



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 10:27 AM UTC

PDB ID : 6QVI / pdb_00006qvi
Title : Crystal structure of competence-associated pilin ComZ from *Thermus thermophilus*
Authors : Karuppiah, V.; Derrick, J.P.
Deposited on : 2019-03-01
Resolution : 2.72 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

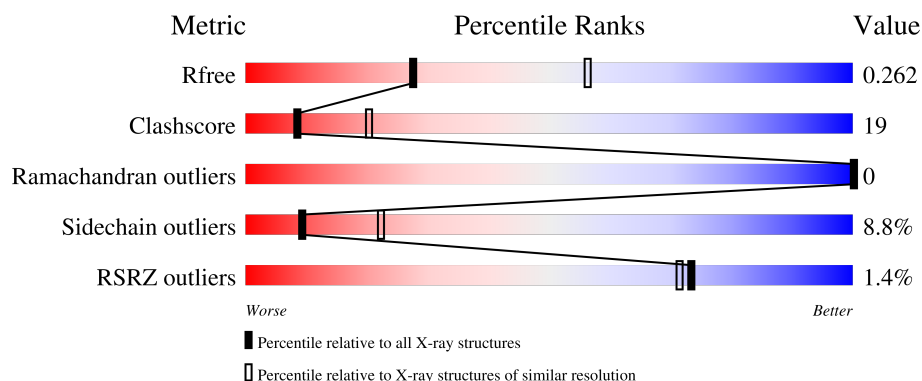
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.72 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	4348 (2.74-2.70)
Clashscore	190562	4665 (2.74-2.70)
Ramachandran outliers	187476	4584 (2.74-2.70)
Sidechain outliers	187428	4585 (2.74-2.70)
RSRZ outliers	180081	4348 (2.74-2.70)

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	534	
1	B	534	
1	C	534	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11520 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 ComZ.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	518	Total	C	N	O	S	0	0	0
			3933	2457	699	763	14			
1	B	498	Total	C	N	O	S	0	0	0
			3761	2348	662	737	14			
1	C	507	Total	C	N	O	S	0	0	0
			3811	2390	664	743	14			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	initiating methionine	UNP Q72JC1
A	0	ALA	-	expression tag	UNP Q72JC1
A	525	LEU	-	expression tag	UNP Q72JC1
A	526	GLU	-	expression tag	UNP Q72JC1
A	527	HIS	-	expression tag	UNP Q72JC1
A	528	HIS	-	expression tag	UNP Q72JC1
A	529	HIS	-	expression tag	UNP Q72JC1
A	530	HIS	-	expression tag	UNP Q72JC1
A	531	HIS	-	expression tag	UNP Q72JC1
A	532	HIS	-	expression tag	UNP Q72JC1
B	-1	MET	-	initiating methionine	UNP Q72JC1
B	0	ALA	-	expression tag	UNP Q72JC1
B	525	LEU	-	expression tag	UNP Q72JC1
B	526	GLU	-	expression tag	UNP Q72JC1
B	527	HIS	-	expression tag	UNP Q72JC1
B	528	HIS	-	expression tag	UNP Q72JC1
B	529	HIS	-	expression tag	UNP Q72JC1
B	530	HIS	-	expression tag	UNP Q72JC1
B	531	HIS	-	expression tag	UNP Q72JC1
B	532	HIS	-	expression tag	UNP Q72JC1
C	-1	MET	-	initiating methionine	UNP Q72JC1
C	0	ALA	-	expression tag	UNP Q72JC1
C	525	LEU	-	expression tag	UNP Q72JC1

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Chain	Residue	Modelled	Actual	Comment	Reference
C	526	GLU	-	expression tag	UNP Q72JC1
C	527	HIS	-	expression tag	UNP Q72JC1
C	528	HIS	-	expression tag	UNP Q72JC1
C	529	HIS	-	expression tag	UNP Q72JC1
C	530	HIS	-	expression tag	UNP Q72JC1
C	531	HIS	-	expression tag	UNP Q72JC1
C	532	HIS	-	expression tag	UNP Q72JC1

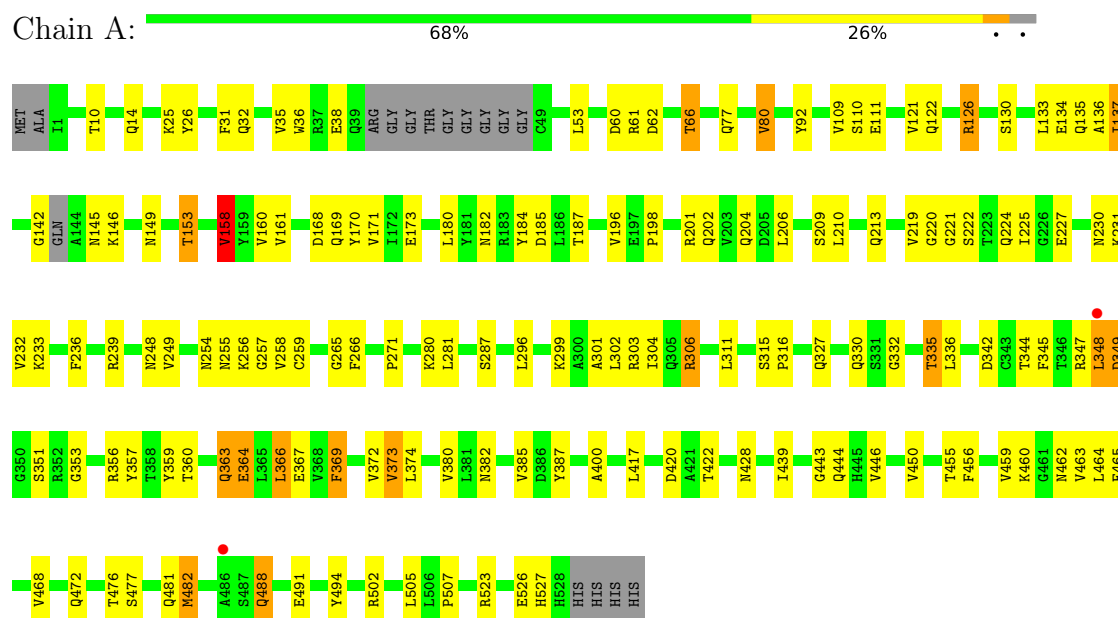
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	7	Total O 7 7	0	0
2	B	4	Total O 4 4	0	0
2	C	4	Total O 4 4	0	0

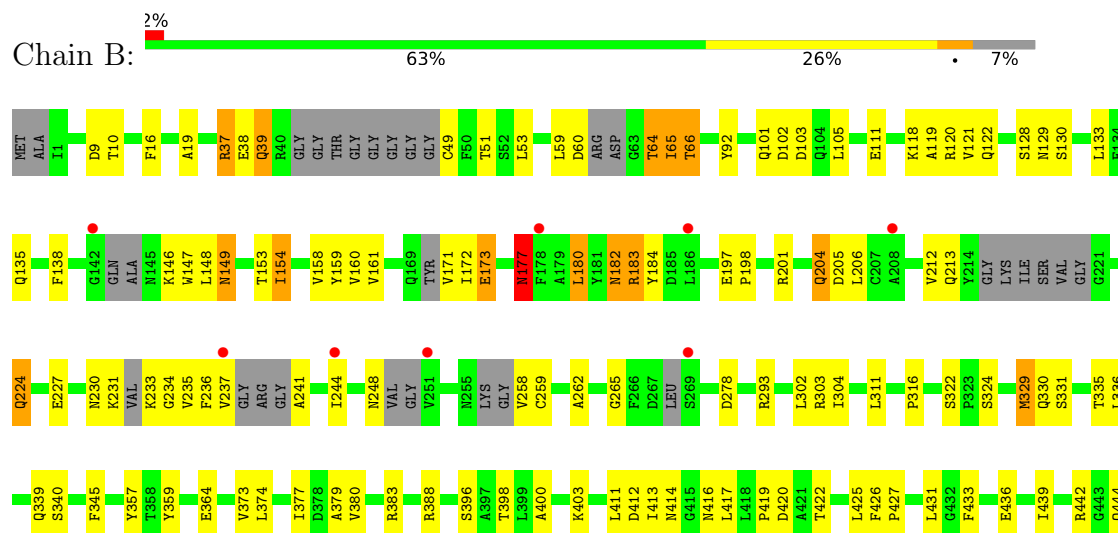
3 Residue-property plots

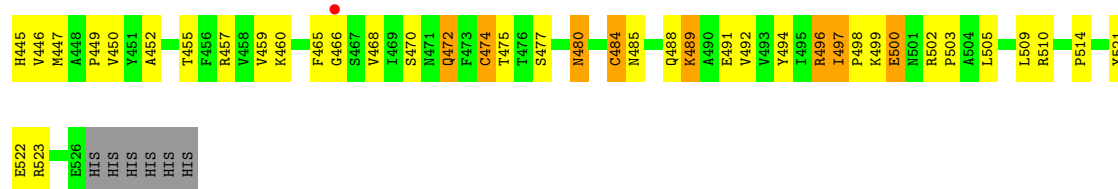
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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

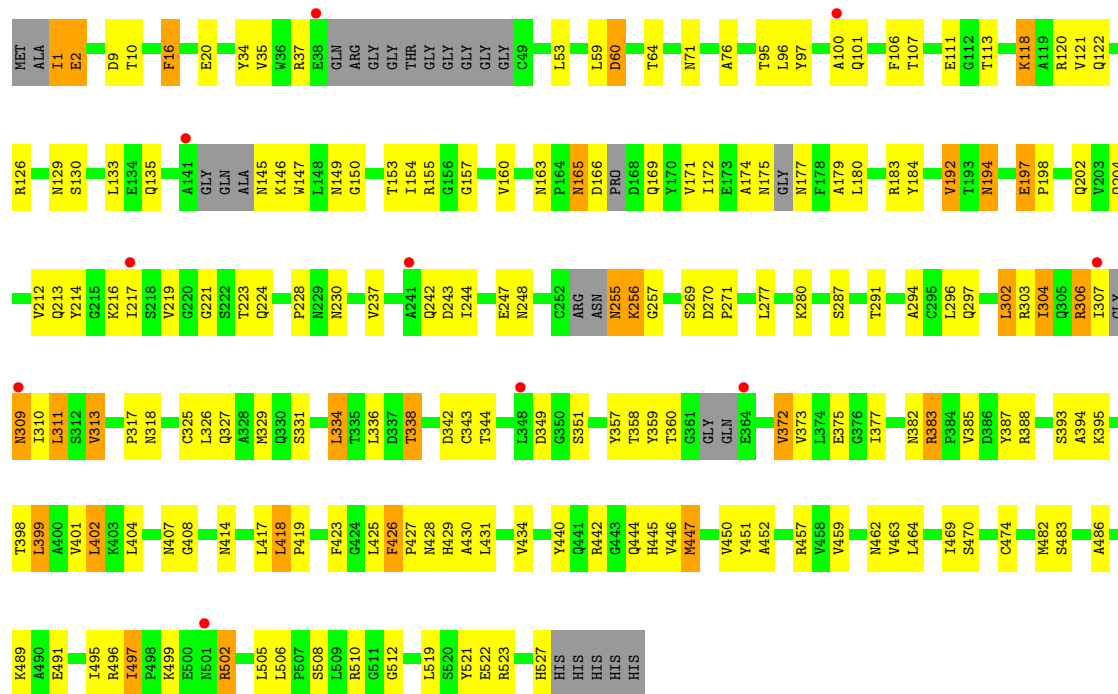


• Molecule 1: ComZ





• Molecule 1: ComZ



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	121.86Å 121.86Å 212.35Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	58.79 – 2.72 58.79 – 2.72	Depositor EDS
% Data completeness (in resolution range)	99.8 (58.79-2.72) 99.8 (58.79-2.72)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.19 (at 2.73Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.206 , 0.267 0.218 , 0.262	Depositor DCC
R_{free} test set	2515 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	75.7	Xtriage
Anisotropy	0.455	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 68.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.025 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11520	wwPDB-VP
Average B, all atoms (Å ²)	96.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.74% 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

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	1.00	6/4006 (0.1%)	0.98	14/5439 (0.3%)
1	B	0.81	1/3821 (0.0%)	0.91	4/5182 (0.1%)
1	C	1.00	1/3877 (0.0%)	1.04	16/5266 (0.3%)
All	All	0.94	8/11704 (0.1%)	0.98	34/15887 (0.2%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	407	ASN	CA-C	-6.13	1.45	1.52
1	B	66	THR	C-O	-5.78	1.20	1.25
1	A	306	ARG	CA-C	-5.59	1.47	1.53
1	A	137	ILE	C-O	-5.46	1.18	1.24
1	A	304	ILE	C-O	-5.36	1.18	1.24
1	A	374	LEU	C-O	-5.26	1.17	1.24
1	A	158	VAL	C-O	-5.21	1.18	1.24
1	A	111	GLU	C-O	-5.03	1.17	1.24

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	408	GLY	N-CA-C	9.04	124.91	112.52
1	B	489	LYS	N-CA-C	8.42	122.62	108.90
1	C	76	ALA	CA-C-N	-8.02	110.39	122.74
1	C	76	ALA	C-N-CA	-8.02	110.39	122.74
1	C	428	ASN	N-CA-C	-7.58	103.92	113.02
1	A	271	PRO	CA-C-N	7.04	127.64	120.38
1	A	271	PRO	C-N-CA	7.04	127.64	120.38
1	C	197	GLU	CA-C-N	7.02	128.61	119.84
1	C	197	GLU	C-N-CA	7.02	128.61	119.84
1	A	428	ASN	N-CA-C	-6.31	105.86	112.93
1	B	484	CYS	N-CA-C	6.23	119.90	112.93
1	C	383	ARG	CA-C-N	6.13	126.49	119.93
1	C	383	ARG	C-N-CA	6.13	126.49	119.93

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	481	GLN	N-CA-C	6.12	119.58	110.52
1	C	192	VAL	CB-CA-C	-6.04	104.95	110.63
1	C	313	VAL	CB-CA-C	-5.97	105.47	111.80
1	C	304	ILE	CB-CA-C	-5.88	101.74	110.33
1	A	80	VAL	N-CA-C	5.74	117.26	108.71
1	B	177	ASN	N-CA-C	-5.59	104.83	111.03
1	A	126	ARG	N-CA-C	-5.58	99.31	108.41
1	C	326	LEU	N-CA-C	5.58	118.13	111.71
1	A	66	THR	N-CA-C	-5.57	99.85	109.15
1	A	153	THR	N-CA-C	-5.51	98.21	108.02
1	A	363	GLN	CB-CG-CD	-5.42	103.38	112.60
1	A	301	ALA	N-CA-C	-5.33	104.90	111.40
1	C	224	GLN	N-CA-C	5.30	116.97	107.80
1	C	16	PHE	N-CA-C	-5.29	105.52	111.28
1	B	330	GLN	N-CA-C	5.26	118.48	111.75
1	C	338	THR	N-CA-C	5.23	121.95	110.80
1	A	281	LEU	CA-C-N	-5.18	115.53	122.42
1	A	281	LEU	C-N-CA	-5.18	115.53	122.42
1	A	488	GLN	N-CA-CB	5.16	118.38	110.49
1	C	394	ALA	N-CA-C	5.03	116.60	110.41
1	A	364	GLU	N-CA-C	5.02	121.50	110.80

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	3933	0	3830	111	0
1	B	3761	0	3634	155	0
1	C	3811	0	3676	165	0
2	A	7	0	0	0	0
2	B	4	0	0	0	0
2	C	4	0	0	0	0
All	All	11520	0	11140	431	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 19.

All (431) 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:256:LYS:HG3	1:C:257:GLY:H	1.07	1.12
1:C:372:VAL:HG22	1:C:399:LEU:HD22	1.33	1.05
1:C:256:LYS:CG	1:C:257:GLY:H	1.73	1.01
1:B:496:ARG:HH11	1:B:496:ARG:HG2	1.24	1.01
1:B:153:THR:CG2	1:B:489:LYS:HD2	1.92	0.99
1:C:256:LYS:CG	1:C:257:GLY:N	2.23	0.96
1:C:414:ASN:HD21	1:C:442:ARG:HH11	1.09	0.96
1:B:497:ILE:HD12	1:B:497:ILE:H	1.31	0.95
1:B:455:THR:HG23	1:B:472:GLN:HB3	1.49	0.93
1:C:160:VAL:HG11	1:C:171:VAL:HG21	1.51	0.93
1:A:248:ASN:HB3	1:A:257:GLY:HA2	1.51	0.93
1:C:306:ARG:HG2	1:C:306:ARG:HH11	1.34	0.91
1:C:296:LEU:HD23	1:C:505:LEU:HD23	1.53	0.91
1:B:329:MET:HE2	1:B:377:ILE:HD13	1.56	0.87
1:B:182:ASN:O	1:B:182:ASN:ND2	2.07	0.87
1:A:373:VAL:HB	1:A:400:ALA:HB3	1.57	0.87
1:C:256:LYS:HG3	1:C:257:GLY:N	1.84	0.86
1:B:183:ARG:HG2	1:B:183:ARG:HH11	1.40	0.86
1:B:431:LEU:HD23	1:B:433:PHE:CE2	2.11	0.85
1:B:153:THR:HG23	1:B:489:LYS:HD2	1.55	0.85
1:B:146:LYS:HZ1	1:B:173:GLU:CG	1.89	0.85
1:C:497:ILE:HD12	1:C:497:ILE:H	1.40	0.85
1:C:446:VAL:CG2	1:C:464:LEU:HD13	2.06	0.84
1:A:335:THR:HB	1:A:380:VAL:HG22	1.60	0.82
1:A:347:ARG:HG3	1:A:347:ARG:HH11	1.43	0.82
1:B:227:GLU:O	1:B:230:ASN:O	1.97	0.81
1:B:414:ASN:HD21	1:B:442:ARG:HH11	1.29	0.81
1:C:113:THR:HG22	1:C:118:LYS:HG3	1.65	0.79
1:A:349:ASP:OD2	1:A:349:ASP:N	2.12	0.79
1:B:496:ARG:HG2	1:B:496:ARG:NH1	1.96	0.79
1:A:348:LEU:HD23	1:A:348:LEU:H	1.47	0.79
1:C:334:LEU:HD23	1:C:334:LEU:O	1.83	0.78
1:B:153:THR:HG23	1:B:489:LYS:CD	2.14	0.77
1:C:426:PHE:HE2	1:C:430:ALA:CB	1.97	0.77
1:B:153:THR:CG2	1:B:489:LYS:CD	2.63	0.76
1:A:336:LEU:HB3	1:A:357:TYR:CE2	2.20	0.76
1:C:372:VAL:HG22	1:C:399:LEU:CD2	2.13	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:256:LYS:HG2	1:C:257:GLY:N	2.02	0.75
1:C:342:ASP:OD1	1:C:344:THR:HG22	1.86	0.75
1:A:347:ARG:HG3	1:A:347:ARG:NH1	2.02	0.74
1:B:154:ILE:HG23	1:B:492:VAL:HG13	1.69	0.74
1:B:120:ARG:NH2	1:B:122:GLN:OE1	2.21	0.74
1:C:174:ALA:HB1	1:C:223:THR:HG21	1.69	0.74
1:B:180:LEU:C	1:B:180:LEU:HD12	2.13	0.73
1:A:219:VAL:HG12	1:A:248:ASN:HB2	1.69	0.73
1:B:224:GLN:O	1:B:224:GLN:HG2	1.88	0.73
1:B:416:ASN:HD21	1:B:444:GLN:H	1.36	0.73
1:C:414:ASN:ND2	1:C:442:ARG:HH11	1.84	0.72
1:C:446:VAL:HG23	1:C:464:LEU:HD13	1.71	0.72
1:C:256:LYS:HB3	1:C:256:LYS:HZ2	1.55	0.72
1:B:278:ASP:OD1	1:B:293:ARG:NH1	2.23	0.72
1:B:177:ASN:OD1	1:B:177:ASN:N	2.20	0.71
1:B:480:ASN:N	1:B:480:ASN:OD1	2.19	0.71
1:C:96:LEU:HD11	1:C:106:PHE:HB3	1.71	0.71
1:C:414:ASN:HD21	1:C:442:ARG:NH1	1.85	0.70
1:B:248:ASN:HB3	1:B:258:VAL:HG12	1.73	0.70
1:C:194:ASN:OD1	1:C:194:ASN:N	2.18	0.70
1:C:426:PHE:CE2	1:C:430:ALA:HB2	2.27	0.70
1:A:446:VAL:HG13	1:A:464:LEU:HD13	1.74	0.70
1:B:180:LEU:HD12	1:B:180:LEU:O	1.92	0.70
1:B:414:ASN:ND2	1:B:442:ARG:HH11	1.90	0.70
1:C:426:PHE:HE2	1:C:430:ALA:HB2	1.55	0.69
1:B:416:ASN:HD21	1:B:444:GLN:HB2	1.57	0.69
1:C:497:ILE:HD12	1:C:497:ILE:N	2.08	0.69
1:B:65:ILE:N	1:B:65:ILE:HD12	2.07	0.69
1:A:231:LYS:HD2	1:A:259:CYS:HB3	1.74	0.69
1:B:447:MET:HG2	1:B:465:PHE:HB2	1.73	0.69
1:B:497:ILE:HD12	1:B:497:ILE:N	2.07	0.68
1:C:256:LYS:HB3	1:C:256:LYS:NZ	2.07	0.68
1:B:121:VAL:HG12	1:B:523:ARG:HG3	1.76	0.68
1:B:147:TRP:CE3	1:B:171:VAL:HG23	2.27	0.68
1:A:348:LEU:HD23	1:A:348:LEU:N	2.07	0.68
1:B:183:ARG:HB2	1:B:204:GLN:HA	1.75	0.68
1:C:197:GLU:HG3	1:C:198:PRO:HD2	1.74	0.68
1:B:359:TYR:HA	1:B:364:GLU:HG3	1.76	0.68
1:A:366:LEU:HD12	1:A:366:LEU:C	2.16	0.68
1:A:280:LYS:NZ	1:A:287:SER:O	2.27	0.67
1:B:452:ALA:HB3	1:B:470:SER:HB2	1.77	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:204:GLN:O	1:C:230:ASN:ND2	2.25	0.67
1:C:302:LEU:C	1:C:302:LEU:HD13	2.20	0.67
1:B:65:ILE:HD12	1:B:65:ILE:H	1.58	0.67
1:C:334:LEU:HD23	1:C:334:LEU:C	2.19	0.66
1:C:130:SER:OG	1:C:510:ARG:NH2	2.27	0.66
1:B:146:LYS:HZ1	1:B:173:GLU:HG3	1.58	0.66
1:B:153:THR:OG1	1:B:491:GLU:HB2	1.95	0.66
1:B:183:ARG:HH11	1:B:183:ARG:CG	2.09	0.66
1:C:426:PHE:CE1	1:C:447:MET:HG3	2.31	0.66
1:C:306:ARG:HH11	1:C:306:ARG:CG	2.09	0.65
1:A:198:PRO:HA	1:A:201:ARG:HE	1.61	0.65
1:B:234:GLY:HA2	1:B:262:ALA:O	1.97	0.65
1:B:146:LYS:HZ1	1:B:173:GLU:HG2	1.60	0.64
1:A:38:GLU:OE2	1:A:130:SER:N	2.30	0.64
1:A:221:GLY:HA2	1:A:248:ASN:HD21	1.62	0.64
1:B:500:GLU:O	1:B:500:GLU:HG2	1.95	0.64
1:A:204:GLN:O	1:A:230:ASN:ND2	2.27	0.64
1:A:296:LEU:HD23	1:A:505:LEU:HD13	1.78	0.64
1:B:146:LYS:NZ	1:B:173:GLU:HG3	2.13	0.64
1:B:426:PHE:CG	1:B:447:MET:HE3	2.34	0.63
1:C:372:VAL:CG2	1:C:399:LEU:HD22	2.20	0.63
1:B:398:THR:HG21	1:B:505:LEU:HB2	1.80	0.63
1:B:102:ASP:HB3	1:B:105:LEU:HG	1.81	0.63
1:B:322:SER:HG	1:B:324:SER:HG	1.45	0.63
1:B:339:GLN:HE21	1:B:340:SER:H	1.47	0.63
1:A:463:VAL:HG23	1:A:491:GLU:HG2	1.81	0.62
1:B:147:TRP:CD2	1:B:171:VAL:HG23	2.34	0.62
1:B:403:LYS:HD2	1:B:436:GLU:HG2	1.80	0.62
1:B:431:LEU:HD23	1:B:433:PHE:HE2	1.65	0.62
1:C:426:PHE:CD1	1:C:447:MET:CG	2.83	0.62
1:A:348:LEU:H	1:A:348:LEU:CD2	2.05	0.62
1:C:248:ASN:ND2	1:C:256:LYS:HD2	2.15	0.61
1:C:242:GLN:HG2	1:C:243:ASP:OD1	1.99	0.61
1:C:280:LYS:NZ	1:C:287:SER:O	2.34	0.61
1:A:160:VAL:HG11	1:A:171:VAL:CG1	2.30	0.60
1:C:146:LYS:HG3	1:C:149:ASN:HD21	1.65	0.60
1:C:457:ARG:NH1	1:C:482:MET:O	2.34	0.60
1:A:502:ARG:NH2	1:A:507:PRO:O	2.34	0.60
1:B:416:ASN:HD21	1:B:444:GLN:N	1.99	0.60
1:B:9:ASP:N	1:B:9:ASP:OD1	2.34	0.60
1:B:16:PHE:CE1	1:B:523:ARG:HG2	2.36	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:101:GLN:HG2	1:B:102:ASP:N	2.16	0.60
1:C:388:ARG:HB3	1:C:423:PHE:CD2	2.36	0.60
1:B:416:ASN:ND2	1:B:444:GLN:H	1.99	0.59
1:C:434:VAL:HG13	1:C:451:TYR:CD2	2.37	0.59
1:B:426:PHE:HB2	1:B:447:MET:CE	2.32	0.59
1:C:418:LEU:HD22	1:C:418:LEU:N	2.17	0.59
1:C:452:ALA:HB3	1:C:470:SER:HB2	1.83	0.59
1:C:329:MET:HB3	1:C:377:ILE:HD13	1.84	0.59
1:C:426:PHE:CD1	1:C:447:MET:HG3	2.37	0.59
1:A:335:THR:HB	1:A:380:VAL:CG2	2.31	0.59
1:A:36:TRP:HB2	1:A:61:ARG:NH2	2.17	0.59
1:B:431:LEU:CD2	1:B:433:PHE:CZ	2.86	0.59
1:C:388:ARG:NE	1:C:423:PHE:HD2	2.01	0.59
1:C:172:ILE:HG22	1:C:172:ILE:O	2.02	0.58
1:A:227:GLU:O	1:A:231:LYS:HB3	2.02	0.58
1:C:418:LEU:HB3	1:C:419:PRO:HD2	1.85	0.58
1:C:388:ARG:HE	1:C:423:PHE:HD2	1.51	0.57
1:A:158:VAL:HG13	1:A:180:LEU:HD11	1.85	0.57
1:A:168:ASP:OD1	1:A:239:ARG:NH2	2.29	0.57
1:A:204:GLN:N	1:A:204:GLN:OE1	2.38	0.57
1:C:401:VAL:O	1:C:401:VAL:HG23	2.03	0.57
1:B:447:MET:CG	1:B:465:PHE:HB2	2.35	0.57
1:A:158:VAL:HG22	1:A:210:LEU:HD13	1.87	0.56
1:C:96:LEU:HD12	1:C:107:THR:O	2.05	0.56
1:C:150:GLY:HA2	1:C:486:ALA:HB3	1.87	0.56
1:C:338:THR:HG23	1:C:338:THR:O	2.04	0.56
1:C:497:ILE:HD13	1:C:502:ARG:NH1	2.19	0.56
1:B:133:LEU:CD1	1:B:159:TYR:CE1	2.89	0.56
1:C:327:GLN:OE1	1:C:327:GLN:HA	2.06	0.56
1:A:369:PHE:CD1	1:A:369:PHE:N	2.73	0.56
1:B:241:ALA:HA	1:B:244:ILE:HG22	1.87	0.56
1:B:153:THR:HG22	1:B:489:LYS:HD2	1.87	0.56
1:B:118:LYS:HG2	1:B:119:ALA:N	2.20	0.55
1:C:111:GLU:HG3	1:C:120:ARG:HG3	1.88	0.55
1:B:447:MET:HG3	1:B:465:PHE:HD1	1.72	0.55
1:B:60:ASP:HA	1:B:64:THR:HG23	1.87	0.55
1:B:302:LEU:HD21	1:B:304:ILE:HD13	1.89	0.55
1:C:177:ASN:O	1:C:489:LYS:HE3	2.07	0.55
1:A:185:ASP:OD2	1:A:187:THR:OG1	2.18	0.55
1:A:220:GLY:O	1:A:248:ASN:ND2	2.40	0.55
1:A:417:LEU:HB3	1:A:446:VAL:HG23	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:431:LEU:HD23	1:B:433:PHE:CZ	2.41	0.55
1:A:145:ASN:O	1:A:146:LYS:HD3	2.07	0.55
1:B:130:SER:HB3	1:B:510:ARG:HD3	1.89	0.55
1:B:133:LEU:CD1	1:B:159:TYR:HE1	2.19	0.55
1:B:414:ASN:HD21	1:B:442:ARG:NH1	2.03	0.55
1:C:34:TYR:HD2	1:C:129:ASN:HD21	1.53	0.55
1:C:155:ARG:NH1	1:C:183:ARG:O	2.37	0.55
1:C:418:LEU:HD12	1:C:447:MET:HE3	1.88	0.55
1:B:417:LEU:O	1:B:446:VAL:HA	2.05	0.55
1:A:446:VAL:O	1:A:464:LEU:HD12	2.05	0.54
1:B:379:ALA:HB3	1:B:411:LEU:HD12	1.89	0.54
1:C:171:VAL:O	1:C:171:VAL:HG22	2.06	0.54
1:C:217:ILE:HG13	1:C:217:ILE:O	2.07	0.54
1:C:60:ASP:OD1	1:C:60:ASP:N	2.41	0.54
1:C:126:ARG:HB2	1:C:519:LEU:HD11	1.88	0.54
1:B:38:GLU:HG2	1:B:129:ASN:OD1	2.07	0.54
1:C:248:ASN:HD21	1:C:256:LYS:HD2	1.73	0.54
1:B:184:TYR:CD1	1:B:184:TYR:C	2.85	0.54
1:B:241:ALA:HA	1:B:244:ILE:CG2	2.37	0.54
1:B:419:PRO:CD	1:B:447:MET:HE2	2.37	0.54
1:C:375:GLU:HG2	1:C:402:LEU:HD12	1.90	0.54
1:A:32:GLN:HG2	1:A:61:ARG:NH2	2.23	0.54
1:C:329:MET:HB3	1:C:377:ILE:CD1	2.37	0.54
1:B:450:VAL:CG2	1:B:468:VAL:HG23	2.38	0.54
1:B:502:ARG:HG3	1:B:503:PRO:HD2	1.89	0.54
1:A:142:GLY:HA2	1:A:472:GLN:HG3	1.89	0.54
1:B:426:PHE:HB2	1:B:447:MET:HE1	1.89	0.54
1:C:216:LYS:HE3	1:C:243:ASP:HA	1.90	0.54
1:C:247:GLU:OE2	1:C:247:GLU:N	2.32	0.54
1:C:183:ARG:CB	1:C:204:GLN:HA	2.37	0.54
1:C:291:THR:HG23	1:C:294:ALA:H	1.72	0.54
1:B:427:PRO:HG2	1:B:498:PRO:HD3	1.90	0.53
1:A:249:VAL:HA	1:A:258:VAL:HG22	1.89	0.53
1:C:145:ASN:C	1:C:146:LYS:HD3	2.32	0.53
1:C:219:VAL:HG13	1:C:219:VAL:O	2.09	0.53
1:C:306:ARG:HG2	1:C:306:ARG:NH1	2.13	0.53
1:A:456:PHE:C	1:A:456:PHE:CD2	2.85	0.53
1:B:160:VAL:HB	1:B:212:VAL:HG22	1.90	0.53
1:B:235:VAL:HG23	1:B:235:VAL:O	2.07	0.53
1:A:306:ARG:CZ	1:A:332:GLY:CA	2.87	0.53
1:C:359:TYR:C	1:C:359:TYR:CD2	2.86	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:184:TYR:CD1	1:A:206:LEU:HD21	2.44	0.53
1:B:153:THR:OG1	1:B:153:THR:O	2.27	0.53
1:C:395:LYS:HB3	1:C:429:HIS:HD2	1.74	0.53
1:B:419:PRO:HD3	1:B:447:MET:HE2	1.91	0.53
1:A:444:GLN:HA	1:A:462:ASN:ND2	2.23	0.52
1:B:227:GLU:N	1:B:230:ASN:O	2.42	0.52
1:C:313:VAL:HG12	1:C:313:VAL:O	2.09	0.52
1:B:183:ARG:CG	1:B:183:ARG:NH1	2.72	0.52
1:B:39:GLN:O	1:B:39:GLN:HG2	2.10	0.52
1:A:347:ARG:HH11	1:A:347:ARG:CG	2.18	0.52
1:C:446:VAL:HG22	1:C:464:LEU:HD13	1.88	0.52
1:B:303:ARG:C	1:B:304:ILE:HD12	2.35	0.52
1:C:95:THR:HG21	1:C:97:TYR:CE2	2.46	0.51
1:C:372:VAL:CG2	1:C:399:LEU:CD2	2.84	0.51
1:B:171:VAL:HG13	1:B:172:ILE:HD12	1.93	0.51
1:B:431:LEU:HD21	1:B:433:PHE:CZ	2.45	0.51
1:A:306:ARG:NH2	1:A:332:GLY:HA3	2.25	0.51
1:C:135:GLN:HB2	1:C:157:GLY:O	2.10	0.51
1:C:163:ASN:HB3	1:C:166:ASP:HB3	1.92	0.51
1:A:225:ILE:HD12	1:A:258:VAL:HG12	1.93	0.51
1:B:153:THR:HG22	1:B:489:LYS:CD	2.39	0.51
1:B:161:VAL:HG13	1:B:213:GLN:HE21	1.76	0.51
1:A:169:GLN:OE1	1:A:169:GLN:HA	2.11	0.51
1:C:213:GLN:HG2	1:C:214:TYR:CD1	2.46	0.51
1:A:306:ARG:CZ	1:A:332:GLY:HA3	2.41	0.51
1:B:431:LEU:CD2	1:B:433:PHE:CE2	2.87	0.51
1:C:447:MET:O	1:C:447:MET:HG2	2.10	0.51
1:C:446:VAL:HG23	1:C:464:LEU:CD1	2.39	0.51
1:A:316:PRO:HG3	1:A:345:PHE:CD1	2.46	0.50
1:B:135:GLN:NE2	1:B:138:PHE:HD2	2.10	0.50
1:B:236:PHE:HB3	1:B:265:GLY:O	2.11	0.50
1:B:447:MET:HG3	1:B:465:PHE:CD1	2.46	0.50
1:C:336:LEU:HD22	1:C:357:TYR:CD2	2.47	0.50
1:C:417:LEU:HB3	1:C:446:VAL:HG12	1.93	0.50
1:A:460:LYS:HG3	1:A:476:THR:O	2.12	0.50
1:C:192:VAL:HG12	1:C:192:VAL:O	2.09	0.50
1:B:413:ILE:HG21	1:B:446:VAL:HG11	1.93	0.50
1:A:303:ARG:HB3	1:A:315:SER:HB2	1.93	0.50
1:A:439:ILE:HD12	1:A:450:VAL:HG13	1.94	0.50
1:B:416:ASN:ND2	1:B:445:HIS:H	2.10	0.50
1:A:10:THR:O	1:A:14:GLN:HG3	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:443:GLY:C	1:A:462:ASN:HD22	2.20	0.50
1:A:185:ASP:OD1	1:A:202:GLN:HB3	2.12	0.50
1:B:316:PRO:HG3	1:B:345:PHE:CG	2.46	0.49
1:A:60:ASP:OD1	1:A:62:ASP:HB3	2.13	0.49
1:B:509:LEU:HD12	1:B:509:LEU:H	1.77	0.49
1:C:302:LEU:C	1:C:302:LEU:CD1	2.85	0.49
1:A:196:VAL:HG23	1:A:201:ARG:HD3	1.94	0.49
1:C:153:THR:HG22	1:C:489:LYS:NZ	2.27	0.49
1:A:184:TYR:CE1	1:A:206:LEU:HD21	2.47	0.49
1:C:248:ASN:CG	1:C:256:LYS:HD2	2.37	0.49
1:A:92:TYR:CD1	1:A:92:TYR:C	2.91	0.49
1:C:306:ARG:CG	1:C:306:ARG:NH1	2.72	0.49
1:A:463:VAL:HG13	1:A:465:PHE:CE2	2.48	0.49
1:B:450:VAL:HG23	1:B:468:VAL:HG23	1.94	0.49
1:C:165:ASN:OD1	1:C:165:ASN:N	2.46	0.49
1:B:227:GLU:O	1:B:231:LYS:HB3	2.13	0.48
1:C:317:PRO:C	1:C:318:ASN:HD22	2.20	0.48
1:A:160:VAL:HG11	1:A:171:VAL:HG11	1.95	0.48
1:C:497:ILE:N	1:C:497:ILE:CD1	2.73	0.48
1:A:158:VAL:HG13	1:A:180:LEU:CD1	2.43	0.48
1:B:388:ARG:NH1	1:B:420:ASP:OD2	2.46	0.48
1:C:387:TYR:CD1	1:C:387:TYR:C	2.91	0.48
1:B:198:PRO:HA	1:B:201:ARG:HG3	1.95	0.48
1:C:1:ILE:HD12	1:C:2:GLU:H	1.78	0.48
1:C:444:GLN:C	1:C:445:HIS:HD2	2.21	0.48
1:A:359:TYR:CD1	1:A:360:THR:N	2.82	0.48
1:C:426:PHE:CD2	1:C:427:PRO:HA	2.49	0.48
1:A:169:GLN:O	1:A:171:VAL:HG23	2.13	0.48
1:C:130:SER:HA	1:C:512:GLY:O	2.13	0.48
1:C:426:PHE:CD1	1:C:447:MET:SD	3.07	0.48
1:B:133:LEU:HD11	1:B:159:TYR:HE1	1.78	0.48
1:C:213:GLN:HG2	1:C:214:TYR:CE1	2.49	0.48
1:B:416:ASN:ND2	1:B:444:GLN:N	2.61	0.47
1:B:426:PHE:CD1	1:B:427:PRO:HA	2.49	0.47
1:B:460:LYS:HD2	1:B:488:GLN:HB3	1.95	0.47
1:C:237:VAL:CG2	1:C:244:ILE:HG13	2.44	0.47
1:A:198:PRO:CA	1:A:201:ARG:HE	2.26	0.47
1:A:231:LYS:HD2	1:A:259:CYS:C	2.40	0.47
1:B:135:GLN:HE22	1:B:138:PHE:HD2	1.61	0.47
1:B:304:ILE:HD12	1:B:304:ILE:N	2.29	0.47
1:B:237:VAL:O	1:B:237:VAL:HG12	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:34:TYR:CD2	1:C:129:ASN:ND2	2.82	0.47
1:A:444:GLN:HA	1:A:462:ASN:HD22	1.80	0.47
1:B:359:TYR:HE2	1:B:383:ARG:HD3	1.80	0.47
1:C:37:ARG:CB	1:C:129:ASN:OD1	2.62	0.47
1:C:383:ARG:CZ	1:C:383:ARG:HB3	2.45	0.47
1:A:236:PHE:HB3	1:A:265:GLY:O	2.15	0.47
1:B:37:ARG:HH11	1:B:37:ARG:CG	2.28	0.47
1:B:336:LEU:HB3	1:B:357:TYR:CE2	2.50	0.47
1:C:219:VAL:HG13	1:C:248:ASN:HB2	1.97	0.47
1:A:170:TYR:HB3	1:A:173:GLU:HG2	1.96	0.47
1:C:153:THR:HG22	1:C:489:LYS:HZ3	1.78	0.47
1:C:228:PRO:HG2	1:C:255:ASN:HD21	1.80	0.46
1:B:502:ARG:HG3	1:B:503:PRO:CD	2.46	0.46
1:C:64:THR:O	1:C:64:THR:HG23	2.14	0.46
1:C:297:GLN:HB3	1:C:303:ARG:CZ	2.45	0.46
1:B:234:GLY:HA3	1:B:236:PHE:CZ	2.51	0.46
1:A:254:ASN:O	1:A:255:ASN:HB3	2.15	0.46
1:B:53:LEU:HD12	1:B:53:LEU:HA	1.47	0.46
1:C:434:VAL:HG13	1:C:451:TYR:HD2	1.79	0.46
1:A:209:SER:HA	1:A:232:VAL:HG22	1.97	0.46
1:C:146:LYS:HD3	1:C:146:LYS:N	2.30	0.46
1:B:180:LEU:C	1:B:180:LEU:CD1	2.85	0.46
1:C:121:VAL:HG12	1:C:523:ARG:HB3	1.98	0.46
1:B:133:LEU:HD13	1:B:159:TYR:CD1	2.51	0.45
1:A:142:GLY:HA2	1:A:472:GLN:CG	2.47	0.45
1:C:121:VAL:HG12	1:C:523:ARG:CB	2.47	0.45
1:A:31:PHE:O	1:A:35:VAL:HG13	2.16	0.45
1:A:455:THR:HG22	1:A:472:GLN:HB2	1.98	0.45
1:C:53:LEU:HD12	1:C:53:LEU:HA	1.63	0.45
1:B:184:TYR:CD1	1:B:184:TYR:O	2.70	0.45
1:B:339:GLN:HE21	1:B:340:SER:N	2.13	0.45
1:A:233:LYS:HE3	1:A:233:LYS:HB3	1.59	0.45
1:A:446:VAL:HG13	1:A:464:LEU:CD1	2.44	0.45
1:B:329:MET:HE1	1:B:374:LEU:HD13	1.98	0.45
1:C:426:PHE:HE2	1:C:430:ALA:HB1	1.81	0.45
1:B:302:LEU:CD2	1:B:304:ILE:HD13	2.46	0.45
1:B:420:ASP:OD1	1:B:422:THR:HG22	2.17	0.45
1:B:502:ARG:HA	1:B:503:PRO:HD3	1.78	0.45
1:C:508:SER:O	1:C:508:SER:OG	2.33	0.45
1:B:431:LEU:O	1:B:449:PRO:HD2	2.17	0.45
1:B:457:ARG:HG2	1:B:474:CYS:SG	2.57	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:136:ALA:HB2	1:A:494:TYR:HB2	1.99	0.45
1:C:357:TYR:O	1:C:357:TYR:CG	2.69	0.45
1:C:463:VAL:HA	1:C:491:GLU:O	2.17	0.45
1:B:133:LEU:HD13	1:B:159:TYR:CE1	2.51	0.44
1:B:231:LYS:HB2	1:B:259:CYS:HB3	1.99	0.44
1:B:426:PHE:HB2	1:B:447:MET:HE3	1.97	0.44
1:B:466:GLY:O	1:B:494:TYR:HA	2.17	0.44
1:C:329:MET:HG2	1:C:377:ILE:HD13	2.00	0.44
1:A:36:TRP:HB2	1:A:61:ARG:HH21	1.82	0.44
1:A:456:PHE:CD2	1:A:456:PHE:O	2.70	0.44
1:C:497:ILE:HD13	1:C:502:ARG:HH11	1.82	0.44
1:A:121:VAL:HG12	1:A:523:ARG:CB	2.48	0.44
1:B:303:ARG:HA	1:B:373:VAL:O	2.18	0.44
1:B:505:LEU:N	1:B:505:LEU:HD22	2.33	0.44
1:C:426:PHE:HA	1:C:427:PRO:HA	1.73	0.44
1:A:146:LYS:HB3	1:A:149:ASN:HD21	1.83	0.44
1:A:316:PRO:HG3	1:A:345:PHE:CG	2.52	0.44
1:A:459:VAL:HG23	1:A:462:ASN:CG	2.42	0.44
1:B:153:THR:HG23	1:B:489:LYS:HD3	1.99	0.44
1:B:227:GLU:C	1:B:230:ASN:O	2.60	0.44
1:C:135:GLN:N	1:C:135:GLN:CD	2.76	0.44
1:A:460:LYS:CE	1:A:488:GLN:HB2	2.47	0.44
1:A:53:LEU:HA	1:A:53:LEU:HD12	1.74	0.44
1:C:9:ASP:CG	1:C:10:THR:N	2.76	0.44
1:C:237:VAL:HG22	1:C:244:ILE:HG13	1.99	0.44
1:A:196:VAL:CG2	1:A:201:ARG:HD3	2.48	0.44
1:A:222:SER:H	1:A:256:LYS:HG3	1.83	0.43
1:B:497:ILE:H	1:B:497:ILE:CD1	2.12	0.43
1:A:182:ASN:O	1:A:182:ASN:ND2	2.51	0.43
1:B:233:LYS:O	1:B:233:LYS:HG2	2.18	0.43
1:C:440:TYR:CE2	1:C:442:ARG:HG2	2.54	0.43
1:A:231:LYS:CD	1:A:259:CYS:HB3	2.44	0.43
1:A:342:ASP:CG	1:A:356:ARG:HD3	2.44	0.43
1:B:197:GLU:HG2	1:B:425:LEU:HB2	2.00	0.43
1:B:335:THR:HA	1:B:380:VAL:HG22	2.01	0.43
1:C:212:VAL:HB	1:C:237:VAL:HB	2.01	0.43
1:A:201:ARG:HB2	1:A:202:GLN:OE1	2.18	0.43
1:A:344:THR:HA	1:A:353:GLY:O	2.18	0.43
1:B:122:GLN:HG2	1:B:522:GLU:HG2	2.00	0.43
1:C:71:ASN:OD1	1:C:100:ALA:HB2	2.18	0.43
1:C:277:LEU:O	1:C:291:THR:OG1	2.36	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:366:LEU:HD12	1:A:367:GLU:N	2.34	0.43
1:A:459:VAL:HG23	1:A:462:ASN:OD1	2.19	0.43
1:C:147:TRP:HB2	1:C:171:VAL:HG23	2.01	0.43
1:C:248:ASN:OD1	1:C:256:LYS:HD2	2.19	0.43
1:C:184:TYR:CD1	1:C:184:TYR:C	2.96	0.43
1:A:137:ILE:HG12	1:A:468:VAL:HG22	2.00	0.43
1:A:184:TYR:CD2	1:A:184:TYR:C	2.97	0.43
1:A:224:GLN:NE2	1:A:227:GLU:HG2	2.34	0.43
1:A:482:MET:HB3	1:A:482:MET:HE3	1.41	0.43
1:A:364:GLU:N	1:A:364:GLU:CD	2.76	0.42
1:C:133:LEU:N	1:C:133:LEU:HD23	2.32	0.42
1:A:134:GLU:HB3	1:A:494:TYR:CE1	2.54	0.42
1:B:103:ASP:OD1	1:B:103:ASP:N	2.50	0.42
1:B:521:TYR:CD1	1:B:521:TYR:C	2.97	0.42
1:C:309:ASN:OD1	1:C:309:ASN:N	2.51	0.42
1:B:182:ASN:HD21	1:B:206:LEU:H	1.68	0.42
1:B:426:PHE:CB	1:B:447:MET:HE3	2.50	0.42
1:C:154:ILE:O	1:C:180:LEU:HA	2.20	0.42
1:C:146:LYS:CG	1:C:149:ASN:HD21	2.32	0.42
1:C:427:PRO:C	1:C:429:HIS:H	2.27	0.42
1:C:306:ARG:HD2	1:C:311:LEU:HD11	2.02	0.42
1:A:306:ARG:O	1:A:306:ARG:HG2	2.20	0.42
1:A:460:LYS:HE3	1:A:477:SER:HA	2.01	0.42
1:C:349:ASP:OD1	1:C:351:SER:HB3	2.19	0.42
1:A:526:GLU:O	1:A:527:HIS:ND1	2.53	0.41
1:B:148:LEU:C	1:B:148:LEU:HD13	2.45	0.41
1:C:297:GLN:HA	1:C:303:ARG:HD2	2.02	0.41
1:C:440:TYR:CD1	1:C:457:ARG:HG3	2.55	0.41
1:A:133:LEU:HB3	1:A:135:GLN:OE1	2.20	0.41
1:B:19:ALA:HB1	1:B:121:VAL:HG22	2.02	0.41
1:C:296:LEU:CD2	1:C:505:LEU:HD23	2.35	0.41
1:A:387:TYR:CD1	1:A:387:TYR:C	2.98	0.41
1:B:474:CYS:HB2	1:B:484:CYS:HB3	1.93	0.41
1:A:25:LYS:HD3	1:A:26:TYR:CE1	2.55	0.41
1:B:92:TYR:HA	1:B:111:GLU:O	2.21	0.41
1:B:128:SER:O	1:B:514:PRO:HA	2.21	0.41
1:B:329:MET:HE3	1:B:329:MET:HB3	1.69	0.41
1:B:380:VAL:HG12	1:B:412:ASP:HB2	2.02	0.41
1:A:382:ASN:O	1:A:382:ASN:ND2	2.47	0.41
1:C:20:GLU:OE1	1:C:521:TYR:OH	2.32	0.41
1:A:213:GLN:HA	1:A:266:PHE:CD1	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:373:VAL:HG22	1:B:400:ALA:HB3	2.03	0.41
1:C:179:ALA:HB3	1:C:489:LYS:HZ3	1.85	0.41
1:C:325:CYS:HB2	1:C:343:CYS:HA	2.03	0.41
1:C:459:VAL:HG22	1:C:462:ASN:ND2	2.36	0.41
1:C:499:LYS:HA	1:C:499:LYS:HD2	1.74	0.41
1:A:360:THR:HG21	1:A:363:GLN:O	2.21	0.41
1:B:135:GLN:HE21	1:B:159:TYR:HD1	1.68	0.41
1:B:148:LEU:HD22	1:B:149:ASN:H	1.86	0.41
1:C:1:ILE:HD12	1:C:1:ILE:HA	1.75	0.41
1:C:150:GLY:CA	1:C:486:ALA:HB3	2.50	0.41
1:C:175:ASN:C	1:C:221:GLY:HA3	2.45	0.41
1:C:197:GLU:HG3	1:C:198:PRO:CD	2.47	0.41
1:C:447:MET:HE2	1:C:447:MET:HB2	1.82	0.41
1:C:497:ILE:H	1:C:497:ILE:CD1	2.22	0.41
1:B:59:LEU:H	1:B:59:LEU:HD12	1.86	0.40
1:C:111:GLU:CG	1:C:120:ARG:HG3	2.51	0.40
1:C:271:PRO:O	1:C:271:PRO:HG2	2.21	0.40
1:C:457:ARG:HA	1:C:474:CYS:O	2.21	0.40
1:A:109:VAL:HG22	1:A:122:GLN:HB2	2.03	0.40
1:C:451:TYR:HA	1:C:469:ILE:O	2.20	0.40
1:A:327:GLN:OE1	1:A:330:GLN:NE2	2.55	0.40
1:C:122:GLN:HG2	1:C:522:GLU:HG2	2.03	0.40
1:C:219:VAL:O	1:C:219:VAL:CG1	2.69	0.40
1:C:360:THR:O	1:C:360:THR:CG2	2.69	0.40
1:C:426:PHE:CE2	1:C:430:ALA:CB	2.86	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	512/534 (96%)	478 (93%)	34 (7%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	476/534 (89%)	441 (93%)	35 (7%)	0	100	100
1	C	491/534 (92%)	460 (94%)	31 (6%)	0	100	100
All	All	1479/1602 (92%)	1379 (93%)	100 (7%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	415/424 (98%)	392 (94%)	23 (6%)	19	43
1	B	396/424 (93%)	361 (91%)	35 (9%)	9	22
1	C	399/424 (94%)	351 (88%)	48 (12%)	5	12
All	All	1210/1272 (95%)	1104 (91%)	106 (9%)	9	22

All (106) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	66	THR
1	A	77	GLN
1	A	80	VAL
1	A	110	SER
1	A	126	ARG
1	A	153	THR
1	A	158	VAL
1	A	161	VAL
1	A	299	LYS
1	A	302	LEU
1	A	311	LEU
1	A	335	THR
1	A	348	LEU
1	A	349	ASP
1	A	351	SER
1	A	366	LEU

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Mol	Chain	Res	Type
1	A	369	PHE
1	A	372	VAL
1	A	373	VAL
1	A	385	VAL
1	A	420	ASP
1	A	422	THR
1	A	482	MET
1	B	10	THR
1	B	37	ARG
1	B	39	GLN
1	B	49	CYS
1	B	51	THR
1	B	64	THR
1	B	65	ILE
1	B	66	THR
1	B	149	ASN
1	B	154	ILE
1	B	158	VAL
1	B	173	GLU
1	B	177	ASN
1	B	180	LEU
1	B	182	ASN
1	B	183	ARG
1	B	204	GLN
1	B	205	ASP
1	B	224	GLN
1	B	311	LEU
1	B	329	MET
1	B	331	SER
1	B	396	SER
1	B	439	ILE
1	B	459	VAL
1	B	472	GLN
1	B	474	CYS
1	B	475	THR
1	B	477	SER
1	B	480	ASN
1	B	485	ASN
1	B	496	ARG
1	B	497	ILE
1	B	499	LYS
1	B	500	GLU

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Mol	Chain	Res	Type
1	C	1	ILE
1	C	2	GLU
1	C	16	PHE
1	C	35	VAL
1	C	59	LEU
1	C	60	ASP
1	C	101	GLN
1	C	118	LYS
1	C	165	ASN
1	C	169	GLN
1	C	194	ASN
1	C	202	GLN
1	C	255	ASN
1	C	256	LYS
1	C	269	SER
1	C	270	ASP
1	C	302	LEU
1	C	304	ILE
1	C	306	ARG
1	C	307	ILE
1	C	309	ASN
1	C	310	ILE
1	C	311	LEU
1	C	331	SER
1	C	334	LEU
1	C	358	THR
1	C	372	VAL
1	C	373	VAL
1	C	382	ASN
1	C	385	VAL
1	C	393	SER
1	C	398	THR
1	C	399	LEU
1	C	402	LEU
1	C	404	LEU
1	C	418	LEU
1	C	425	LEU
1	C	426	PHE
1	C	431	LEU
1	C	447	MET
1	C	450	VAL
1	C	483	SER

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Mol	Chain	Res	Type
1	C	495	ILE
1	C	496	ARG
1	C	497	ILE
1	C	502	ARG
1	C	506	LEU
1	C	527	HIS

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

Mol	Chain	Res	Type
1	A	71	ASN
1	A	149	ASN
1	A	248	ASN
1	A	254	ASN
1	A	330	GLN
1	A	444	GLN
1	B	14	GLN
1	B	70	ASN
1	B	104	GLN
1	B	149	ASN
1	B	165	ASN
1	B	175	ASN
1	B	248	ASN
1	B	318	ASN
1	B	327	GLN
1	B	339	GLN
1	B	407	ASN
1	B	414	ASN
1	B	416	ASN
1	B	445	HIS
1	B	488	GLN
1	C	14	GLN
1	C	70	ASN
1	C	85	ASN
1	C	129	ASN
1	C	149	ASN
1	C	177	ASN
1	C	204	GLN
1	C	255	ASN
1	C	305	GLN
1	C	318	ASN
1	C	390	GLN

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Mol	Chain	Res	Type
1	C	414	ASN
1	C	416	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 [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	518/534 (97%)	-0.12	2 (0%) 88 88	45, 82, 133, 186	0
1	B	498/534 (93%)	0.01	9 (1%) 67 65	53, 95, 163, 197	0
1	C	507/534 (94%)	0.11	10 (1%) 65 62	67, 102, 145, 177	0
All	All	1523/1602 (95%)	-0.00	21 (1%) 73 71	45, 94, 148, 197	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	364	GLU	5.2
1	B	186	LEU	4.5
1	C	501	ASN	3.5
1	B	142	GLY	3.4
1	C	100	ALA	2.8
1	B	178	PHE	2.6
1	B	269	SER	2.6
1	B	237	VAL	2.5
1	C	217	ILE	2.4
1	A	486	ALA	2.4
1	C	241	ALA	2.4
1	B	251	VAL	2.3
1	C	348	LEU	2.3
1	C	307	ILE	2.2
1	A	348	LEU	2.2
1	B	244	ILE	2.2
1	C	38	GLU	2.2
1	B	466	GLY	2.1
1	B	208	ALA	2.0
1	C	141	ALA	2.0
1	C	309	ASN	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.