



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 12:35 PM UTC

PDB ID : 9PR8 / pdb_00009pr8
Title : Crystal Structure of the Clostridiodes difficile CspC-CspA heterodimer.
Authors : Heldwein, E.E.; Gonzalez-Del Pino, G.L.
Deposited on : 2025-07-23
Resolution : 3.35 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

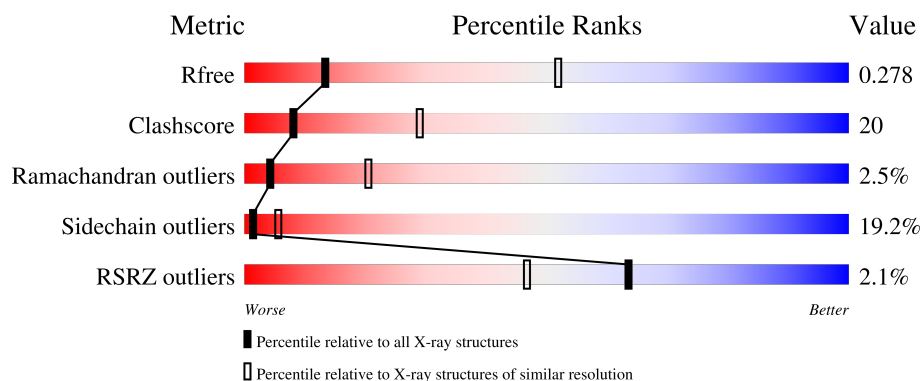
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.35 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	1099 (3.40-3.32)
Clashscore	190562	1116 (3.40-3.32)
Ramachandran outliers	187476	1101 (3.40-3.32)
Sidechain outliers	187428	1101 (3.40-3.32)
RSRZ outliers	180081	1099 (3.40-3.32)

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	551	2% 48% 38% 10% .
1	B	551	4% 49% 41% 7% .
2	C	565	2% 52% 39% 7% ..
2	D	565	% 54% 37% 6% .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 15999 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 Germination-specific protease.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	528	Total	C	N	O	S	0	0	0
			3886	2465	634	777	10			
1	B	530	Total	C	N	O	S	0	0	0
			3838	2446	622	760	10			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP A0A9R0BLF4
A	363	GLU	PHE	conflict	UNP A0A9R0BLF4
A	511	GLU	TYR	conflict	UNP A0A9R0BLF4
B	1	MET	-	initiating methionine	UNP A0A9R0BLF4
B	363	GLU	PHE	conflict	UNP A0A9R0BLF4
B	511	GLU	TYR	conflict	UNP A0A9R0BLF4

- Molecule 2 is a protein called Germination-specific protease.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	556	Total	C	N	O	S	0	0	0
			4171	2652	680	832	7			
2	D	551	Total	C	N	O	S	0	0	0
			4102	2601	664	830	7			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	558	LEU	-	expression tag	UNP A0A9R0BLE1
C	559	GLU	-	expression tag	UNP A0A9R0BLE1
C	560	HIS	-	expression tag	UNP A0A9R0BLE1
C	561	HIS	-	expression tag	UNP A0A9R0BLE1
C	562	HIS	-	expression tag	UNP A0A9R0BLE1
C	563	HIS	-	expression tag	UNP A0A9R0BLE1

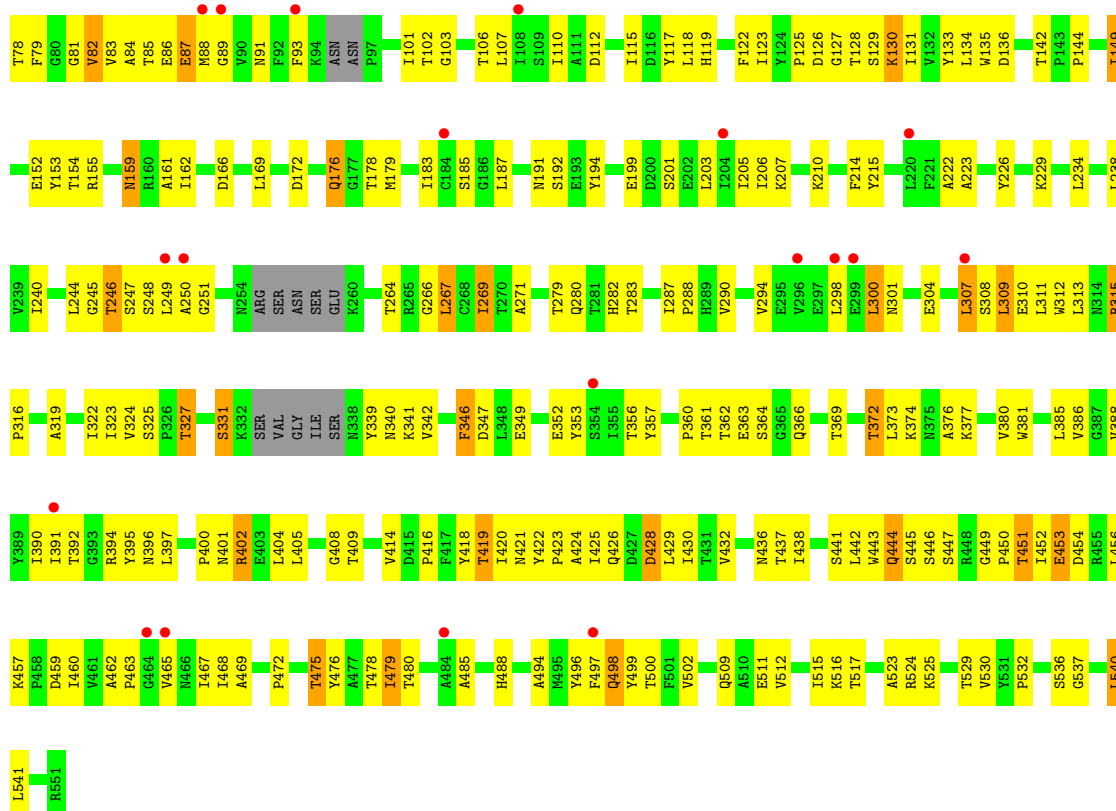
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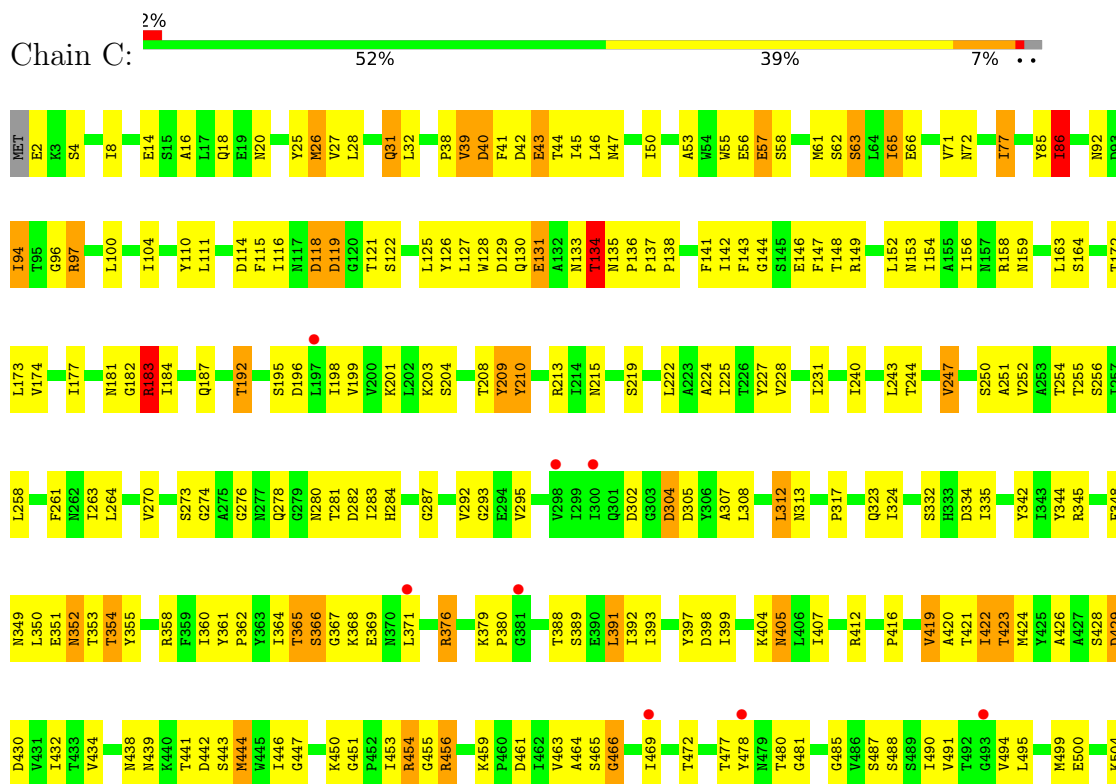
Chain	Residue	Modelled	Actual	Comment	Reference
C	564	HIS	-	expression tag	UNP A0A9R0BLE1
C	565	HIS	-	expression tag	UNP A0A9R0BLE1
D	558	LEU	-	expression tag	UNP A0A9R0BLE1
D	559	GLU	-	expression tag	UNP A0A9R0BLE1
D	560	HIS	-	expression tag	UNP A0A9R0BLE1
D	561	HIS	-	expression tag	UNP A0A9R0BLE1
D	562	HIS	-	expression tag	UNP A0A9R0BLE1
D	563	HIS	-	expression tag	UNP A0A9R0BLE1
D	564	HIS	-	expression tag	UNP A0A9R0BLE1
D	565	HIS	-	expression tag	UNP A0A9R0BLE1

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	2	Total O 2 2	0	0

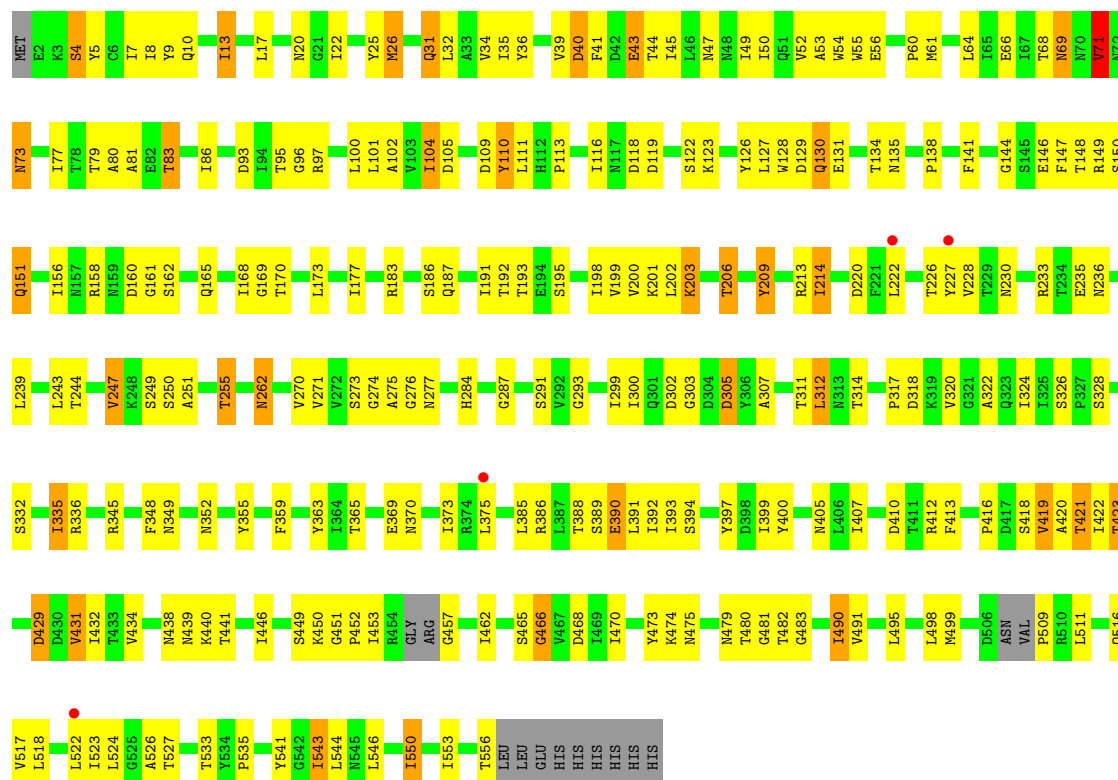


● Molecule 2: Germination-specific protease





• Molecule 2: Germination-specific protease



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	71.14Å 105.50Å 311.32Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	99.92 – 3.35 99.92 – 3.35	Depositor EDS
% Data completeness (in resolution range)	99.7 (99.92-3.35) 99.7 (99.92-3.35)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.24 (at 3.33Å)	Xtriage
Refinement program	PHENIX 1.21.2_5419	Depositor
R, R_{free}	0.246 , 0.278 0.246 , 0.278	Depositor DCC
R_{free} test set	1735 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	86.2	Xtriage
Anisotropy	0.485	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 34.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	15999	wwPDB-VP
Average B, all atoms (Å ²)	81.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.07% 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 [i](#)

5.1 Standard geometry [i](#)

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	0.59	0/3960	0.86	0/5407
1	B	0.50	0/3915	0.78	0/5355
2	C	0.56	0/4247	0.80	0/5810
2	D	0.51	0/4172	0.76	0/5711
All	All	0.54	0/16294	0.80	0/22283

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	3886	0	3672	168	0
1	B	3838	0	3574	156	0
2	C	4171	0	4043	163	0
2	D	4102	0	3939	141	0
3	A	2	0	0	0	0
All	All	15999	0	15228	610	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 20.

All (610) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:246:THR:HG22	1:A:248:SER:H	1.28	0.97
2:C:420:ALA:H	2:C:451:GLY:HA3	1.36	0.90
2:D:420:ALA:H	2:D:451:GLY:HA3	1.36	0.89
1:A:9:ILE:HD11	1:A:62:THR:HB	1.57	0.87
2:D:138:PRO:HG3	2:D:146:GLU:HB2	1.57	0.86
1:A:324:VAL:HB	1:A:382:LYS:HB3	1.59	0.84
1:B:341:LYS:HG3	1:B:356:THR:HG22	1.60	0.83
1:A:265:ARG:HH21	2:C:53:ALA:HA	1.44	0.80
2:C:32:LEU:HD21	2:C:222:LEU:HD13	1.62	0.80
1:A:240:ILE:HB	1:A:269:ILE:HG23	1.64	0.79
2:C:419:VAL:HG12	2:C:451:GLY:HA2	1.64	0.78
1:A:22:ALA:HA	1:A:27:SER:HB2	1.67	0.76
1:B:363:GLU:HG2	1:B:364:SER:H	1.49	0.76
2:C:116:ILE:HG23	2:C:156:ILE:HD13	1.68	0.75
1:A:465:VAL:HA	1:A:480:THR:HG23	1.67	0.75
1:B:82:VAL:HG13	1:B:467:ILE:HD13	1.70	0.74
2:D:32:LEU:HD21	2:D:222:LEU:HD13	1.69	0.73
1:B:452:ILE:H	1:B:452:ILE:HD12	1.51	0.73
2:D:77:ILE:HD13	2:D:177:ILE:HA	1.71	0.72
2:C:213:ARG:NH1	2:C:407:ILE:O	2.23	0.72
2:C:128:TRP:HE1	2:C:144:GLY:HA3	1.54	0.71
2:C:428:SER:C	2:C:430:ASP:H	1.98	0.71
2:C:312:LEU:HD12	2:C:399:ILE:HG12	1.73	0.71
2:C:118:ASP:HA	2:C:183:ARG:HH22	1.56	0.71
2:C:181:ASN:OD1	2:C:182:GLY:N	2.24	0.70
2:C:131:GLU:HG3	2:C:203:LYS:HA	1.74	0.69
2:D:64:LEU:HD22	2:D:276:GLY:HA2	1.75	0.69
2:C:428:SER:O	2:C:430:ASP:N	2.25	0.69
1:B:428:ASP:N	1:B:428:ASP:OD1	2.25	0.69
2:D:129:ASP:OD2	2:D:201:LYS:HD3	1.91	0.69
1:A:84:ALA:HB3	1:A:488:HIS:NE2	2.08	0.68
1:B:279:THR:O	1:B:402:ARG:NH2	2.26	0.68
2:D:420:ALA:H	2:D:451:GLY:CA	2.06	0.68
2:D:213:ARG:NH1	2:D:407:ILE:O	2.27	0.68
1:A:334:VAL:HG11	1:A:340:ASN:HB3	1.75	0.67
2:D:86:ILE:HG12	2:D:546:LEU:HD21	1.77	0.67
1:A:280:GLN:HB3	1:A:414:VAL:HG21	1.76	0.67
1:B:33:ILE:HG23	1:B:149:ILE:HG12	1.77	0.67
1:B:451:THR:HG22	1:B:453:GLU:H	1.58	0.67
1:B:194:TYR:HD2	1:B:476:TYR:CD2	2.12	0.67
1:A:451:THR:HG22	1:A:453:GLU:H	1.58	0.67
1:A:452:ILE:HD12	1:A:452:ILE:H	1.59	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:65:ILE:HD11	2:C:244:THR:HB	1.77	0.67
1:A:132:VAL:HG22	1:A:202:GLU:HG2	1.77	0.66
2:D:17:LEU:HB3	2:D:22:ILE:HB	1.78	0.66
2:D:129:ASP:O	2:D:131:GLU:N	2.27	0.66
1:A:255:ARG:HD3	2:C:430:ASP:HB2	1.78	0.65
1:A:179:MET:HE1	1:A:477:ALA:HB3	1.79	0.65
2:C:295:VAL:HG22	2:C:388:THR:OG1	1.96	0.65
2:C:251:ALA:HB2	2:C:420:ALA:HB1	1.77	0.65
1:A:526:ASP:OD1	1:A:527:SER:N	2.28	0.65
2:D:102:ALA:HB2	2:D:239:LEU:HD11	1.79	0.65
2:D:128:TRP:HE1	2:D:144:GLY:HA3	1.62	0.65
1:A:446:SER:O	1:A:448:ARG:HG3	1.97	0.65
2:C:353:THR:HG23	2:C:376:ARG:O	1.97	0.65
2:D:44:THR:HA	2:D:47:ASN:HD22	1.62	0.64
1:A:325:SER:HB3	1:A:346:PHE:CE2	2.32	0.64
2:D:8:ILE:HG23	2:D:31:GLN:HB3	1.77	0.64
2:D:420:ALA:N	2:D:451:GLY:HA3	2.11	0.64
1:A:206:ILE:HD13	1:A:223:ALA:HB1	1.79	0.64
1:A:122:PHE:O	1:A:130:LYS:HG3	1.98	0.63
2:C:129:ASP:HB2	2:C:164:SER:HB2	1.79	0.63
1:A:153:TYR:CE1	1:A:169:LEU:HD22	2.34	0.63
1:A:402:ARG:C	1:A:404:LEU:H	2.05	0.63
1:A:13:HIS:HB2	1:A:17:PHE:HB2	1.80	0.63
2:C:8:ILE:HD12	2:C:31:GLN:HB2	1.79	0.62
2:C:270:VAL:HG21	2:C:518:LEU:HD12	1.80	0.62
2:C:459:LYS:NZ	2:C:461:ASP:OD2	2.32	0.62
2:D:104:ILE:HG23	2:D:200:VAL:HB	1.81	0.62
2:D:122:SER:HB3	2:D:156:ILE:HD11	1.81	0.62
1:B:54:LEU:O	1:B:59:ILE:HG12	2.00	0.62
1:B:443:TRP:O	1:B:446:SER:N	2.27	0.62
1:B:245:GLY:HA2	1:B:423:PRO:HG3	1.81	0.62
2:D:419:VAL:HG12	2:D:451:GLY:HA2	1.81	0.62
2:D:491:VAL:HG22	2:D:544:LEU:HD22	1.81	0.62
1:A:336:ILE:HG22	1:A:337:SER:H	1.65	0.62
1:B:159:ASN:HA	1:B:162:ILE:HD12	1.79	0.62
2:C:292:VAL:HG23	2:C:392:ILE:HG12	1.80	0.62
2:C:116:ILE:O	2:C:183:ARG:NH1	2.33	0.62
2:D:363:TYR:CE2	2:D:365:THR:HB	2.35	0.62
1:A:341:LYS:HG3	1:A:356:THR:HG23	1.82	0.62
2:C:210:TYR:HB3	2:C:213:ARG:HG3	1.82	0.61
1:B:74:ILE:HD11	1:B:176:GLN:HG3	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:84:ALA:C	1:B:86:GLU:H	2.08	0.61
1:B:443:TRP:O	1:B:445:SER:N	2.33	0.61
1:B:149:ILE:HG13	1:B:149:ILE:O	2.00	0.61
1:A:82:VAL:HG13	1:A:467:ILE:HD13	1.81	0.61
1:A:102:THR:OG1	1:A:200:ASP:HB2	2.00	0.61
1:A:246:THR:HA	1:A:401:ASN:HD21	1.64	0.61
2:C:273:SER:OG	2:C:274:GLY:N	2.34	0.61
1:B:280:GLN:HB3	1:B:414:VAL:HG21	1.83	0.60
1:B:70:LEU:HD11	1:B:215:TYR:HE2	1.66	0.60
2:C:548:ASN:O	2:C:552:GLN:HG2	2.00	0.60
1:A:263:PHE:HB3	2:C:8:ILE:HG21	1.84	0.60
2:D:169:GLY:HA3	2:D:473:TYR:CE1	2.35	0.60
1:A:248:SER:OG	1:A:249:LEU:N	2.34	0.60
2:D:275:ALA:O	2:D:449:SER:OG	2.20	0.60
2:D:73:ASN:N	2:D:73:ASN:OD1	2.34	0.60
2:D:418:SER:O	2:D:421:THR:OG1	2.20	0.60
1:A:53:VAL:O	1:A:56:LEU:HB2	2.02	0.60
1:B:271:ALA:HB3	1:B:429:LEU:HD13	1.83	0.60
1:B:12:VAL:HA	1:B:36:SER:O	2.02	0.60
1:B:191:ASN:HB3	1:B:476:TYR:HE2	1.67	0.59
1:B:436:ASN:OD1	1:B:438:ILE:HB	2.01	0.59
2:D:429:ASP:N	2:D:429:ASP:OD1	2.35	0.59
1:A:496:TYR:CD2	1:A:546:MET:HE2	2.38	0.59
1:B:74:ILE:HG13	1:B:479:ILE:HD12	1.85	0.59
2:D:302:ASP:OD1	2:D:303:GLY:N	2.34	0.59
2:D:61:MET:HG2	2:D:247:VAL:HG13	1.84	0.59
1:A:387:GLY:HA3	1:A:390:ILE:HD11	1.83	0.59
2:C:317:PRO:HG2	2:C:393:ILE:HG13	1.84	0.59
2:C:443:SER:O	2:C:444:MET:HB2	2.02	0.59
1:B:9:ILE:HD11	1:B:62:THR:HB	1.84	0.58
1:A:84:ALA:C	1:A:86:GLU:H	2.11	0.58
2:C:122:SER:N	2:C:153:ASN:OD1	2.36	0.58
1:A:300:LEU:HD12	1:A:302:GLU:O	2.02	0.58
2:D:287:GLY:HA3	2:D:397:TYR:CE1	2.39	0.58
1:B:22:ALA:HA	1:B:27:SER:OG	2.04	0.58
2:D:81:ALA:HB2	2:D:490:ILE:HD13	1.86	0.58
2:C:138:PRO:HB2	2:C:141:PHE:CD2	2.39	0.57
1:A:509:GLN:HA	1:A:514:LYS:HD2	1.86	0.57
2:D:138:PRO:HB2	2:D:141:PHE:HD2	1.69	0.57
1:B:465:VAL:HA	1:B:480:THR:HG23	1.86	0.57
2:D:324:ILE:HB	2:D:348:PHE:HZ	1.70	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:130:LYS:HZ3	1:B:185:SER:HB3	1.68	0.57
1:B:401:ASN:ND2	1:B:421:ASN:HD21	2.02	0.57
2:C:323:GLN:NE2	2:C:334:ASP:OD1	2.38	0.57
2:D:45:ILE:O	2:D:49:ILE:HG13	2.05	0.57
1:B:400:PRO:HB2	1:B:405:LEU:HD21	1.86	0.57
2:C:495:LEU:O	2:C:499:MET:HG3	2.03	0.57
2:C:270:VAL:HG13	2:C:515:THR:HG23	1.85	0.56
2:D:527:THR:N	2:D:543:ILE:O	2.35	0.56
2:D:270:VAL:HG21	2:D:518:LEU:HD12	1.87	0.56
1:A:526:ASP:CG	1:A:527:SER:H	2.13	0.56
1:B:115:ILE:HD13	1:B:131:ILE:HD11	1.87	0.56
1:B:247:SER:OG	1:B:400:PRO:HA	2.06	0.56
1:B:308:SER:HB2	1:B:372:THR:HG23	1.87	0.56
2:C:41:PHE:CE1	2:C:46:LEU:HG	2.40	0.56
1:B:459:ASP:OD2	1:B:516:LYS:HD3	2.05	0.56
2:D:336:ARG:NH1	2:D:390:GLU:OE1	2.39	0.56
1:A:532:PRO:HB3	1:A:538:TYR:CE1	2.41	0.56
2:C:302:ASP:OD1	2:C:305:ASP:HB2	2.06	0.56
1:B:55:GLN:NE2	2:D:509:PRO:HD2	2.21	0.56
2:D:535:PRO:HB3	2:D:541:TYR:CE1	2.41	0.56
2:C:509:PRO:O	2:C:511:LEU:N	2.39	0.56
1:A:514:LYS:HD3	2:C:362:PRO:HD2	1.87	0.56
1:B:144:PRO:HG3	1:B:152:GLU:HB2	1.88	0.56
2:D:335:ILE:O	2:D:335:ILE:HD12	2.06	0.56
2:D:419:VAL:HG12	2:D:452:PRO:HD3	1.87	0.55
1:B:56:LEU:C	1:B:58:SER:H	2.14	0.55
2:D:113:PRO:HB3	2:D:183:ARG:HD2	1.88	0.55
2:C:138:PRO:HB2	2:C:141:PHE:HD2	1.72	0.55
2:C:293:GLY:N	2:C:389:SER:OG	2.38	0.55
2:D:101:LEU:HD11	2:D:192:THR:HG21	1.87	0.55
2:D:230:ASN:HA	2:D:233:ARG:HD3	1.88	0.55
1:A:30:PHE:CZ	1:A:38:GLY:HA3	2.41	0.55
1:B:136:ASP:HA	1:B:207:LYS:CB	2.36	0.55
2:D:17:LEU:HD22	2:D:22:ILE:HD13	1.87	0.55
1:B:250:ALA:HA	1:B:425:ILE:HD13	1.89	0.55
2:D:209:TYR:HB3	2:D:307:ALA:O	2.07	0.55
1:A:535:ASN:OD1	1:A:535:ASN:N	2.39	0.55
1:A:68:MET:HE1	1:A:422:TYR:HD2	1.71	0.55
1:A:300:LEU:O	1:A:302:GLU:N	2.40	0.55
2:D:160:ASP:C	2:D:162:SER:H	2.15	0.55
1:A:113:THR:HG21	1:A:174:VAL:HB	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:356:THR:O	1:A:369:THR:HA	2.08	0.54
1:A:521:ALA:HB2	2:C:317:PRO:HB2	1.90	0.54
2:C:16:ALA:O	2:C:20:ASN:ND2	2.29	0.54
2:C:174:VAL:HG11	2:C:244:THR:HG21	1.89	0.54
2:D:322:ALA:HA	2:D:386:ARG:O	2.07	0.54
1:A:252:LEU:HA	1:A:363:GLU:HG3	1.88	0.54
1:A:276:GLU:OE2	1:A:401:ASN:HB3	2.07	0.54
1:B:288:PRO:HD3	1:B:294:VAL:HG23	1.88	0.54
1:B:422:TYR:O	1:B:423:PRO:C	2.50	0.54
1:B:496:TYR:OH	1:B:509:GLN:O	2.22	0.54
1:A:508:ASN:O	1:A:514:LYS:NZ	2.36	0.54
2:C:240:ILE:HD13	2:C:495:LEU:HB3	1.89	0.54
2:D:206:THR:HG23	2:D:214:ILE:HG23	1.89	0.54
2:C:494:VAL:HG13	2:C:546:LEU:HD13	1.89	0.54
1:A:342:VAL:HG23	1:A:355:ILE:O	2.08	0.54
2:C:192:THR:HB	2:C:195:SER:HB3	1.89	0.54
1:B:313:LEU:O	1:B:366:GLN:NE2	2.41	0.54
1:A:6:LEU:HG	1:A:65:THR:O	2.07	0.54
1:A:183:ILE:O	1:A:197:ILE:HG13	2.08	0.54
2:C:86:ILE:HD11	2:C:94:ILE:HG13	1.90	0.54
2:C:281:THR:HB	2:C:404:LYS:HB2	1.90	0.54
1:A:351:THR:HG22	1:A:375:ASN:O	2.08	0.53
2:D:322:ALA:HB3	2:D:335:ILE:HD11	1.89	0.53
1:A:304:GLU:HG2	1:A:307:LEU:HD21	1.91	0.53
1:B:178:THR:HB	1:B:472:PRO:HD3	1.89	0.53
1:B:430:ILE:HG23	1:B:516:LYS:HG3	1.91	0.53
1:A:540:LEU:HG	1:A:541:LEU:N	2.22	0.53
2:D:249:SER:HA	2:D:423:THR:HA	1.90	0.53
1:A:334:VAL:HB	1:A:342:VAL:HG22	1.91	0.53
1:B:86:GLU:O	1:B:88:MET:N	2.42	0.53
1:B:287:ILE:HD12	1:B:313:LEU:HD21	1.90	0.53
2:C:304:ASP:H	2:C:380:PRO:HG3	1.74	0.53
2:D:66:GLU:O	2:D:481:GLY:HA2	2.08	0.53
2:C:47:ASN:OD1	2:C:55:TRP:NE1	2.28	0.53
2:C:138:PRO:HA	2:C:146:GLU:OE1	2.09	0.53
1:A:138:THR:HG22	1:A:209:GLY:HA2	1.91	0.52
1:B:153:TYR:CG	1:B:169:LEU:HD23	2.44	0.52
2:C:416:PRO:HD2	2:D:416:PRO:HD2	1.92	0.52
1:A:16:GLU:O	1:A:18:GLU:N	2.42	0.52
2:D:68:THR:OG1	2:D:468:ASP:OD1	2.24	0.52
1:A:425:ILE:HG13	1:A:451:THR:OG1	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:130:LYS:NZ	1:B:185:SER:HB3	2.25	0.52
2:C:26:MET:SD	2:C:141:PHE:HD1	2.33	0.52
1:B:462:ALA:HB3	1:B:485:ALA:HB1	1.92	0.52
2:C:251:ALA:HB1	2:C:454:ARG:HG2	1.92	0.52
1:B:315:ARG:HB3	1:B:316:PRO:HD3	1.91	0.52
2:D:546:LEU:O	2:D:550:ILE:HG13	2.09	0.52
2:C:8:ILE:HA	2:C:31:GLN:O	2.10	0.52
1:A:505:ARG:O	1:A:507:PRO:HD3	2.10	0.51
1:B:246:THR:HG23	1:B:422:TYR:HB3	1.92	0.51
1:B:310:GLU:O	1:B:397:LEU:HD12	2.10	0.51
2:D:4:SER:HB3	2:D:36:TYR:CE1	2.45	0.51
2:D:293:GLY:N	2:D:389:SER:OG	2.41	0.51
1:A:313:LEU:O	1:A:366:GLN:HB2	2.10	0.51
1:A:319:ALA:HB1	1:A:386:VAL:O	2.10	0.51
1:B:357:TYR:HA	1:B:369:THR:HG22	1.92	0.51
1:B:422:TYR:O	1:B:424:ALA:N	2.43	0.51
2:C:128:TRP:CE2	2:C:130:GLN:HG2	2.45	0.51
1:A:107:LEU:HD13	1:A:204:ILE:HD11	1.93	0.51
2:C:287:GLY:HA3	2:C:397:TYR:CZ	2.46	0.51
1:A:313:LEU:HB2	1:A:367:GLN:HB3	1.93	0.51
2:C:114:ASP:OD2	2:C:478:TYR:OH	2.20	0.51
2:C:115:PHE:HE2	2:C:172:THR:HG23	1.76	0.51
2:D:110:TYR:CZ	2:D:161:GLY:HA2	2.46	0.51
2:C:129:ASP:HA	2:C:201:LYS:HB3	1.93	0.50
2:D:255:THR:O	2:D:365:THR:HG23	2.11	0.50
1:A:518:PHE:CZ	2:C:362:PRO:HD3	2.46	0.50
2:C:56:GLU:HG2	2:C:57:GLU:N	2.26	0.50
1:B:135:TRP:HB3	1:B:206:ILE:HG23	1.94	0.50
2:C:172:THR:O	2:C:472:THR:OG1	2.30	0.50
2:D:138:PRO:HB2	2:D:141:PHE:CD2	2.46	0.50
1:B:12:VAL:HG12	1:B:60:ILE:HD12	1.93	0.50
1:B:416:PRO:O	1:B:419:THR:OG1	2.23	0.50
1:A:418:TYR:N	1:A:449:GLY:HA3	2.26	0.50
2:C:119:ASP:OD1	2:C:119:ASP:N	2.37	0.50
1:B:452:ILE:HD12	1:B:452:ILE:N	2.23	0.50
1:B:161:ALA:HB1	1:B:166:ASP:HB3	1.94	0.50
1:B:194:TYR:CD2	1:B:476:TYR:CD2	2.99	0.50
1:B:524:ARG:O	1:B:540:LEU:HG	2.12	0.50
1:B:287:ILE:CD1	1:B:313:LEU:HD21	2.41	0.50
1:B:339:TYR:HA	1:B:357:TYR:O	2.11	0.50
2:C:434:VAL:HG21	2:C:488:SER:HA	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:74:ILE:HG13	1:A:479:ILE:HD12	1.93	0.50
1:B:442:LEU:O	1:B:444:GLN:N	2.45	0.50
1:A:208:LEU:HD11	1:A:220:LEU:HD13	1.94	0.49
1:A:82:VAL:HG13	1:A:467:ILE:HA	1.94	0.49
1:A:400:PRO:HB2	1:A:405:LEU:HD13	1.93	0.49
1:A:467:ILE:O	1:A:478:THR:HA	2.12	0.49
1:B:452:ILE:H	1:B:452:ILE:CD1	2.23	0.49
2:C:342:TYR:CD2	2:C:344:TYR:CE1	3.01	0.49
2:C:349:ASN:O	2:C:350:LEU:HB3	2.12	0.49
1:A:26:MET:C	1:A:28:GLN:H	2.20	0.49
1:A:120:PRO:C	1:A:122:PHE:H	2.21	0.49
1:B:91:ASN:C	1:B:93:PHE:H	2.20	0.49
2:D:262:ASN:OD1	2:D:262:ASN:N	2.45	0.49
1:A:33:ILE:HA	1:A:149:ILE:HD12	1.94	0.49
1:B:133:TYR:HB3	1:B:226:TYR:OH	2.12	0.49
2:C:354:THR:O	2:C:376:ARG:HG2	2.13	0.49
2:D:434:VAL:HG13	2:D:462:ILE:HG13	1.95	0.49
1:A:154:THR:O	1:A:155:ARG:C	2.55	0.49
1:A:246:THR:HG22	1:A:248:SER:N	2.12	0.49
1:B:70:LEU:HA	1:B:244:LEU:HD22	1.94	0.49
2:C:438:ASN:OD1	2:C:441:THR:N	2.46	0.49
2:C:332:SER:HB2	2:C:355:TYR:CE2	2.48	0.49
2:D:100:LEU:HD13	2:D:198:ILE:HD11	1.94	0.49
2:D:109:ASP:C	2:D:111:LEU:H	2.21	0.49
2:D:326:SER:C	2:D:328:SER:H	2.21	0.49
1:B:479:ILE:HG13	1:B:480:THR:N	2.28	0.49
2:D:71:VAL:HG13	2:D:470:ILE:HG23	1.94	0.49
1:A:119:HIS:HB3	1:A:122:PHE:HD2	1.77	0.49
1:A:179:MET:CE	1:A:477:ALA:HB3	2.43	0.49
1:A:443:TRP:NE1	1:A:445:SER:HB2	2.28	0.49
1:B:322:ILE:HG23	1:B:331:SER:O	2.12	0.49
1:A:160:ARG:HB3	1:A:160:ARG:CZ	2.42	0.48
1:B:33:ILE:CG2	1:B:149:ILE:HG12	2.42	0.48
2:D:314:THR:OG1	2:D:369:GLU:N	2.35	0.48
1:A:269:ILE:HB	1:A:429:LEU:HG	1.95	0.48
1:A:532:PRO:HB3	1:A:538:TYR:CZ	2.48	0.48
2:C:42:ASP:O	2:C:44:THR:N	2.47	0.48
2:D:345:ARG:HA	2:D:355:TYR:O	2.14	0.48
1:A:310:GLU:O	1:A:397:LEU:HD12	2.12	0.48
1:B:327:THR:OG1	1:B:349:GLU:HG2	2.13	0.48
2:C:77:ILE:HD13	2:C:177:ILE:HA	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:436:ASN:HB3	1:A:439:ASN:OD1	2.14	0.48
1:A:83:VAL:HG11	1:A:194:TYR:HA	1.94	0.48
1:B:325:SER:HB3	1:B:381:TRP:CZ3	2.48	0.48
1:B:363:GLU:HG2	1:B:364:SER:N	2.23	0.48
2:C:77:ILE:CD1	2:C:177:ILE:HA	2.44	0.48
2:C:369:GLU:OE1	2:C:371:LEU:HB2	2.13	0.48
2:D:498:LEU:HD21	2:D:550:ILE:HG12	1.96	0.48
1:A:237:PRO:HG3	1:A:501:PHE:HD2	1.79	0.48
1:A:443:TRP:O	1:A:445:SER:N	2.43	0.48
1:A:292:GLY:O	1:A:386:VAL:HA	2.14	0.48
1:B:37:LEU:HD22	1:B:222:ALA:HB2	1.96	0.48
1:B:102:THR:OG1	1:B:103:GLY:N	2.47	0.48
2:C:499:MET:HE3	2:C:513:LEU:HD22	1.96	0.48
2:D:41:PHE:HE1	2:D:43:GLU:HA	1.79	0.48
2:D:128:TRP:NE1	2:D:144:GLY:HA3	2.29	0.48
1:A:11:VAL:HG12	1:A:62:THR:HG22	1.95	0.48
1:A:107:LEU:HD21	1:A:132:VAL:HG21	1.95	0.48
1:A:452:ILE:HD12	1:A:452:ILE:N	2.28	0.48
2:C:243:LEU:HB2	2:C:273:SER:HA	1.96	0.48
2:C:444:MET:HB3	2:C:444:MET:HE3	1.70	0.48
1:A:56:LEU:HD23	1:A:57:PRO:HD2	1.96	0.47
2:C:254:THR:HB	2:C:365:THR:O	2.14	0.47
1:B:264:THR:HG22	1:B:267:LEU:HB2	1.97	0.47
2:D:127:LEU:HD13	2:D:199:VAL:HB	1.96	0.47
2:D:243:LEU:HD12	2:D:273:SER:HB2	1.95	0.47
2:D:440:LYS:HE3	2:D:468:ASP:HB2	1.95	0.47
2:D:522:LEU:O	2:D:526:ALA:HB2	2.14	0.47
1:A:144:PRO:HB2	1:A:147:PHE:HD2	1.79	0.47
2:D:438:ASN:HB3	2:D:441:THR:OG1	2.13	0.47
1:B:517:THR:HG23	2:D:317:PRO:HB3	1.96	0.47
2:D:95:THR:HG23	2:D:97:ARG:H	1.79	0.47
1:B:159:ASN:O	1:B:162:ILE:HB	2.15	0.47
1:A:67:LYS:O	1:A:246:THR:OG1	2.24	0.47
2:C:292:VAL:CG2	2:C:392:ILE:HG12	2.42	0.47
2:C:531:ILE:HG23	2:C:532:TYR:CD2	2.49	0.47
2:D:36:TYR:OH	2:D:220:ASP:OD1	2.28	0.47
1:B:459:ASP:CG	1:B:516:LYS:HD3	2.40	0.47
2:C:61:MET:HG2	2:C:247:VAL:HG13	1.97	0.47
2:C:438:ASN:O	2:C:442:ASP:N	2.48	0.47
2:C:527:THR:N	2:C:543:ILE:O	2.45	0.47
1:B:179:MET:SD	1:B:479:ILE:HB	2.54	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:498:GLN:O	1:B:499:TYR:C	2.58	0.47
2:C:464:ALA:HB3	2:C:487:SER:HB3	1.97	0.47
2:D:203:LYS:NZ	2:D:220:ASP:OD1	2.48	0.47
2:D:250:SER:OG	2:D:251:ALA:N	2.48	0.47
1:A:176:GLN:O	1:A:180:LEU:HG	2.15	0.47
1:B:282:HIS:NE2	1:B:396:ASN:OD1	2.45	0.47
1:B:319:ALA:HA	1:B:388:VAL:HG23	1.96	0.47
2:D:7:ILE:HG23	2:D:52:VAL:HG13	1.97	0.47
2:C:110:TYR:CD2	2:C:152:LEU:HD22	2.50	0.46
1:B:54:LEU:HD22	1:B:62:THR:HG21	1.96	0.46
1:B:93:PHE:HD2	1:B:101:ILE:HG13	1.80	0.46
1:A:248:SER:C	1:A:250:ALA:H	2.24	0.46
1:B:451:THR:O	1:B:454:ASP:N	2.48	0.46
2:D:81:ALA:HB2	2:D:490:ILE:CD1	2.45	0.46
2:D:43:GLU:HB2	2:D:55:TRP:CH2	2.50	0.46
2:D:80:ALA:HB1	2:D:465:SER:HB3	1.96	0.46
2:D:147:PHE:HA	2:D:151:GLN:OE1	2.14	0.46
1:A:9:ILE:HA	1:A:64:SER:HA	1.97	0.46
1:B:49:SER:N	1:B:52:ARG:H	2.14	0.46
2:C:392:ILE:O	2:C:393:ILE:HD13	2.16	0.46
1:B:240:ILE:HB	1:B:269:ILE:HG23	1.98	0.46
1:B:59:ILE:HG13	1:B:59:ILE:O	2.15	0.46
2:D:271:VAL:HB	2:D:431:VAL:HG22	1.98	0.46
2:D:320:VAL:O	2:D:359:PHE:HE2	1.99	0.46
1:A:336:ILE:HG22	1:A:337:SER:N	2.30	0.46
1:B:53:VAL:O	1:B:56:LEU:HB2	2.16	0.46
2:C:154:ILE:O	2:C:158:ARG:HG3	2.15	0.46
1:A:170:SER:O	1:A:171:GLN:HG3	2.16	0.46
1:A:248:SER:C	1:A:250:ALA:N	2.74	0.46
1:A:300:LEU:C	1:A:302:GLU:H	2.24	0.46
1:A:402:ARG:O	1:A:403:GLU:HB2	2.14	0.46
1:B:364:SER:HB2	1:B:366:GLN:OE1	2.16	0.46
2:C:379:LYS:HA	2:C:379:LYS:HD3	1.63	0.45
2:D:40:ASP:OD1	2:D:40:ASP:N	2.39	0.45
2:D:324:ILE:HD13	2:D:375:LEU:HD21	1.98	0.45
1:A:84:ALA:C	1:A:86:GLU:N	2.74	0.45
1:B:7:SER:O	1:B:42:ILE:HG12	2.16	0.45
2:C:46:LEU:HD23	2:C:46:LEU:HA	1.67	0.45
2:C:317:PRO:O	2:C:361:TYR:OH	2.26	0.45
2:D:277:ASN:HD21	2:D:483:GLY:HA3	1.81	0.45
1:A:229:LYS:O	1:A:233:GLU:HG3	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:284:HIS:NE2	2:C:398:ASP:OD2	2.46	0.45
2:C:304:ASP:N	2:C:380:PRO:HG3	2.30	0.45
2:C:350:LEU:C	2:C:352:ASN:H	2.25	0.45
2:D:102:ALA:HB1	2:D:228:VAL:HG11	1.98	0.45
2:D:389:SER:HB2	2:D:392:ILE:HG23	1.98	0.45
1:A:275:ASN:OD1	1:A:275:ASN:N	2.49	0.45
1:B:245:GLY:HA2	1:B:423:PRO:CG	2.45	0.45
2:C:227:TYR:O	2:C:231:ILE:HG12	2.17	0.45
1:B:125:PRO:C	1:B:127:GLY:H	2.24	0.45
2:C:134:THR:HB	2:C:135:ASN:H	1.48	0.45
2:C:439:ASN:HB2	2:C:465:SER:OG	2.16	0.45
2:C:548:ASN:OD1	2:C:548:ASN:N	2.48	0.45
2:D:126:TYR:HB3	2:D:227:TYR:OH	2.15	0.45
2:D:275:ALA:HB1	2:D:449:SER:O	2.16	0.45
1:A:312:TRP:HA	1:A:367:GLN:O	2.17	0.45
1:A:451:THR:O	1:A:454:ASP:N	2.50	0.45
1:B:443:TRP:O	1:B:444:GLN:C	2.59	0.45
2:D:8:ILE:HA	2:D:31:GLN:O	2.17	0.45
1:B:119:HIS:HB3	1:B:122:PHE:HD2	1.81	0.45
1:B:467:ILE:O	1:B:478:THR:HA	2.16	0.45
2:C:65:ILE:HD13	2:C:485:GLY:HA3	1.99	0.45
2:C:96:GLY:O	2:C:97:ARG:C	2.59	0.45
2:C:142:ILE:HG13	2:C:143:PHE:HD1	1.82	0.45
2:C:422:ILE:HG22	2:C:426:ALA:HB3	1.97	0.45
2:D:26:MET:SD	2:D:141:PHE:HD1	2.40	0.45
2:D:129:ASP:HA	2:D:201:LYS:HB3	1.97	0.45
2:D:466:GLY:O	2:D:482:THR:HA	2.16	0.45
1:A:398:TYR:OH	1:A:452:ILE:HD11	2.16	0.45
2:C:224:ALA:O	2:C:228:VAL:HG22	2.16	0.45
2:D:332:SER:HB2	2:D:355:TYR:CE2	2.52	0.45
1:A:447:SER:O	1:A:448:ARG:HB2	2.17	0.45
1:B:206:ILE:HD13	1:B:223:ALA:HB1	1.99	0.45
2:D:300:ILE:HG12	2:D:413:PHE:CE1	2.52	0.45
1:A:70:LEU:HA	1:A:244:LEU:CD2	2.47	0.45
1:A:418:TYR:HA	1:A:450:PRO:O	2.17	0.45
1:B:280:GLN:CB	1:B:414:VAL:HG21	2.46	0.45
1:B:307:LEU:HD23	1:B:307:LEU:HA	1.67	0.45
1:A:261:ALA:HB3	1:A:428:ASP:HB2	1.99	0.44
2:C:256:SER:HB2	2:C:424:MET:HE2	1.98	0.44
2:D:318:ASP:OD1	2:D:318:ASP:N	2.50	0.44
1:B:283:THR:HG23	1:B:395:TYR:HE1	1.83	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:62:SER:OG	2:C:63:SER:N	2.47	0.44
2:C:491:VAL:O	2:C:495:LEU:N	2.46	0.44
2:D:96:GLY:HA3	2:D:195:SER:HB3	1.99	0.44
2:D:324:ILE:HB	2:D:348:PHE:CZ	2.51	0.44
2:C:358:ARG:O	2:C:371:LEU:HD12	2.18	0.44
2:C:429:ASP:O	2:C:516:GLN:NE2	2.50	0.44
2:D:54:TRP:CZ2	2:D:363:TYR:HE1	2.36	0.44
2:C:350:LEU:HD12	2:C:351:GLU:H	1.81	0.44
2:C:412:ARG:HH21	2:C:416:PRO:HD3	1.82	0.44
2:D:251:ALA:HB2	2:D:420:ALA:HB1	1.99	0.44
1:B:107:LEU:O	1:B:238:LEU:HA	2.18	0.44
2:C:66:GLU:O	2:C:481:GLY:HA2	2.18	0.44
2:C:453:ILE:O	2:C:456:ARG:HB2	2.17	0.44
1:A:14:THR:O	1:A:15:PRO:C	2.59	0.44
2:D:284:HIS:CD2	2:D:400:TYR:CZ	3.05	0.44
1:A:132:VAL:HG23	1:A:203:LEU:O	2.18	0.44
1:B:300:LEU:HG	1:B:409:THR:HG23	1.99	0.44
1:B:523:ALA:O	1:B:525:LYS:HD3	2.18	0.44
2:C:278:GLN:HB3	2:C:283:ILE:HG12	1.99	0.44
1:A:14:THR:O	1:A:16:GLU:N	2.51	0.44
1:A:133:TYR:HB3	1:A:226:TYR:OH	2.18	0.44
1:B:426:GLN:O	1:B:457:LYS:HE3	2.17	0.44
2:C:43:GLU:HB3	2:C:55:TRP:CH2	2.53	0.44
2:C:149:ARG:HH22	2:C:196:ASP:CG	2.22	0.44
1:B:56:LEU:HD23	1:B:56:LEU:HA	1.77	0.44
1:B:418:TYR:N	1:B:449:GLY:HA3	2.33	0.43
2:C:404:LYS:HA	2:C:407:ILE:HG12	2.00	0.43
2:D:191:ILE:HD11	2:D:490:ILE:HD13	2.00	0.43
2:D:303:GLY:O	2:D:305:ASP:N	2.47	0.43
2:C:416:PRO:HG2	2:D:416:PRO:HD2	2.01	0.43
2:D:116:ILE:HG23	2:D:156:ILE:HD13	2.00	0.43
2:D:397:TYR:HE2	2:D:399:ILE:HD11	1.83	0.43
2:C:38:PRO:O	2:C:40:ASP:N	2.52	0.43
1:B:323:ILE:HD12	1:B:353:TYR:HB2	2.00	0.43
1:A:295:GLU:OE1	1:A:382:LYS:HD2	2.19	0.43
1:A:405:LEU:HB3	1:A:406:LYS:H	1.55	0.43
2:C:187:GLN:H	2:C:187:GLN:HG3	1.68	0.43
1:A:334:VAL:CG1	1:A:340:ASN:HB3	2.43	0.43
1:A:351:THR:HG21	1:A:381:TRP:CZ2	2.53	0.43
1:A:463:PRO:HD3	1:A:540:LEU:HD11	2.00	0.43
1:B:247:SER:HG	1:B:400:PRO:HA	1.82	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:114:ASP:HA	2:C:181:ASN:OD1	2.19	0.43
2:C:243:LEU:HD23	2:C:243:LEU:HA	1.70	0.43
1:A:344:GLY:HA3	1:A:353:TYR:CZ	2.54	0.43
1:A:353:TYR:HA	1:A:372:THR:O	2.19	0.43
1:A:297:GLU:HB2	1:A:412:ARG:HB2	2.01	0.43
1:A:341:LYS:HB2	1:A:341:LYS:HE2	1.62	0.43
1:A:443:TRP:CE2	1:A:445:SER:HB2	2.54	0.43
1:B:304:GLU:H	1:B:376:ALA:HB3	1.83	0.43
1:B:315:ARG:NH2	2:D:516:GLN:OE1	2.49	0.43
1:B:316:PRO:HB3	2:D:523:ILE:CG2	2.49	0.43
1:B:346:PHE:HD1	1:B:346:PHE:HA	1.73	0.43
2:D:138:PRO:CG	2:D:146:GLU:HB2	2.37	0.43
1:A:165:ASN:O	1:A:167:PRO:HD3	2.19	0.43
1:A:249:LEU:O	1:A:251:GLY:N	2.47	0.43
1:B:87:GLU:C	1:B:89:GLY:H	2.27	0.43
1:B:462:ALA:HB1	1:B:463:PRO:HD2	2.00	0.43
1:A:514:LYS:HA	2:C:362:PRO:HG2	2.01	0.43
1:B:185:SER:HB2	1:B:203:LEU:HD11	2.00	0.43
1:B:360:PRO:HB2	2:D:517:VAL:HG13	2.01	0.43
2:D:110:TYR:CE1	2:D:161:GLY:HA2	2.53	0.43
2:D:495:LEU:O	2:D:499:MET:HG3	2.18	0.43
1:A:209:GLY:HA3	1:A:219:MET:HE2	2.00	0.42
1:A:276:GLU:CD	1:A:401:ASN:HB3	2.44	0.42
1:A:336:ILE:CG2	1:A:337:SER:H	2.28	0.42
1:A:346:PHE:HB2	1:A:351:THR:H	1.84	0.42
1:B:294:VAL:HG11	1:B:395:TYR:CE2	2.54	0.42
1:B:316:PRO:HG2	1:B:391:ILE:HB	2.01	0.42
1:B:352:GLU:O	1:B:373:LEU:HD23	2.19	0.42
2:C:41:PHE:CZ	2:C:45:ILE:HB	2.54	0.42
2:C:136:PRO:HA	2:C:137:PRO:HD3	1.92	0.42
2:C:247:VAL:HG23	2:C:423:THR:HG23	2.00	0.42
2:D:9:TYR:HB2	2:D:13:ILE:HB	2.01	0.42
1:A:22:ALA:CA	1:A:27:SER:HB2	2.45	0.42
1:A:175:GLY:O	1:A:176:GLN:C	2.62	0.42
2:C:282:ASP:OD2	2:D:412:ARG:NH2	2.51	0.42
2:C:304:ASP:HA	2:C:380:PRO:HG3	2.01	0.42
2:C:454:ARG:O	2:C:456:ARG:N	2.53	0.42
2:D:60:PRO:O	2:D:247:VAL:HG12	2.19	0.42
2:D:273:SER:OG	2:D:274:GLY:N	2.52	0.42
2:C:510:ARG:C	2:C:512:SER:H	2.28	0.42
2:D:41:PHE:CE1	2:D:43:GLU:HA	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:186:SER:OG	2:D:187:GLN:NE2	2.50	0.42
2:D:524:LEU:HA	2:D:524:LEU:HD23	1.84	0.42
1:B:312:TRP:HB3	1:B:366:GLN:NE2	2.35	0.42
1:A:117:TYR:OH	1:A:166:ASP:O	2.28	0.42
1:A:426:GLN:O	1:A:457:LYS:NZ	2.52	0.42
2:C:127:LEU:HA	2:C:199:VAL:O	2.20	0.42
2:C:419:VAL:HG12	2:C:451:GLY:CA	2.43	0.42
2:D:69:ASN:HA	2:D:479:ASN:OD1	2.19	0.42
1:A:460:ILE:HG13	1:A:461:VAL:N	2.34	0.42
1:B:347:ASP:C	1:B:349:GLU:H	2.28	0.42
1:A:107:LEU:O	1:A:238:LEU:HA	2.19	0.42
1:A:120:PRO:O	1:A:122:PHE:N	2.52	0.42
1:B:78:THR:O	1:B:81:GLY:N	2.53	0.42
1:B:199:GLU:C	1:B:201:SER:H	2.28	0.42
1:B:316:PRO:HB3	2:D:523:ILE:HG21	2.01	0.42
1:B:532:PRO:HA	1:B:537:GLY:O	2.20	0.42
2:D:5:TYR:HB2	2:D:35:ILE:HG12	2.02	0.42
2:D:373:ILE:HD12	2:D:385:LEU:HD11	2.00	0.42
1:A:283:THR:HG21	1:A:296:VAL:HG11	2.01	0.42
2:C:14:GLU:O	2:C:18:GLN:HG3	2.19	0.42
2:C:92:ASN:O	2:C:94:ILE:N	2.52	0.42
2:C:500:GLU:O	2:C:504:LYS:N	2.48	0.42
2:D:302:ASP:OD1	2:D:305:ASP:HB2	2.19	0.42
2:D:453:ILE:HG13	2:D:457:GLY:HA3	2.01	0.42
1:A:347:ASP:C	1:A:349:GLU:H	2.27	0.42
1:A:426:GLN:HB2	1:A:429:LEU:HD13	2.02	0.42
1:A:521:ALA:HB1	2:C:391:LEU:HD21	2.01	0.42
1:B:460:ILE:HD12	1:B:541:LEU:HD13	2.01	0.42
2:C:213:ARG:HD2	2:C:405:ASN:O	2.20	0.42
1:B:91:ASN:C	1:B:93:PHE:N	2.78	0.41
1:B:134:LEU:HA	1:B:205:ILE:O	2.20	0.41
1:B:267:LEU:O	1:B:512:VAL:HG22	2.20	0.41
2:C:58:SER:HB2	2:C:219:SER:H	1.85	0.41
2:C:61:MET:HE2	2:C:61:MET:HB3	1.98	0.41
2:C:510:ARG:C	2:C:512:SER:N	2.78	0.41
2:D:498:LEU:CD2	2:D:550:ILE:HG12	2.50	0.41
1:B:183:ILE:HG12	1:B:469:ALA:HB2	2.02	0.41
1:B:418:TYR:HA	1:B:450:PRO:O	2.19	0.41
1:B:498:GLN:HE21	1:B:499:TYR:N	2.18	0.41
1:A:261:ALA:CB	1:A:428:ASP:HB2	2.50	0.41
1:A:519:MET:O	1:A:523:ALA:HB2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:77:GLY:C	1:B:79:PHE:H	2.28	0.41
1:B:123:ILE:HG23	1:B:162:ILE:HD13	2.03	0.41
2:C:92:ASN:C	2:C:94:ILE:N	2.76	0.41
2:C:252:VAL:HG21	2:C:313:ASN:HB3	2.02	0.41
2:C:454:ARG:C	2:C:456:ARG:H	2.28	0.41
1:A:16:GLU:O	1:A:17:PHE:C	2.62	0.41
1:A:499:TYR:HE2	1:A:551:ARG:HD3	1.86	0.41
1:B:172:ASP:OD2	1:B:178:THR:OG1	2.22	0.41
1:B:194:TYR:HD2	1:B:476:TYR:CE2	2.37	0.41
1:B:341:LYS:NZ	1:B:356:THR:HG21	2.36	0.41
2:C:284:HIS:HE2	2:C:398:ASP:HB3	1.84	0.41
2:C:348:PHE:CE1	2:C:355:TYR:HD2	2.38	0.41
2:D:312:LEU:HD12	2:D:399:ILE:HG12	2.02	0.41
1:A:461:VAL:O	1:A:540:LEU:HA	2.21	0.41
1:B:55:GLN:HE22	2:D:509:PRO:HD2	1.85	0.41
1:B:106:THR:HG21	1:B:494:ALA:O	2.20	0.41
1:B:301:ASN:HB2	1:B:408:GLY:HA3	2.02	0.41
1:B:475:THR:HG22	1:B:476:TYR:H	1.86	0.41
1:B:511:GLU:O	1:B:515:ILE:HG13	2.20	0.41
2:C:100:LEU:HD13	2:C:198:ILE:HD11	2.03	0.41
2:C:149:ARG:NH2	2:C:196:ASP:OD1	2.37	0.41
2:C:250:SER:OG	2:C:251:ALA:N	2.50	0.41
1:A:34:SER:HA	1:A:147:PHE:HA	2.02	0.41
1:A:172:ASP:OD1	1:A:174:VAL:O	2.38	0.41
1:B:498:GLN:O	1:B:500:THR:N	2.53	0.41
2:C:366:SER:O	2:C:368:LYS:N	2.54	0.41
1:A:255:ARG:HB2	1:A:255:ARG:NH1	2.35	0.41
1:A:300:LEU:C	1:A:302:GLU:N	2.79	0.41
1:A:402:ARG:C	1:A:404:LEU:N	2.76	0.41
1:B:498:GLN:O	1:B:502:VAL:O	2.39	0.41
2:C:308:LEU:HD12	2:C:308:LEU:HA	1.88	0.41
2:C:554:ALA:O	2:C:556:THR:N	2.53	0.41
1:B:84:ALA:HB3	1:B:488:HIS:NE2	2.36	0.41
1:B:107:LEU:HD23	1:B:107:LEU:HA	1.94	0.41
2:C:125:LEU:HG	2:C:126:TYR:CE2	2.56	0.41
2:C:280:ASN:OD1	2:C:447:GLY:HA2	2.21	0.41
2:C:463:VAL:HG23	2:C:543:ILE:HA	2.02	0.41
2:D:127:LEU:O	2:D:146:GLU:HA	2.21	0.41
1:A:177:GLY:O	1:A:178:THR:C	2.63	0.41
1:A:427:ASP:C	1:A:429:LEU:H	2.28	0.41
1:B:264:THR:HG21	1:B:267:LEU:HD22	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:298:LEU:HD13	1:B:309:LEU:HD21	2.02	0.41
1:B:541:LEU:HA	1:B:541:LEU:HD12	1.84	0.41
2:C:225:ILE:HD13	2:C:225:ILE:HA	1.86	0.41
2:C:466:GLY:HA2	2:C:469:ILE:HD13	2.03	0.41
2:D:105:ASP:OD1	2:D:170:THR:OG1	2.27	0.41
2:D:109:ASP:OD2	2:D:111:LEU:HB2	2.21	0.41
2:C:258:LEU:HD22	2:C:264:LEU:CD1	2.51	0.41
2:C:523:ILE:HG23	2:C:541:TYR:HB3	2.03	0.41
2:C:553:ILE:O	2:C:557:LEU:HG	2.21	0.41
2:D:535:PRO:HB3	2:D:541:TYR:CD1	2.56	0.41
1:A:334:VAL:HB	1:A:342:VAL:CG2	2.51	0.40
1:A:516:LYS:O	1:A:520:GLN:HG3	2.20	0.40
2:D:544:LEU:HA	2:D:544:LEU:HD12	1.78	0.40
1:A:74:ILE:HD13	1:A:174:VAL:HG12	2.02	0.40
1:B:240:ILE:HG21	1:B:240:ILE:HD13	1.78	0.40
1:B:443:TRP:HB2	1:B:465:VAL:CG2	2.50	0.40
2:C:135:ASN:HB2	2:C:147:PHE:CE1	2.56	0.40
2:C:261:PHE:C	2:C:263:ILE:H	2.30	0.40
2:C:404:LYS:HD2	2:C:407:ILE:HD11	2.03	0.40
2:D:123:LYS:C	2:D:149:ARG:HH21	2.30	0.40
1:A:33:ILE:HA	1:A:149:ILE:CD1	2.52	0.40
1:A:460:ILE:HD12	1:A:541:LEU:HD13	2.04	0.40
1:B:129:SER:N	1:B:159:ASN:OD1	2.54	0.40
2:C:209:TYR:HB3	2:C:307:ALA:O	2.21	0.40
2:C:422:ILE:H	2:C:422:ILE:HG12	1.55	0.40
2:D:130:GLN:OE1	2:D:202:LEU:HA	2.21	0.40
1:A:10:ASP:N	1:A:63:VAL:O	2.55	0.40
1:A:77:GLY:C	1:A:79:PHE:H	2.29	0.40
1:A:117:TYR:OH	1:A:167:PRO:HA	2.21	0.40
1:A:285:GLY:O	1:A:394:ARG:HA	2.21	0.40
1:A:354:SER:OG	1:A:372:THR:HB	2.22	0.40
1:A:388:VAL:H	1:A:388:VAL:HG23	1.64	0.40
1:A:447:SER:O	1:A:448:ARG:CB	2.70	0.40
1:B:266:GLY:HA2	1:B:497:PHE:CE1	2.57	0.40
2:D:83:THR:O	2:D:86:ILE:N	2.52	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	520/551 (94%)	409 (79%)	98 (19%)	13 (2%)	4	20
1	B	520/551 (94%)	411 (79%)	98 (19%)	11 (2%)	5	23
2	C	554/565 (98%)	469 (85%)	64 (12%)	21 (4%)	2	14
2	D	545/565 (96%)	478 (88%)	59 (11%)	8 (2%)	8	29
All	All	2139/2232 (96%)	1767 (83%)	319 (15%)	53 (2%)	4	20

All (53) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	17	PHE
1	A	305	GLU
1	B	444	GLN
2	C	39	VAL
2	C	40	ASP
2	C	304	ASP
2	C	509	PRO
2	C	510	ARG
2	D	69	ASN
2	D	134	THR
1	A	301	ASN
1	A	388	VAL
1	A	526	ASP
1	B	117	TYR
2	C	43	GLU
2	C	134	THR
2	C	183	ARG
2	C	367	GLY
2	C	429	ASP
2	C	455	GLY
1	A	327	THR
1	B	44	ASN

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Mol	Chain	Res	Type
1	B	85	THR
1	B	112	ASP
1	B	377	LYS
2	C	85	TYR
2	C	352	ASN
2	C	466	GLY
2	C	555	ASN
2	D	53	ALA
2	D	130	GLN
2	D	236	ASN
1	A	85	THR
1	A	333	SER
2	C	210	TYR
2	D	110	TYR
2	D	466	GLY
1	A	112	ASP
1	A	336	ILE
1	B	29	ASN
1	B	142	THR
1	B	447	SER
2	C	86	ILE
2	C	444	MET
2	C	454	ARG
1	B	315	ARG
2	C	184	ILE
2	C	276	GLY
1	A	251	GLY
1	A	502	VAL
1	A	142	THR
2	D	71	VAL
1	B	251	GLY

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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	398/460 (86%)	303 (76%)	95 (24%)	1 2
1	B	383/460 (83%)	302 (79%)	81 (21%)	1 4
2	C	448/491 (91%)	380 (85%)	68 (15%)	3 11
2	D	439/491 (89%)	363 (83%)	76 (17%)	2 8
All	All	1668/1902 (88%)	1348 (81%)	320 (19%)	1 5

All (320) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	LEU
1	A	9	ILE
1	A	11	VAL
1	A	12	VAL
1	A	14	THR
1	A	24	LEU
1	A	33	ILE
1	A	34	SER
1	A	35	ASP
1	A	36	SER
1	A	41	SER
1	A	44	ASN
1	A	53	VAL
1	A	58	SER
1	A	59	ILE
1	A	60	ILE
1	A	63	VAL
1	A	64	SER
1	A	66	THR
1	A	70	LEU
1	A	76	ARG
1	A	82	VAL
1	A	83	VAL
1	A	100	ASN
1	A	102	THR
1	A	126	ASP
1	A	128	THR
1	A	132	VAL
1	A	134	LEU
1	A	138	THR

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Mol	Chain	Res	Type
1	A	142	THR
1	A	149	ILE
1	A	151	THR
1	A	160	ARG
1	A	174	VAL
1	A	184	CYS
1	A	192	SER
1	A	197	ILE
1	A	202	GLU
1	A	210	LYS
1	A	219	MET
1	A	243	SER
1	A	247	SER
1	A	253	THR
1	A	255	ARG
1	A	267	LEU
1	A	269	ILE
1	A	286	ILE
1	A	290	VAL
1	A	293	SER
1	A	294	VAL
1	A	297	GLU
1	A	300	LEU
1	A	304	GLU
1	A	306	GLU
1	A	309	LEU
1	A	313	LEU
1	A	333	SER
1	A	334	VAL
1	A	340	ASN
1	A	342	VAL
1	A	345	LEU
1	A	348	LEU
1	A	352	GLU
1	A	356	THR
1	A	374	LYS
1	A	380	VAL
1	A	382	LYS
1	A	385	LEU
1	A	392	THR
1	A	401	ASN
1	A	405	LEU

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Mol	Chain	Res	Type
1	A	407	SER
1	A	412	ARG
1	A	419	THR
1	A	425	ILE
1	A	426	GLN
1	A	427	ASP
1	A	429	LEU
1	A	430	ILE
1	A	437	THR
1	A	439	ASN
1	A	451	THR
1	A	456	LEU
1	A	460	ILE
1	A	475	THR
1	A	479	ILE
1	A	502	VAL
1	A	527	SER
1	A	528	ASN
1	A	529	THR
1	A	534	THR
1	A	535	ASN
1	A	536	SER
1	A	540	LEU
1	B	9	ILE
1	B	18	GLU
1	B	27	SER
1	B	36	SER
1	B	37	LEU
1	B	40	LEU
1	B	41	SER
1	B	44	ASN
1	B	45	THR
1	B	54	LEU
1	B	59	ILE
1	B	60	ILE
1	B	63	VAL
1	B	64	SER
1	B	69	THR
1	B	70	LEU
1	B	73	GLU
1	B	82	VAL
1	B	83	VAL

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Mol	Chain	Res	Type
1	B	87	GLU
1	B	110	ILE
1	B	118	LEU
1	B	126	ASP
1	B	128	THR
1	B	130	LYS
1	B	149	ILE
1	B	154	THR
1	B	155	ARG
1	B	159	ASN
1	B	176	GLN
1	B	187	LEU
1	B	192	SER
1	B	210	LYS
1	B	214	PHE
1	B	229	LYS
1	B	234	LEU
1	B	246	THR
1	B	248	SER
1	B	249	LEU
1	B	267	LEU
1	B	269	ILE
1	B	290	VAL
1	B	300	LEU
1	B	307	LEU
1	B	309	LEU
1	B	311	LEU
1	B	324	VAL
1	B	327	THR
1	B	331	SER
1	B	340	ASN
1	B	342	VAL
1	B	346	PHE
1	B	361	THR
1	B	362	THR
1	B	372	THR
1	B	374	LYS
1	B	380	VAL
1	B	385	LEU
1	B	386	VAL
1	B	390	ILE
1	B	392	THR

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Mol	Chain	Res	Type
1	B	394	ARG
1	B	402	ARG
1	B	404	LEU
1	B	419	THR
1	B	420	ILE
1	B	428	ASP
1	B	432	VAL
1	B	437	THR
1	B	441	SER
1	B	451	THR
1	B	453	GLU
1	B	456	LEU
1	B	468	ILE
1	B	475	THR
1	B	479	ILE
1	B	498	GLN
1	B	529	THR
1	B	530	VAL
1	B	536	SER
1	B	540	LEU
2	C	2	GLU
2	C	4	SER
2	C	25	TYR
2	C	26	MET
2	C	27	VAL
2	C	28	LEU
2	C	31	GLN
2	C	39	VAL
2	C	50	ILE
2	C	57	GLU
2	C	63	SER
2	C	65	ILE
2	C	71	VAL
2	C	72	ASN
2	C	77	ILE
2	C	86	ILE
2	C	94	ILE
2	C	97	ARG
2	C	104	ILE
2	C	111	LEU
2	C	118	ASP
2	C	119	ASP

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Mol	Chain	Res	Type
2	C	121	THR
2	C	131	GLU
2	C	133	ASN
2	C	134	THR
2	C	148	THR
2	C	159	ASN
2	C	163	LEU
2	C	173	LEU
2	C	183	ARG
2	C	192	THR
2	C	204	SER
2	C	208	THR
2	C	209	TYR
2	C	215	ASN
2	C	247	VAL
2	C	255	THR
2	C	312	LEU
2	C	324	ILE
2	C	335	ILE
2	C	345	ARG
2	C	354	THR
2	C	360	ILE
2	C	364	ILE
2	C	365	THR
2	C	366	SER
2	C	376	ARG
2	C	391	LEU
2	C	405	ASN
2	C	419	VAL
2	C	421	THR
2	C	422	ILE
2	C	423	THR
2	C	432	ILE
2	C	446	ILE
2	C	450	LYS
2	C	456	ARG
2	C	477	THR
2	C	480	THR
2	C	490	ILE
2	C	510	ARG
2	C	512	SER
2	C	531	ILE

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Mol	Chain	Res	Type
2	C	543	ILE
2	C	544	LEU
2	C	550	ILE
2	C	553	ILE
2	D	4	SER
2	D	10	GLN
2	D	13	ILE
2	D	20	ASN
2	D	25	TYR
2	D	26	MET
2	D	31	GLN
2	D	34	VAL
2	D	39	VAL
2	D	40	ASP
2	D	43	GLU
2	D	50	ILE
2	D	56	GLU
2	D	71	VAL
2	D	73	ASN
2	D	79	THR
2	D	83	THR
2	D	93	ASP
2	D	104	ILE
2	D	118	ASP
2	D	119	ASP
2	D	135	ASN
2	D	148	THR
2	D	150	SER
2	D	151	GLN
2	D	158	ARG
2	D	165	GLN
2	D	168	ILE
2	D	173	LEU
2	D	193	THR
2	D	203	LYS
2	D	206	THR
2	D	209	TYR
2	D	214	ILE
2	D	226	THR
2	D	235	GLU
2	D	244	THR
2	D	247	VAL

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Mol	Chain	Res	Type
2	D	255	THR
2	D	262	ASN
2	D	291	SER
2	D	299	ILE
2	D	305	ASP
2	D	311	THR
2	D	312	LEU
2	D	335	ILE
2	D	349	ASN
2	D	352	ASN
2	D	370	ASN
2	D	388	THR
2	D	390	GLU
2	D	391	LEU
2	D	393	ILE
2	D	394	SER
2	D	405	ASN
2	D	410	ASP
2	D	419	VAL
2	D	421	THR
2	D	422	ILE
2	D	423	THR
2	D	429	ASP
2	D	431	VAL
2	D	432	ILE
2	D	439	ASN
2	D	446	ILE
2	D	450	LYS
2	D	474	LYS
2	D	475	ASN
2	D	480	THR
2	D	490	ILE
2	D	511	LEU
2	D	533	THR
2	D	543	ILE
2	D	550	ILE
2	D	553	ILE
2	D	556	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	44	ASN
1	A	176	GLN
1	A	366	GLN
1	A	401	ASN
1	A	421	ASN
1	B	55	GLN
1	B	176	GLN
1	B	301	ASN
1	B	421	ASN
1	B	498	GLN
1	B	520	GLN
2	C	70	ASN
2	C	73	ASN
2	C	323	GLN
2	C	545	ASN
2	D	236	ASN
2	D	439	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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	528/551 (95%)	0.03	10 (1%) 66 49	62, 72, 88, 98	0
1	B	530/551 (96%)	0.18	23 (4%) 40 28	77, 93, 108, 118	0
2	C	556/565 (98%)	0.02	9 (1%) 70 55	63, 76, 92, 102	0
2	D	551/565 (97%)	-0.00	4 (0%) 84 73	65, 81, 98, 105	0
All	All	2165/2232 (96%)	0.06	46 (2%) 63 47	62, 80, 101, 118	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	197	LEU	4.0
1	B	298	LEU	3.9
2	C	478	TYR	3.8
2	D	227	TYR	3.7
1	A	434	ALA	3.6
1	B	53	VAL	3.2
1	B	497	PHE	2.9
1	B	484	ALA	2.8
1	B	354	SER	2.8
1	B	204	ILE	2.8
2	D	222	LEU	2.7
2	D	522	LEU	2.7
1	B	23	GLY	2.7
1	B	464	GLY	2.7
1	A	149	ILE	2.7
1	B	184	CYS	2.7
2	C	469	ILE	2.5
1	B	307	LEU	2.5
2	C	381	GLY	2.5
1	B	465	VAL	2.5
1	A	389	TYR	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	296	VAL	2.4
1	A	108	ILE	2.4
2	C	371	LEU	2.4
1	A	110	ILE	2.3
2	C	300	ILE	2.3
1	B	250	ALA	2.3
2	D	375	LEU	2.3
2	C	493	GLY	2.3
1	B	299	GLU	2.2
1	B	108	ILE	2.2
1	A	218	ALA	2.2
1	B	88	MET	2.2
1	A	490	ALA	2.2
1	A	296	VAL	2.2
1	B	50	ILE	2.2
1	B	89	GLY	2.2
1	B	249	LEU	2.1
2	C	557	LEU	2.1
1	B	220	LEU	2.1
2	C	298	VAL	2.1
1	B	32	LYS	2.1
1	A	63	VAL	2.0
1	B	391	ILE	2.0
1	A	422	TYR	2.0
1	B	93	PHE	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.