
1	D	338	PHE	3.0
1	D	365	VAL	3.0
1	D	360	ILE	2.9
1	D	347	LEU	2.9
1	C	335	ALA	2.9
1	D	339	VAL	2.8
1	D	341	TYR	2.7
1	D	310	LEU	2.6
1	D	190	GLY	2.6
1	D	218	VAL	2.6
1	D	327	HIS	2.5
1	C	343	PRO	2.5
1	D	330	SER	2.5
1	C	296	ALA	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	341	TYR	2.4
1	D	374	LEU	2.4
1	D	239	GLY	2.4
1	D	311	ALA	2.4
1	C	240	HIS	2.4
1	C	300	LEU	2.4
1	D	308	VAL	2.3
1	D	35	GLY	2.3
1	D	369	ALA	2.3
1	C	290	VAL	2.3
1	D	343	PRO	2.3
1	C	337	VAL	2.2
1	D	400	VAL	2.2
1	D	212	LEU	2.2
1	D	348	GLY	2.2
1	D	309	VAL	2.2
1	D	224	ALA	2.2
1	D	322	LEU	2.1
1	C	277	LEU	2.0
1	C	336	LEU	2.0
1	D	219	LEU	2.0
1	D	336	LEU	2.0
1	C	192	ARG	2.0
1	D	319	GLY	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	C	445	5/5	0.56	0.09	120,121,121,121	0
2	SO4	A	436	5/5	0.60	0.11	135,135,135,136	0
2	SO4	A	451	5/5	0.61	0.15	145,145,146,146	0
2	SO4	C	438	5/5	0.62	0.17	137,137,138,138	0
2	SO4	A	447	5/5	0.65	0.14	129,129,130,130	0
2	SO4	B	442	5/5	0.66	0.15	119,119,119,120	0
2	SO4	B	436	5/5	0.66	0.15	125,125,125,125	0
2	SO4	B	441	5/5	0.66	0.12	138,138,138,138	0
2	SO4	D	438	5/5	0.66	0.10	146,146,147,147	0
2	SO4	C	440	5/5	0.67	0.13	132,132,132,133	0
2	SO4	B	447	5/5	0.67	0.13	130,131,131,131	0
2	SO4	B	435	5/5	0.67	0.13	112,112,112,112	0
2	SO4	B	444	5/5	0.68	0.13	149,150,150,150	0
2	SO4	A	453	5/5	0.68	0.19	153,153,153,153	0
2	SO4	D	437	5/5	0.69	0.10	135,135,135,136	0
2	SO4	D	439	5/5	0.69	0.10	139,139,139,139	0
2	SO4	B	449	5/5	0.70	0.14	112,112,112,113	0
2	SO4	A	452	5/5	0.71	0.12	129,129,129,129	0
2	SO4	D	434	5/5	0.72	0.14	151,151,152,152	0
3	BTB	A	456	14/14	0.73	0.15	60,67,70,70	0
2	SO4	B	440	5/5	0.74	0.12	140,141,141,141	0
2	SO4	A	446	5/5	0.75	0.14	144,144,144,144	0
2	SO4	B	448	5/5	0.76	0.18	131,131,132,132	0
2	SO4	A	435	5/5	0.76	0.14	171,171,172,172	0
2	SO4	A	445	5/5	0.77	0.10	127,127,127,127	0
2	SO4	A	450	5/5	0.77	0.12	123,123,124,124	0
2	SO4	C	443	5/5	0.77	0.11	93,93,94,94	0
2	SO4	C	436	5/5	0.78	0.10	117,117,117,117	0
2	SO4	B	433	5/5	0.78	0.14	114,114,115,115	0
2	SO4	A	438	5/5	0.78	0.16	114,114,115,115	0
2	SO4	D	435	5/5	0.78	0.13	97,98,99,99	0
2	SO4	A	439	5/5	0.79	0.15	134,134,135,135	0
2	SO4	A	433	5/5	0.79	0.11	130,131,131,131	0
2	SO4	A	454	5/5	0.79	0.10	110,110,111,111	0
2	SO4	C	441	5/5	0.80	0.18	133,134,134,134	0
2	SO4	B	443	5/5	0.81	0.12	117,117,118,118	0
2	SO4	A	434	5/5	0.81	0.13	127,127,127,127	0
2	SO4	B	445	5/5	0.81	0.18	104,104,105,105	0
2	SO4	B	437	5/5	0.82	0.15	97,98,98,98	0
2	SO4	A	437	5/5	0.82	0.10	126,126,126,126	0
2	SO4	A	455	5/5	0.83	0.13	136,136,136,136	0
2	SO4	C	433	5/5	0.84	0.10	98,100,100,100	0
2	SO4	A	441	5/5	0.85	0.11	110,111,111,111	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	C	437	5/5	0.86	0.10	126,126,126,126	0
2	SO4	A	443	5/5	0.86	0.10	96,96,97,97	0
2	SO4	A	440	5/5	0.87	0.23	99,100,101,101	0
2	SO4	D	436	5/5	0.87	0.08	119,119,120,120	0
2	SO4	C	444	5/5	0.88	0.18	127,128,128,128	0
2	SO4	B	432	5/5	0.88	0.11	92,93,93,94	0
2	SO4	D	433	5/5	0.88	0.08	99,100,100,101	0
2	SO4	C	442	5/5	0.88	0.09	90,90,91,91	0
2	SO4	B	446	5/5	0.88	0.13	96,96,97,97	0
2	SO4	B	434	5/5	0.89	0.09	75,75,76,76	0
2	SO4	C	432	5/5	0.90	0.09	84,85,86,87	0
2	SO4	A	442	5/5	0.90	0.12	104,104,104,104	0
2	SO4	C	434	5/5	0.90	0.08	93,93,93,93	0
2	SO4	B	438	5/5	0.90	0.11	70,70,71,71	0
2	SO4	C	435	5/5	0.91	0.11	90,90,90,90	0
2	SO4	A	449	5/5	0.91	0.17	84,84,85,85	0
2	SO4	D	432	5/5	0.93	0.09	80,81,81,82	0
2	SO4	B	450	5/5	0.93	0.09	51,52,54,56	0
2	SO4	C	439	5/5	0.93	0.07	63,64,65,65	0
4	ZN	D	440	1/1	0.94	0.16	86,86,86,86	0
4	ZN	B	451	1/1	0.95	0.14	76,76,76,76	0
2	SO4	A	448	5/5	0.95	0.07	70,71,72,73	0
2	SO4	A	444	5/5	0.96	0.06	47,49,50,50	0
2	SO4	B	439	5/5	0.96	0.07	62,63,64,64	0
4	ZN	A	457	1/1	0.97	0.14	50,50,50,50	0
4	ZN	C	446	1/1	0.98	0.10	76,76,76,76	0
2	SO4	A	432	5/5	0.98	0.06	36,37,38,42	0

6.5 Other polymers [i](#)

There are no such residues in this entry.