



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:03 PM UTC

PDB ID : 3TQI / pdb_00003tqi
Title : Structure of the GMP synthase (guaA) from Coxiella burnetii
Authors : Franklin, M.C.; Cheung, J.; Rudolph, M.; Cassidy, M.; Gary, E.; Burshteyn, F.; Love, J.
Deposited on : 2011-09-09
Resolution : 2.84 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

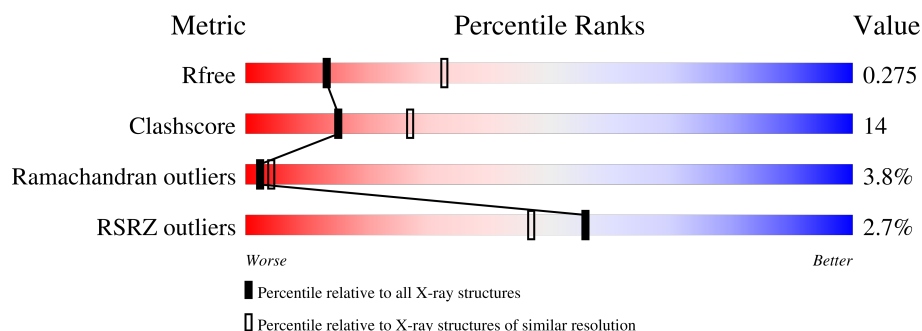
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.84 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	1520 (2.86-2.82)
Clashscore	190562	1559 (2.86-2.82)
Ramachandran outliers	187476	1517 (2.86-2.82)
RSRZ outliers	180081	1521 (2.86-2.82)

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	527	2% 67% 18% • 12%
1	B	527	2% 67% 19% • 12%
1	C	527	3% 65% 21% • 12%
1	D	527	2% 62% 23% • 12%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 14708 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 GMP synthase [glutamine-hydrolyzing].

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	464	Total	C	N	O	S	0	0	0
			3677	2371	626	663	17			
1	B	464	Total	C	N	O	S	0	0	0
			3677	2371	626	663	17			
1	C	464	Total	C	N	O	S	0	0	0
			3677	2371	626	663	17			
1	D	464	Total	C	N	O	S	0	0	0
			3677	2371	626	663	17			

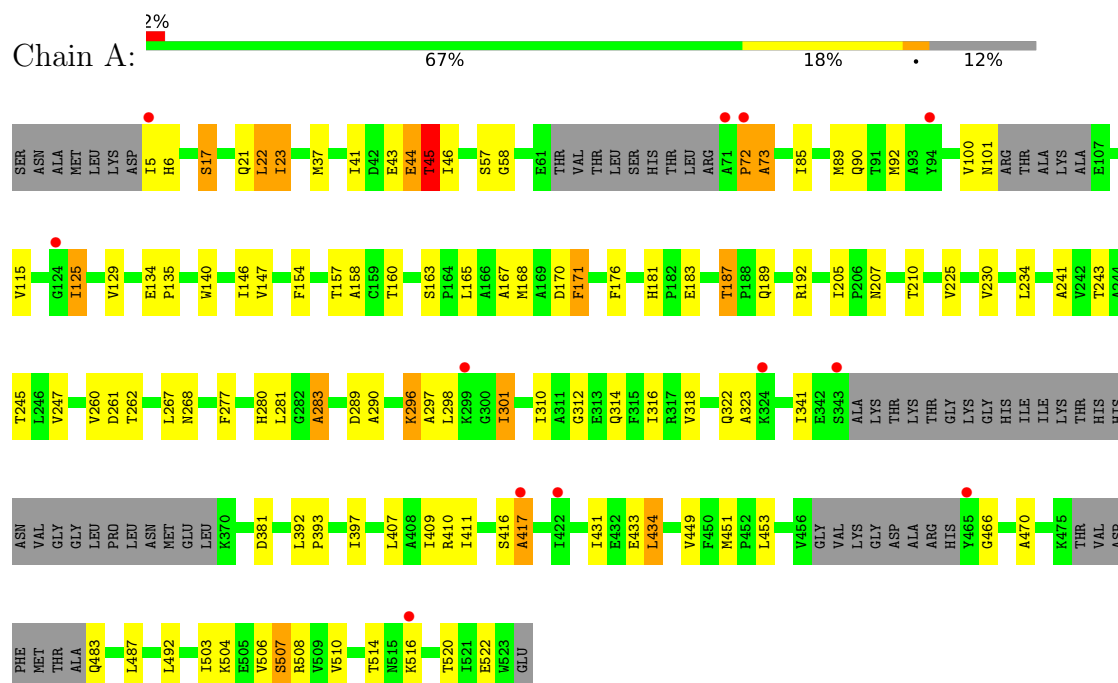
There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	SER	-	expression tag	UNP Q83BZ6
C	-1	ASN	-	expression tag	UNP Q83BZ6
C	0	ALA	-	expression tag	UNP Q83BZ6
A	-2	SER	-	expression tag	UNP Q83BZ6
A	-1	ASN	-	expression tag	UNP Q83BZ6
A	0	ALA	-	expression tag	UNP Q83BZ6
D	-2	SER	-	expression tag	UNP Q83BZ6
D	-1	ASN	-	expression tag	UNP Q83BZ6
D	0	ALA	-	expression tag	UNP Q83BZ6
B	-2	SER	-	expression tag	UNP Q83BZ6
B	-1	ASN	-	expression tag	UNP Q83BZ6
B	0	ALA	-	expression tag	UNP Q83BZ6

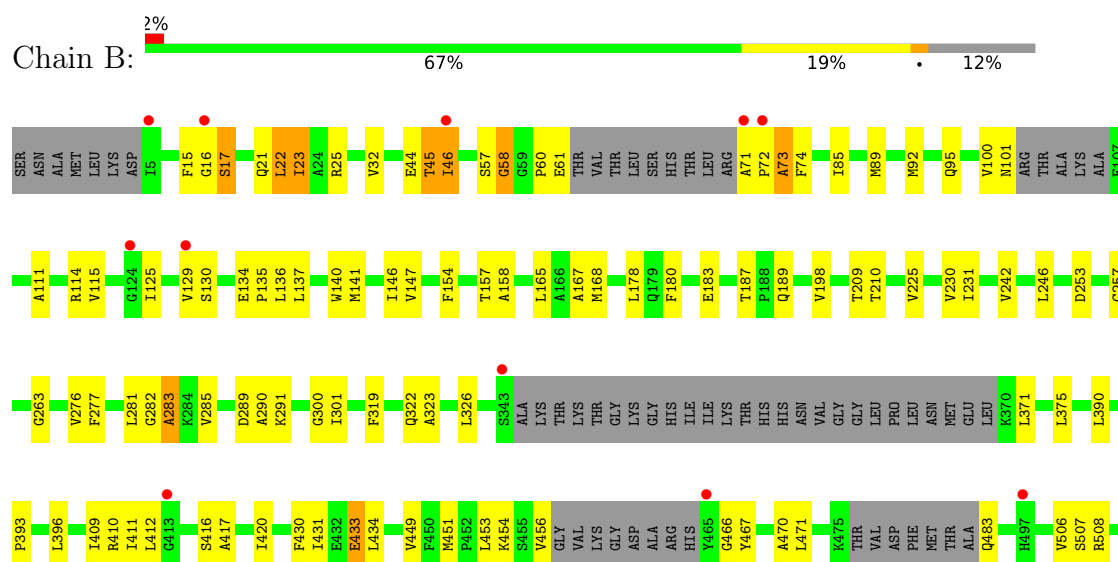
3 Residue-property plots

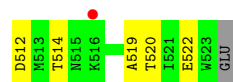
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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

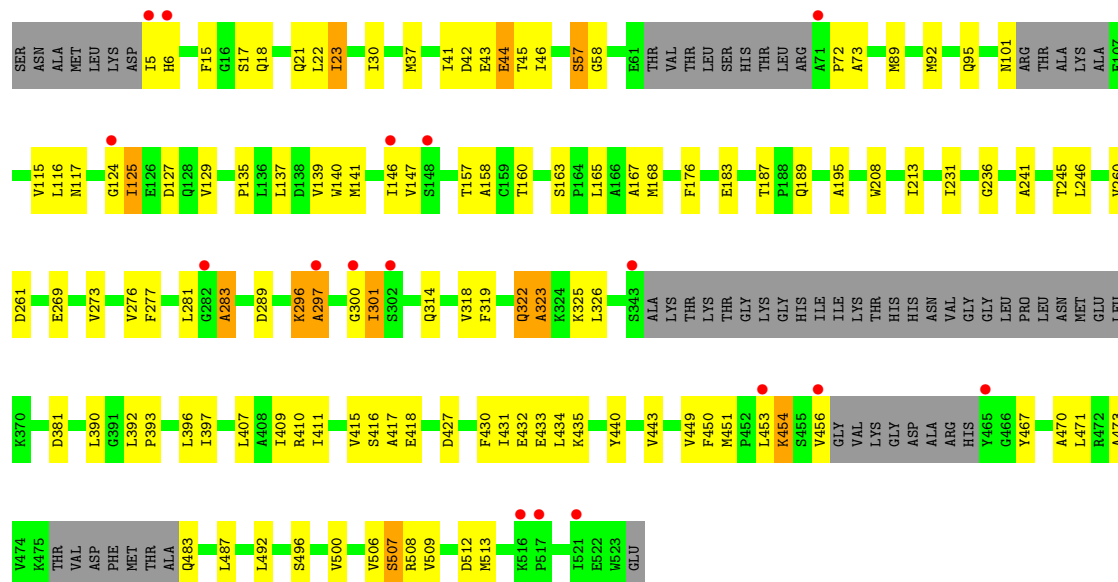


• Molecule 1: GMP synthase [glutamine-hydrolyzing]

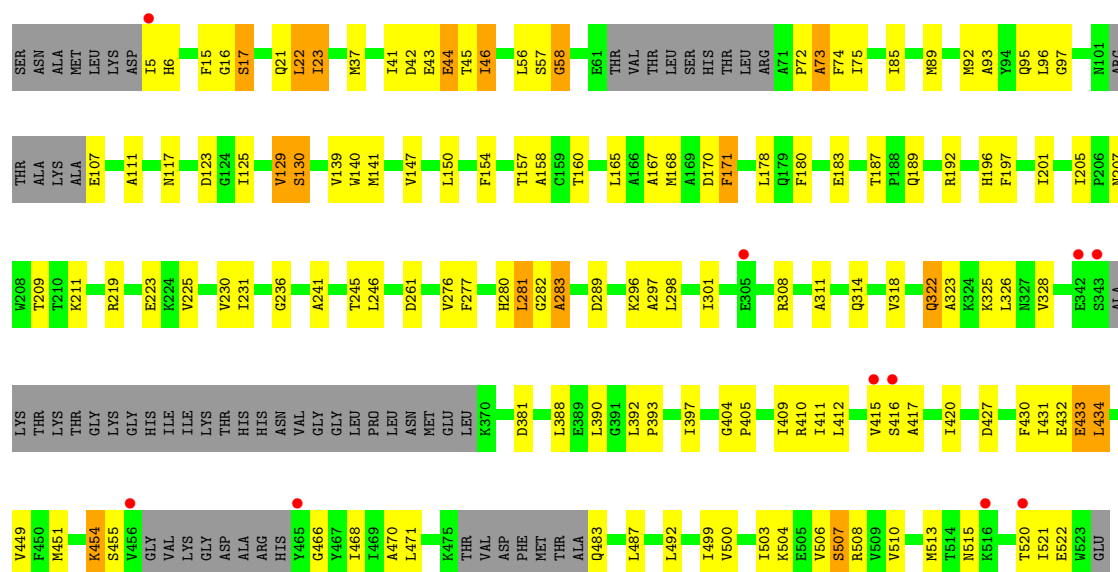




• Molecule 1: GMP synthase [glutamine-hydrolyzing]



• Molecule 1: GMP synthase [glutamine-hydrolyzing]



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	74.91Å 143.95Å 107.89Å 90.00° 97.66° 90.00°	Depositor
Resolution (Å)	39.01 – 2.84 39.01 – 2.84	Depositor EDS
% Data completeness (in resolution range)	98.7 (39.01-2.84) 99.0 (39.01-2.84)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.58 (at 2.86Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.245 , 0.281 (Not available) , 0.275	Depositor DCC
R_{free} test set	2691 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	59.1	Xtriage
Anisotropy	0.167	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 26.1	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	0.91	EDS
Total number of atoms	14708	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.44% 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	0.56	0/3758	0.82	5/5085 (0.1%)
1	B	0.57	0/3758	0.82	5/5085 (0.1%)
1	C	0.56	0/3758	0.81	3/5085 (0.1%)
1	D	0.57	0/3758	0.84	5/5085 (0.1%)
All	All	0.56	0/15032	0.82	18/20340 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	2
1	C	0	3
1	D	0	2
All	All	0	9

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	71	ALA	CA-C-N	6.85	126.32	118.85
1	B	71	ALA	C-N-CA	6.85	126.32	118.85
1	D	23	ILE	N-CA-C	-6.79	103.90	110.42
1	C	23	ILE	N-CA-C	-6.57	104.11	110.42
1	B	433	GLU	N-CA-C	-6.53	101.91	110.53
1	D	433	GLU	N-CA-C	-6.27	101.14	110.23
1	A	23	ILE	N-CA-C	-6.07	104.59	110.42
1	D	322	GLN	N-CA-C	-5.72	102.94	110.43
1	B	23	ILE	N-CA-C	-5.61	105.03	110.42
1	B	58	GLY	N-CA-C	5.37	120.78	112.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	322	GLN	N-CA-C	-5.35	102.36	110.28
1	D	58	GLY	N-CA-C	5.22	120.54	112.45
1	A	187	THR	CA-C-N	5.15	124.59	119.24
1	A	187	THR	C-N-CA	5.15	124.59	119.24
1	A	296	LYS	N-CA-C	-5.14	103.75	110.53
1	D	96	LEU	N-CA-C	-5.07	102.00	109.86
1	C	296	LYS	N-CA-C	-5.03	103.41	110.50
1	A	45	THR	N-CA-C	-5.02	100.10	110.80

There are no chirality outliers.

All (9) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	44	GLU	Peptide
1	A	57	SER	Peptide
1	B	44	GLU	Peptide
1	B	57	SER	Peptide
1	C	432	GLU	Peptide
1	C	44	GLU	Peptide
1	C	57	SER	Peptide
1	D	44	GLU	Peptide
1	D	57	SER	Peptide

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	3677	0	3699	94	0
1	B	3677	0	3699	100	0
1	C	3677	0	3699	111	0
1	D	3677	0	3699	111	0
All	All	14708	0	14796	401	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 14.

All (401) 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:D:449:VAL:HG13	1:D:470:ALA:HB3	1.39	1.00
1:D:451:MET:HE2	1:D:470:ALA:HB2	1.43	0.99
1:C:451:MET:HE2	1:C:470:ALA:HB2	1.43	0.99
1:B:451:MET:CE	1:B:470:ALA:HB2	1.93	0.98
1:D:205:ILE:HD12	1:D:205:ILE:O	1.68	0.93
1:B:483:GLN:N	1:B:483:GLN:OE1	2.03	0.91
1:A:451:MET:HE2	1:A:470:ALA:HB2	1.55	0.88
1:D:246:LEU:HD13	1:D:390:LEU:HD11	1.55	0.87
1:C:241:ALA:O	1:C:245:THR:HG23	1.77	0.85
1:B:451:MET:HE2	1:B:470:ALA:HB2	1.58	0.84
1:B:147:VAL:O	1:B:165:LEU:HD11	1.78	0.83
1:A:147:VAL:O	1:A:165:LEU:HD11	1.77	0.83
1:D:23:ILE:HG12	1:D:183:GLU:HG2	1.60	0.83
1:B:449:VAL:CG2	1:B:470:ALA:HB3	2.09	0.83
1:A:483:GLN:OE1	1:A:483:GLN:N	2.12	0.83
1:B:89:MET:HE3	1:B:168:MET:HG3	1.61	0.82
1:A:158:ALA:HB3	1:A:167:ALA:HB3	1.61	0.81
1:B:147:VAL:HG23	1:B:165:LEU:HD13	1.62	0.81
1:C:411:ILE:HD11	1:C:415:VAL:HG22	1.62	0.80
1:C:513:MET:HE1	1:D:471:LEU:HD12	1.63	0.80
1:C:89:MET:HE3	1:C:168:MET:HG3	1.64	0.79
1:D:520:THR:HG23	1:D:522:GLU:O	1.83	0.79
1:C:487:LEU:HB2	1:C:492:LEU:HD21	1.65	0.79
1:C:147:VAL:O	1:C:165:LEU:HD11	1.82	0.78
1:D:89:MET:HE3	1:D:168:MET:HG3	1.64	0.78
1:A:449:VAL:CG2	1:A:470:ALA:HB3	2.14	0.78
1:A:322:GLN:O	1:A:323:ALA:CB	2.31	0.78
1:B:520:THR:HG23	1:B:522:GLU:O	1.85	0.77
1:C:277:PHE:O	1:C:283:ALA:HB3	1.84	0.77
1:A:92:MET:HE2	1:A:168:MET:HE3	1.67	0.77
1:A:241:ALA:O	1:A:245:THR:HG23	1.84	0.77
1:C:483:GLN:N	1:C:483:GLN:OE1	2.19	0.76
1:B:322:GLN:O	1:B:323:ALA:HB3	1.86	0.75
1:B:23:ILE:HG12	1:B:183:GLU:HG2	1.69	0.75
1:A:277:PHE:O	1:A:283:ALA:HB3	1.86	0.74
1:A:23:ILE:HG12	1:A:183:GLU:HG2	1.70	0.74
1:C:15:PHE:HB2	1:C:58:GLY:HA3	1.69	0.74
1:C:157:THR:HG22	1:C:167:ALA:C	2.14	0.73
1:C:322:GLN:O	1:C:323:ALA:HB3	1.88	0.73
1:B:322:GLN:O	1:B:323:ALA:CB	2.37	0.73
1:C:322:GLN:O	1:C:323:ALA:CB	2.36	0.73
1:B:231:ILE:HD12	1:B:323:ALA:HB2	1.71	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:89:MET:HE2	1:A:168:MET:HG3	1.72	0.72
1:A:147:VAL:O	1:A:165:LEU:CD1	2.37	0.72
1:D:15:PHE:HB2	1:D:58:GLY:HA3	1.72	0.72
1:C:42:ASP:N	1:C:42:ASP:OD1	2.23	0.71
1:D:322:GLN:O	1:D:323:ALA:HB3	1.90	0.71
1:B:15:PHE:HB2	1:B:58:GLY:HA3	1.73	0.71
1:D:231:ILE:HD12	1:D:323:ALA:HB2	1.73	0.70
1:A:322:GLN:O	1:A:323:ALA:HB3	1.90	0.70
1:A:409:ILE:HG22	1:A:410:ARG:H	1.57	0.69
1:A:261:ASP:O	1:A:289:ASP:O	2.11	0.69
1:C:296:LYS:O	1:C:297:ALA:HB3	1.90	0.69
1:D:225:VAL:HG21	1:D:230:VAL:CG2	2.23	0.69
1:D:483:GLN:N	1:D:483:GLN:OE1	2.26	0.68
1:D:111:ALA:HB3	1:D:141:MET:HE2	1.75	0.68
1:A:89:MET:HG3	1:A:168:MET:HE2	1.76	0.68
1:A:281:LEU:O	1:A:281:LEU:HD12	1.91	0.68
1:B:277:PHE:O	1:B:283:ALA:HB3	1.94	0.68
1:C:158:ALA:HB3	1:C:167:ALA:HB3	1.76	0.67
1:B:416:SER:O	1:B:417:ALA:HB3	1.95	0.67
1:C:89:MET:CE	1:C:168:MET:HG3	2.24	0.67
1:C:147:VAL:O	1:C:165:LEU:CD1	2.42	0.66
1:A:449:VAL:HG22	1:A:470:ALA:HB3	1.76	0.66
1:C:411:ILE:HD11	1:C:415:VAL:CG2	2.24	0.65
1:B:89:MET:HG3	1:B:168:MET:HE2	1.78	0.65
1:C:21:GLN:O	1:C:22:LEU:HB2	1.97	0.64
1:B:451:MET:HE3	1:B:470:ALA:HB2	1.79	0.64
1:A:520:THR:HG23	1:A:522:GLU:O	1.97	0.64
1:D:89:MET:HG3	1:D:168:MET:HE2	1.79	0.64
1:C:509:VAL:HG22	1:D:513:MET:HE2	1.79	0.63
1:A:416:SER:O	1:A:417:ALA:CB	2.46	0.63
1:D:141:MET:CE	1:D:160:THR:HG21	2.27	0.63
1:B:140:TRP:N	1:B:187:THR:HG22	2.13	0.63
1:B:300:GLY:C	1:B:301:ILE:HD12	2.24	0.63
1:A:407:LEU:O	1:A:409:ILE:O	2.16	0.63
1:A:92:MET:HE2	1:A:168:MET:CE	2.28	0.63
1:A:21:GLN:NE2	1:A:341:ILE:HD11	2.14	0.63
1:A:21:GLN:O	1:A:22:LEU:CB	2.46	0.63
1:A:296:LYS:O	1:A:297:ALA:HB3	1.98	0.63
1:C:21:GLN:O	1:C:22:LEU:CB	2.47	0.63
1:C:160:THR:HG23	1:C:163:SER:H	1.62	0.63
1:D:141:MET:HE3	1:D:160:THR:HG21	1.81	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:92:MET:HE2	1:C:168:MET:HE3	1.80	0.62
1:B:449:VAL:HG22	1:B:470:ALA:HB3	1.81	0.62
1:B:23:ILE:HD13	1:B:85:ILE:HD12	1.82	0.62
1:B:129:VAL:HG21	1:D:276:VAL:HG22	1.81	0.62
1:C:231:ILE:HD12	1:C:323:ALA:HB2	1.82	0.62
1:A:416:SER:O	1:A:417:ALA:HB3	1.99	0.62
1:D:241:ALA:O	1:D:245:THR:HG23	1.99	0.61
1:D:21:GLN:O	1:D:22:LEU:CB	2.48	0.61
1:D:487:LEU:HB2	1:D:492:LEU:HD21	1.82	0.61
1:A:157:THR:HG22	1:A:167:ALA:C	2.25	0.61
1:A:225:VAL:HG21	1:A:230:VAL:CG2	2.30	0.61
1:A:140:TRP:H	1:A:187:THR:HG22	1.65	0.61
1:B:16:GLY:O	1:B:17:SER:HB3	2.01	0.61
1:D:323:ALA:H	1:D:326:LEU:HD12	1.66	0.60
1:B:100:VAL:HG13	1:B:101:ASN:N	2.16	0.60
1:C:409:ILE:HG22	1:C:410:ARG:H	1.67	0.60
1:A:205:ILE:HD12	1:A:205:ILE:O	2.00	0.60
1:B:276:VAL:HG13	1:B:277:PHE:CD2	2.36	0.60
1:B:431:ILE:HA	1:B:434:LEU:HD12	1.82	0.60
1:B:416:SER:O	1:B:417:ALA:CB	2.49	0.60
1:C:319:PHE:O	1:C:322:GLN:O	2.20	0.60
1:A:21:GLN:O	1:A:22:LEU:HB2	2.02	0.59
1:A:392:LEU:HB2	1:A:397:ILE:HD11	1.83	0.59
1:B:246:LEU:HD13	1:B:390:LEU:HD11	1.83	0.59
1:D:16:GLY:O	1:D:17:SER:HB3	2.01	0.59
1:A:158:ALA:HB3	1:A:167:ALA:CB	2.32	0.59
1:D:427:ASP:O	1:D:431:ILE:HG22	2.03	0.59
1:D:449:VAL:CG1	1:D:470:ALA:HB3	2.25	0.59
1:A:314:GLN:O	1:A:318:VAL:HG23	2.03	0.58
1:D:178:LEU:HD22	1:D:180:PHE:CZ	2.39	0.58
1:C:261:ASP:O	1:C:289:ASP:O	2.21	0.58
1:D:433:GLU:O	1:D:434:LEU:HB2	2.04	0.58
1:A:41:ILE:HD11	1:A:45:THR:HG21	1.85	0.58
1:D:409:ILE:HG22	1:D:410:ARG:H	1.68	0.58
1:C:427:ASP:O	1:C:431:ILE:HG22	2.04	0.58
1:D:157:THR:HG22	1:D:167:ALA:C	2.29	0.58
1:B:158:ALA:HB3	1:B:167:ALA:HB3	1.86	0.57
1:B:129:VAL:HG12	1:B:130:SER:N	2.19	0.57
1:C:5:ILE:HG22	1:C:6:HIS:H	1.70	0.57
1:D:205:ILE:HD12	1:D:205:ILE:C	2.29	0.57
1:B:433:GLU:O	1:B:434:LEU:HB2	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:5:ILE:HD12	1:C:5:ILE:N	2.19	0.57
1:C:281:LEU:HD12	1:C:281:LEU:O	2.05	0.57
1:A:260:VAL:HG23	1:A:262:THR:HG23	1.87	0.56
1:C:381:ASP:OD1	1:C:381:ASP:N	2.37	0.56
1:A:160:THR:HG22	1:A:163:SER:HB3	1.87	0.56
1:C:411:ILE:CD1	1:C:415:VAL:HG22	2.33	0.56
1:A:92:MET:HE1	1:A:176:PHE:C	2.31	0.56
1:D:147:VAL:O	1:D:165:LEU:HD13	2.05	0.56
1:B:140:TRP:H	1:B:187:THR:HG22	1.69	0.56
1:B:454:LYS:HA	1:B:466:GLY:O	2.06	0.56
1:A:23:ILE:HD13	1:A:85:ILE:HD12	1.88	0.56
1:C:269:GLU:O	1:C:273:VAL:HG12	2.06	0.55
1:A:392:LEU:CB	1:A:397:ILE:HD11	2.35	0.55
1:A:503:ILE:HG22	1:A:504:LYS:O	2.07	0.55
1:D:107:GLU:N	1:D:107:GLU:OE1	2.39	0.55
1:A:433:GLU:O	1:A:434:LEU:HB2	2.06	0.55
1:B:449:VAL:HG23	1:B:470:ALA:HB3	1.85	0.55
1:C:296:LYS:O	1:C:297:ALA:CB	2.52	0.55
1:C:89:MET:HG3	1:C:168:MET:HE2	1.89	0.55
1:C:314:GLN:O	1:C:318:VAL:HG23	2.07	0.54
1:B:409:ILE:HG22	1:B:410:ARG:H	1.72	0.54
1:C:141:MET:HE3	1:C:160:THR:HG21	1.89	0.54
1:C:433:GLU:O	1:C:434:LEU:HB2	2.07	0.54
1:D:140:TRP:N	1:D:187:THR:HG22	2.22	0.54
1:B:21:GLN:O	1:B:22:LEU:CB	2.54	0.54
1:C:411:ILE:HG22	1:C:450:PHE:HB3	1.90	0.54
1:D:21:GLN:O	1:D:22:LEU:HB2	2.07	0.54
1:D:140:TRP:H	1:D:187:THR:HG22	1.71	0.54
1:B:115:VAL:HB	1:B:135:PRO:HG2	1.89	0.54
1:A:21:GLN:CD	1:A:341:ILE:HD11	2.33	0.54
1:B:111:ALA:HB3	1:B:141:MET:HG3	1.90	0.53
1:C:116:LEU:O	1:C:157:THR:O	2.25	0.53
1:D:277:PHE:O	1:D:283:ALA:HB3	2.08	0.53
1:C:453:LEU:O	1:C:454:LYS:HB2	2.09	0.53
1:D:500:VAL:HG13	1:D:507:SER:HA	1.90	0.53
1:B:72:PRO:O	1:B:73:ALA:HB3	2.08	0.53
1:B:178:LEU:HD22	1:B:180:PHE:CZ	2.42	0.53
1:B:178:LEU:HD22	1:B:180:PHE:CE2	2.44	0.53
1:C:125:ILE:HG22	1:C:137:LEU:HD21	1.89	0.53
1:A:101:ASN:CB	1:A:146:ILE:HG22	2.39	0.53
1:C:487:LEU:CB	1:C:492:LEU:HD21	2.36	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:322:GLN:O	1:D:323:ALA:CB	2.55	0.53
1:C:410:ARG:HA	1:C:449:VAL:HG13	1.89	0.53
1:C:507:SER:OG	1:C:507:SER:O	2.25	0.53
1:A:129:VAL:O	1:A:134:GLU:O	2.27	0.53
1:B:263:GLY:O	1:B:420:ILE:HD12	2.09	0.53
1:B:393:PRO:HG3	1:D:129:VAL:HG23	1.91	0.53
1:B:147:VAL:O	1:B:165:LEU:CD1	2.52	0.53
1:C:496:SER:HB3	1:D:513:MET:HE3	1.91	0.53
1:D:261:ASP:O	1:D:289:ASP:O	2.27	0.53
1:C:246:LEU:HD13	1:C:390:LEU:HD11	1.91	0.53
1:D:392:LEU:HB2	1:D:397:ILE:HD11	1.91	0.53
1:B:319:PHE:O	1:B:322:GLN:O	2.27	0.53
1:A:90:GLN:HE21	1:A:100:VAL:HG22	1.74	0.52
1:B:225:VAL:HG21	1:B:230:VAL:CG2	2.39	0.52
1:D:72:PRO:HD2	1:D:95:GLN:OE1	2.08	0.52
1:D:381:ASP:OD1	1:D:381:ASP:N	2.32	0.52
1:A:189:GLN:NE2	1:A:192:ARG:HH11	2.08	0.52
1:B:23:ILE:CD1	1:B:85:ILE:HD12	2.40	0.52
1:B:32:VAL:HG21	1:B:198:VAL:CG1	2.40	0.52
1:C:431:ILE:HD13	1:C:440:TYR:CD1	2.44	0.52
1:B:157:THR:HG23	1:B:167:ALA:CB	2.40	0.52
1:D:430:PHE:O	1:D:433:GLU:O	2.27	0.52
1:A:297:ALA:O	1:A:310:ILE:HG21	2.10	0.52
1:C:127:ASP:OD1	1:C:189:GLN:NE2	2.44	0.51
1:C:101:ASN:CB	1:C:146:ILE:HG22	2.40	0.51
1:A:409:ILE:HG22	1:A:410:ARG:N	2.25	0.51
1:C:396:LEU:HD23	1:C:396:LEU:C	2.36	0.51
1:C:507:SER:O	1:C:508:ARG:HB3	2.09	0.51
1:C:129:VAL:CG1	1:C:129:VAL:O	2.58	0.51
1:A:483:GLN:HE22	1:A:516:LYS:CD	2.23	0.51
1:A:520:THR:HG21	1:B:456:VAL:HB	1.92	0.51
1:D:158:ALA:HB3	1:D:167:ALA:HB3	1.92	0.50
1:A:243:THR:O	1:A:247:VAL:HG23	2.11	0.50
1:A:510:VAL:HG12	1:B:512:ASP:HA	1.93	0.50
1:C:512:ASP:HA	1:D:510:VAL:HG12	1.93	0.50
1:A:483:GLN:HE22	1:A:516:LYS:HD2	1.77	0.50
1:D:503:ILE:HG22	1:D:504:LYS:O	2.12	0.50
1:A:431:ILE:HA	1:A:434:LEU:HD12	1.93	0.50
1:B:101:ASN:HB2	1:B:146:ILE:HG22	1.93	0.50
1:C:430:PHE:CD2	1:C:471:LEU:HD22	2.47	0.50
1:D:45:THR:O	1:D:46:ILE:HB	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:157:THR:HG22	1:B:167:ALA:C	2.37	0.50
1:C:409:ILE:HG22	1:C:410:ARG:N	2.27	0.49
1:A:506:VAL:HG12	1:A:507:SER:N	2.26	0.49
1:D:219:ARG:O	1:D:223:GLU:HG3	2.13	0.49
1:D:454:LYS:HA	1:D:466:GLY:O	2.12	0.49
1:C:392:LEU:HB2	1:C:397:ILE:HD11	1.94	0.49
1:C:411:ILE:C	1:C:411:ILE:HD12	2.37	0.49
1:C:30:ILE:HD13	1:C:195:ALA:HA	1.93	0.49
1:D:276:VAL:O	1:D:280:HIS:O	2.29	0.49
1:A:267:LEU:HD22	1:A:268:ASN:ND2	2.28	0.49
1:A:298:LEU:HA	1:A:301:ILE:HD13	1.94	0.49
1:D:433:GLU:O	1:D:434:LEU:CB	2.59	0.49
1:A:17:SER:CB	1:A:58:GLY:HA2	2.43	0.49
1:A:72:PRO:O	1:A:73:ALA:HB3	2.11	0.49
1:A:101:ASN:HB2	1:A:146:ILE:HG22	1.94	0.49
1:B:22:LEU:HD13	1:B:25:ARG:HH21	1.78	0.49
1:A:41:ILE:CD1	1:A:45:THR:HG21	2.43	0.49
1:B:433:GLU:O	1:B:434:LEU:CB	2.60	0.49
1:D:308:ARG:O	1:D:311:ALA:O	2.31	0.49
1:C:276:VAL:HG23	1:C:277:PHE:CD2	2.48	0.48
1:A:90:GLN:NE2	1:A:100:VAL:HG22	2.28	0.48
1:B:92:MET:HE2	1:B:168:MET:HE3	1.95	0.48
1:D:411:ILE:HD11	1:D:420:ILE:HG22	1.95	0.48
1:D:455:SER:HB3	1:D:468:ILE:HD11	1.95	0.48
1:B:506:VAL:HG12	1:B:507:SER:N	2.28	0.48
1:C:140:TRP:H	1:C:187:THR:HG22	1.78	0.48
1:C:396:LEU:HD23	1:C:396:LEU:O	2.13	0.48
1:A:45:THR:O	1:A:46:ILE:HB	2.14	0.48
1:B:129:VAL:HG23	1:D:393:PRO:HG3	1.96	0.48
1:B:431:ILE:C	1:B:433:GLU:O	2.57	0.48
1:D:129:VAL:HG12	1:D:130:SER:N	2.28	0.48
1:D:74:PHE:C	1:D:74:PHE:CD2	2.90	0.48
1:C:45:THR:O	1:C:46:ILE:HB	2.14	0.48
1:D:157:THR:HG23	1:D:167:ALA:CB	2.44	0.48
1:B:16:GLY:O	1:B:17:SER:CB	2.61	0.47
1:B:396:LEU:C	1:B:396:LEU:HD23	2.39	0.47
1:D:23:ILE:HD13	1:D:85:ILE:HD12	1.95	0.47
1:D:41:ILE:HD11	1:D:45:THR:HG21	1.95	0.47
1:D:392:LEU:CB	1:D:397:ILE:HD11	2.43	0.47
1:A:170:ASP:O	1:A:171:PHE:HB2	2.14	0.47
1:B:514:THR:HG21	1:B:519:ALA:HB2	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:467:TYR:O	1:C:506:VAL:O	2.33	0.47
1:C:471:LEU:HD12	1:D:513:MET:HE1	1.96	0.47
1:D:197:PHE:HA	1:D:201:ILE:HD12	1.97	0.47
1:D:506:VAL:O	1:D:507:SER:OG	2.28	0.47
1:D:45:THR:O	1:D:46:ILE:CB	2.63	0.47
1:C:431:ILE:HD12	1:C:431:ILE:C	2.40	0.47
1:B:411:ILE:HD11	1:B:420:ILE:HG22	1.97	0.47
1:C:72:PRO:O	1:C:73:ALA:HB3	2.15	0.47
1:D:92:MET:HE2	1:D:168:MET:CE	2.44	0.47
1:D:506:VAL:HG12	1:D:507:SER:N	2.29	0.47
1:B:147:VAL:CG2	1:B:165:LEU:HD13	2.38	0.47
1:A:160:THR:HG22	1:A:163:SER:CB	2.45	0.47
1:A:433:GLU:O	1:A:434:LEU:CB	2.63	0.47
1:D:280:HIS:O	1:D:281:LEU:CG	2.63	0.47
1:B:72:PRO:O	1:B:73:ALA:CB	2.63	0.47
1:C:416:SER:O	1:C:418:GLU:N	2.47	0.47
1:C:5:ILE:HG22	1:C:6:HIS:N	2.30	0.46
1:D:72:PRO:O	1:D:73:ALA:HB3	2.14	0.46
1:B:125:ILE:HG23	1:B:189:GLN:CG	2.45	0.46
1:D:170:ASP:O	1:D:171:PHE:HB2	2.15	0.46
1:A:92:MET:HE3	1:A:154:PHE:CE1	2.50	0.46
1:B:125:ILE:O	1:B:137:LEU:HD11	2.15	0.46
1:D:139:VAL:HA	1:D:187:THR:HB	1.98	0.46
1:B:45:THR:O	1:B:46:ILE:HB	2.14	0.46
1:C:260:VAL:O	1:C:260:VAL:HG23	2.15	0.46
1:C:411:ILE:CG1	1:C:415:VAL:HG22	2.45	0.46
1:D:298:LEU:HD13	1:D:415:VAL:HG11	1.97	0.46
1:A:466:GLY:C	1:A:506:VAL:O	2.59	0.46
1:B:430:PHE:CD2	1:B:471:LEU:HD22	2.51	0.46
1:B:125:ILE:HG23	1:B:189:GLN:HG3	1.97	0.46
1:C:101:ASN:HB2	1:C:146:ILE:HG22	1.98	0.46
1:C:158:ALA:HB3	1:C:167:ALA:CB	2.45	0.46
1:C:23:ILE:HG12	1:C:183:GLU:HG2	1.98	0.45
1:C:139:VAL:HA	1:C:187:THR:HB	1.98	0.45
1:B:289:ASP:O	1:B:291:LYS:N	2.50	0.45
1:C:509:VAL:HG22	1:D:513:MET:CE	2.44	0.45
1:D:280:HIS:O	1:D:281:LEU:HG	2.17	0.45
1:A:381:ASP:OD1	1:A:381:ASP:N	2.47	0.45
1:C:140:TRP:N	1:C:187:THR:HG22	2.31	0.45
1:D:189:GLN:HE21	1:D:192:ARG:HD3	1.81	0.45
1:D:16:GLY:O	1:D:17:SER:CB	2.62	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:409:ILE:HG22	1:B:410:ARG:N	2.32	0.45
1:D:5:ILE:O	1:D:6:HIS:HB2	2.17	0.45
1:B:100:VAL:CG1	1:B:101:ASN:N	2.78	0.45
1:C:101:ASN:HB3	1:C:146:ILE:HG22	1.99	0.45
1:D:123:ASP:O	1:D:196:HIS:HE1	2.00	0.45
1:D:231:ILE:HG13	1:D:328:VAL:HG11	1.99	0.45
1:A:487:LEU:HB2	1:A:492:LEU:HD21	1.99	0.44
1:B:257:CYS:HB2	1:B:285:VAL:HG22	1.99	0.44
1:C:129:VAL:O	1:C:129:VAL:HG12	2.17	0.44
1:A:205:ILE:HD12	1:A:205:ILE:C	2.41	0.44
1:A:507:SER:O	1:A:507:SER:OG	2.35	0.44
1:C:23:ILE:CG1	1:C:183:GLU:HG2	2.48	0.44
1:C:45:THR:O	1:C:46:ILE:CB	2.66	0.44
1:C:92:MET:HE2	1:C:168:MET:CE	2.46	0.44
1:C:208:TRP:CE2	1:C:213:ILE:HD11	2.53	0.44
1:D:157:THR:HG23	1:D:167:ALA:HB3	1.99	0.44
1:C:72:PRO:HD2	1:C:95:GLN:OE1	2.17	0.44
1:A:44:GLU:C	1:A:45:THR:O	2.60	0.44
1:C:392:LEU:CB	1:C:397:ILE:HD11	2.48	0.44
1:A:115:VAL:HB	1:A:135:PRO:HG2	1.99	0.44
1:C:115:VAL:HB	1:C:135:PRO:HG2	1.99	0.44
1:D:209:THR:HG22	1:D:211:LYS:H	1.83	0.44
1:D:507:SER:O	1:D:508:ARG:HB3	2.18	0.44
1:A:296:LYS:O	1:A:297:ALA:CB	2.63	0.44
1:B:129:VAL:HG12	1:B:130:SER:H	1.83	0.44
1:C:37:MET:SD	1:C:41:ILE:HD13	2.58	0.44
1:C:453:LEU:HD13	1:D:412:LEU:HD22	1.99	0.44
1:D:409:ILE:HG22	1:D:410:ARG:N	2.33	0.43
1:A:514:THR:HG21	1:B:508:ARG:HA	1.99	0.43
1:B:100:VAL:CG2	1:B:147:VAL:HG12	2.48	0.43
1:B:157:THR:HG23	1:B:167:ALA:HB1	2.00	0.43
1:D:483:GLN:N	1:D:515:ASN:OD1	2.51	0.43
1:A:187:THR:O	1:A:187:THR:OG1	2.33	0.43
1:B:45:THR:O	1:B:46:ILE:CB	2.66	0.43
1:D:93:ALA:O	1:D:97:GLY:O	2.36	0.43
1:A:89:MET:CE	1:A:168:MET:HG3	2.46	0.43
1:A:409:ILE:C	1:A:411:ILE:H	2.26	0.43
1:D:37:MET:SD	1:D:41:ILE:HD13	2.59	0.43
1:A:125:ILE:HD13	1:A:192:ARG:CZ	2.49	0.43
1:C:500:VAL:HG13	1:C:507:SER:HA	2.01	0.43
1:B:92:MET:HE3	1:B:154:PHE:CZ	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:451:MET:HE2	1:B:470:ALA:CB	2.38	0.43
1:D:507:SER:OG	1:D:507:SER:O	2.35	0.43
1:B:74:PHE:CD2	1:B:74:PHE:C	2.94	0.43
1:C:393:PRO:O	1:C:397:ILE:HD13	2.19	0.43
1:D:141:MET:HE1	1:D:160:THR:HG21	1.98	0.43
1:B:140:TRP:H	1:B:187:THR:CG2	2.32	0.42
1:B:323:ALA:H	1:B:326:LEU:HD12	1.84	0.42
1:C:431:ILE:HD11	1:C:435:LYS:HD3	2.00	0.42
1:D:296:LYS:O	1:D:297:ALA:HB3	2.19	0.42
1:D:42:ASP:OD1	1:D:42:ASP:N	2.47	0.42
1:C:92:MET:HE1	1:C:176:PHE:C	2.44	0.42
1:D:431:ILE:C	1:D:431:ILE:HD12	2.43	0.42
1:A:312:GLY:O	1:A:316:ILE:HG13	2.18	0.42
1:B:453:LEU:O	1:B:454:LYS:HB2	2.19	0.42
1:C:407:LEU:C	1:C:409:ILE:O	2.62	0.42
1:A:5:ILE:HG23	1:A:6:HIS:N	2.34	0.42
1:B:253:ASP:OD1	1:B:253:ASP:N	2.45	0.42
1:B:431:ILE:O	1:B:433:GLU:O	2.37	0.42
1:C:300:GLY:C	1:C:301:ILE:HD12	2.45	0.42
1:D:314:GLN:O	1:D:318:VAL:HG23	2.20	0.42
1:C:443:VAL:HG21	1:C:473:ALA:HB1	2.01	0.42
1:D:178:LEU:HD22	1:D:180:PHE:CE2	2.55	0.42
1:D:416:SER:O	1:D:417:ALA:HB3	2.20	0.42
1:D:499:ILE:HG23	1:D:503:ILE:HD12	2.02	0.42
1:B:100:VAL:HG23	1:B:147:VAL:HG12	2.01	0.42
1:A:101:ASN:HB3	1:A:146:ILE:HG22	2.00	0.42
1:A:205:ILE:CD1	1:A:207:ASN:ND2	2.83	0.42
1:D:521:ILE:HD12	1:D:521:ILE:H	1.84	0.42
1:B:21:GLN:O	1:B:22:LEU:HB2	2.20	0.42
1:B:209:THR:HG22	1:B:210:THR:N	2.35	0.42
1:B:467:TYR:O	1:B:506:VAL:O	2.38	0.42
1:C:276:VAL:CG2	1:C:277:PHE:CD2	3.03	0.42
1:C:453:LEU:O	1:C:454:LYS:CB	2.68	0.42
1:C:456:VAL:HB	1:D:520:THR:HG21	2.02	0.42
1:C:57:SER:OG	1:C:58:GLY:N	2.48	0.41
1:D:56:LEU:HD21	1:D:75:ILE:CD1	2.50	0.41
1:C:245:THR:HG21	1:C:392:LEU:HD21	2.01	0.41
1:A:393:PRO:O	1:A:397:ILE:HD13	2.20	0.41
1:A:453:LEU:HD13	1:B:412:LEU:HD12	2.01	0.41
1:B:114:ARG:HA	1:B:136:LEU:HD23	2.03	0.41
1:B:411:ILE:CD1	1:B:420:ILE:HG22	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:72:PRO:HD2	1:B:95:GLN:OE1	2.20	0.41
1:B:371:LEU:HD12	1:B:371:LEU:N	2.35	0.41
1:C:17:SER:O	1:C:18:GLN:C	2.63	0.41
1:D:205:ILE:CD1	1:D:207:ASN:ND2	2.83	0.41
1:A:89:MET:CG	1:A:168:MET:HE2	2.46	0.41
1:C:187:THR:O	1:C:187:THR:OG1	2.37	0.41
1:C:124:GLY:O	1:C:125:ILE:C	2.63	0.41
1:D:404:GLY:N	1:D:405:PRO:HD2	2.36	0.41
1:A:43:GLU:O	1:A:44:GLU:CB	2.69	0.41
1:C:41:ILE:HD11	1:C:45:THR:HG21	2.03	0.41
1:D:43:GLU:O	1:D:44:GLU:HB3	2.21	0.41
1:D:92:MET:HE3	1:D:154:PHE:CE1	2.56	0.41
1:D:388:LEU:HD12	1:D:388:LEU:HA	1.96	0.41
1:A:17:SER:HB2	1:A:58:GLY:HA2	2.02	0.41
1:C:323:ALA:H	1:C:326:LEU:HD12	1.85	0.41
1:C:411:ILE:HD12	1:C:411:ILE:O	2.21	0.41
1:D:150:LEU:HG	1:D:165:LEU:HD23	2.02	0.41
1:B:60:PRO:O	1:B:61:GLU:C	2.65	0.40
1:B:130:SER:HB3	1:B:134:GLU:O	2.22	0.40
1:D:411:ILE:HD11	1:D:420:ILE:CG2	2.51	0.40
1:A:37:MET:SD	1:A:41:ILE:HD13	2.62	0.40
1:A:181:HIS:O	1:A:187:THR:CG2	2.69	0.40
1:A:234:LEU:HD23	1:A:234:LEU:HA	1.94	0.40
1:B:242:VAL:HG12	1:B:375:LEU:HD21	2.03	0.40
1:C:43:GLU:O	1:C:44:GLU:CB	2.69	0.40
1:A:280:HIS:O	1:A:281:LEU:HG	2.22	0.40
1:C:396:LEU:HD22	1:C:397:ILE:HD12	2.03	0.40
1:D:92:MET:CE	1:D:168:MET:HE3	2.52	0.40
1:B:281:LEU:O	1:B:281:LEU:CG	2.69	0.40
1:C:125:ILE:HG23	1:C:189:GLN:HG3	2.03	0.40
1:D:281:LEU:O	1:D:282:GLY:C	2.65	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	416/527 (79%)	368 (88%)	33 (8%)	15 (4%)	2	5
1	B	134/527 (25%)	119 (89%)	7 (5%)	8 (6%)	1	1
1	C	383/527 (73%)	343 (90%)	29 (8%)	11 (3%)	3	8
1	D	452/527 (86%)	394 (87%)	40 (9%)	18 (4%)	2	3
All	All	1385/2108 (66%)	1224 (88%)	109 (8%)	52 (4%)	2	4

All (52) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	17	SER
1	A	73	ALA
1	A	210	THR
1	A	417	ALA
1	B	22	LEU
1	B	73	ALA
1	C	125	ILE
1	C	323	ALA
1	C	417	ALA
1	D	46	ILE
1	D	73	ALA
1	D	434	LEU
1	A	22	LEU
1	A	283	ALA
1	C	325	LYS
1	D	22	LEU
1	D	129	VAL
1	A	507	SER
1	B	17	SER
1	B	282	GLY
1	B	290	ALA
1	C	236	GLY
1	D	171	PHE
1	D	281	LEU
1	D	325	LYS
1	A	171	PHE
1	A	290	ALA
1	A	434	LEU
1	B	45	THR
1	C	297	ALA

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Mol	Chain	Res	Type
1	C	454	LYS
1	C	507	SER
1	D	130	SER
1	D	236	GLY
1	A	45	THR
1	A	72	PRO
1	A	508	ARG
1	B	46	ILE
1	B	283	ALA
1	C	283	ALA
1	D	17	SER
1	D	283	ALA
1	D	432	GLU
1	D	507	SER
1	A	125	ILE
1	A	301	ILE
1	C	117	ASN
1	D	117	ASN
1	D	125	ILE
1	D	454	LYS
1	C	301	ILE
1	D	301	ILE

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

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

There are no ligands in this entry.

5.7 Other polymers

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	464/527 (88%)	0.27	12 (2%) 57 48	35, 55, 80, 98	0
1	B	464/527 (88%)	0.26	12 (2%) 57 48	35, 53, 80, 102	0
1	C	464/527 (88%)	0.28	17 (3%) 45 35	36, 56, 84, 96	0
1	D	464/527 (88%)	0.23	10 (2%) 62 53	33, 53, 80, 93	0
All	All	1856/2108 (88%)	0.26	51 (2%) 56 47	33, 54, 81, 102	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	343	SER	5.1
1	B	71	ALA	4.7
1	A	71	ALA	4.4
1	B	5	ILE	4.4
1	D	5	ILE	3.9
1	A	5	ILE	3.9
1	C	5	ILE	3.8
1	C	465	TYR	3.7
1	B	124	GLY	3.4
1	C	71	ALA	3.4
1	D	343	SER	3.3
1	D	465	TYR	3.3
1	C	517	PRO	3.2
1	B	46	ILE	3.2
1	B	343	SER	3.2
1	A	465	TYR	3.1
1	B	465	TYR	3.0
1	A	422	ILE	2.9
1	A	343	SER	2.9
1	B	129	VAL	2.9
1	A	516	LYS	2.8

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Mol	Chain	Res	Type	RSRZ
1	C	148	SER	2.8
1	C	516	LYS	2.8
1	A	124	GLY	2.8
1	A	324	LYS	2.7
1	D	305	GLU	2.7
1	B	516	LYS	2.6
1	C	124	GLY	2.6
1	B	413	GLY	2.5
1	C	297	ALA	2.4
1	C	453	LEU	2.4
1	A	417	ALA	2.3
1	A	72	PRO	2.3
1	D	415	VAL	2.3
1	C	302	SER	2.3
1	D	456	VAL	2.3
1	D	516	LYS	2.3
1	C	456	VAL	2.3
1	C	521	ILE	2.2
1	B	497	HIS	2.2
1	C	146	ILE	2.2
1	D	520	THR	2.2
1	D	342	GLU	2.2
1	B	72	PRO	2.1
1	D	416	SER	2.1
1	C	6	HIS	2.1
1	A	299	LYS	2.1
1	C	300	GLY	2.1
1	B	16	GLY	2.0
1	A	94	TYR	2.0
1	C	282	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.