



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:01 AM UTC

PDB ID : 6TTJ / pdb_00006ttj
Title : Neutral invertase 2 from Arabidopsis thaliana
Authors : Tsirkone, V.G.; Osipov, E.M.; Beelen, S.; Strelkov, S.V.
Deposited on : 2019-12-27
Resolution : 3.39 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

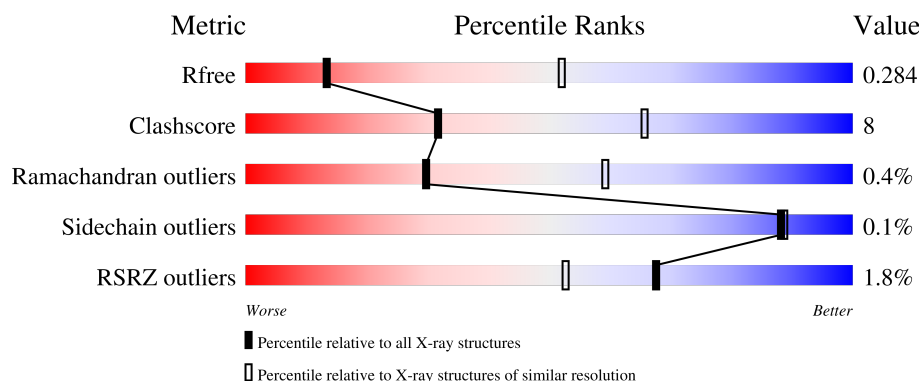
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.39 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	1001 (3.44-3.36)
Clashscore	190562	1022 (3.44-3.36)
Ramachandran outliers	187476	1012 (3.44-3.36)
Sidechain outliers	187428	1012 (3.44-3.36)
RSRZ outliers	180081	1001 (3.44-3.36)

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	% 60% 18% 21%
1	B	551	61% 18% 21%
1	C	551	% 66% 15% 20%
1	D	551	% 63% 17% 20%
1	E	551	% 62% 17% 21%

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Mol	Chain	Length	Quality of chain
1	F	551	2% 64% 16% 20%
1	G	551	% 63% 17% 20%
1	H	551	3% 65% 15% 20%
1	I	551	% 63% 16% 21%
1	J	551	% 64% 16% 20%
1	K	551	2% 62% 17% 20%
1	L	551	2% 65% 15% 21%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 42174 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.

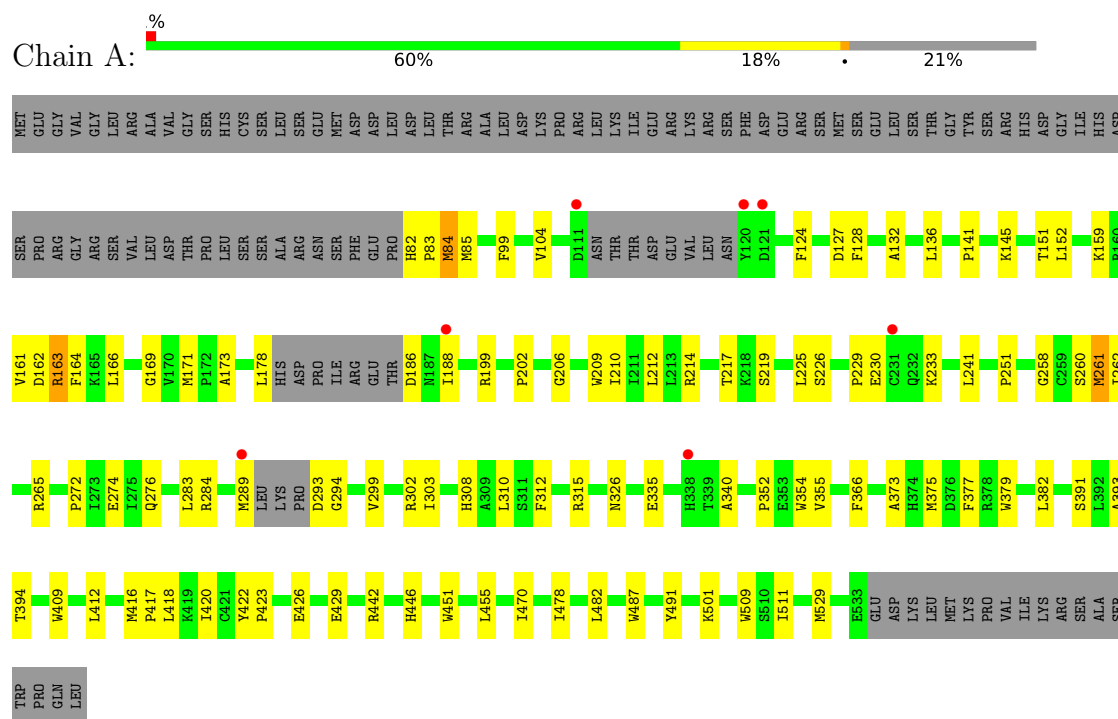
- Molecule 1 is a protein called Alkaline/neutral invertase CINV1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	434	Total	C	N	O	S	0	0	0
			3464	2238	591	607	28			
1	B	438	Total	C	N	O	S	0	0	0
			3499	2259	599	613	28			
1	C	441	Total	C	N	O	S	0	0	0
			3536	2282	604	621	29			
1	D	440	Total	C	N	O	S	0	0	0
			3526	2276	604	618	28			
1	E	438	Total	C	N	O	S	0	0	0
			3511	2264	602	617	28			
1	F	441	Total	C	N	O	S	0	0	0
			3540	2284	607	621	28			
1	G	440	Total	C	N	O	S	0	0	0
			3536	2281	606	620	29			
1	H	439	Total	C	N	O	S	0	0	0
			3525	2275	605	617	28			
1	I	434	Total	C	N	O	S	0	0	0
			3478	2247	596	607	28			
1	J	440	Total	C	N	O	S	0	0	0
			3534	2280	606	620	28			
1	K	439	Total	C	N	O	S	0	0	0
			3523	2273	603	619	28			
1	L	437	Total	C	N	O	S	0	0	0
			3502	2262	598	614	28			

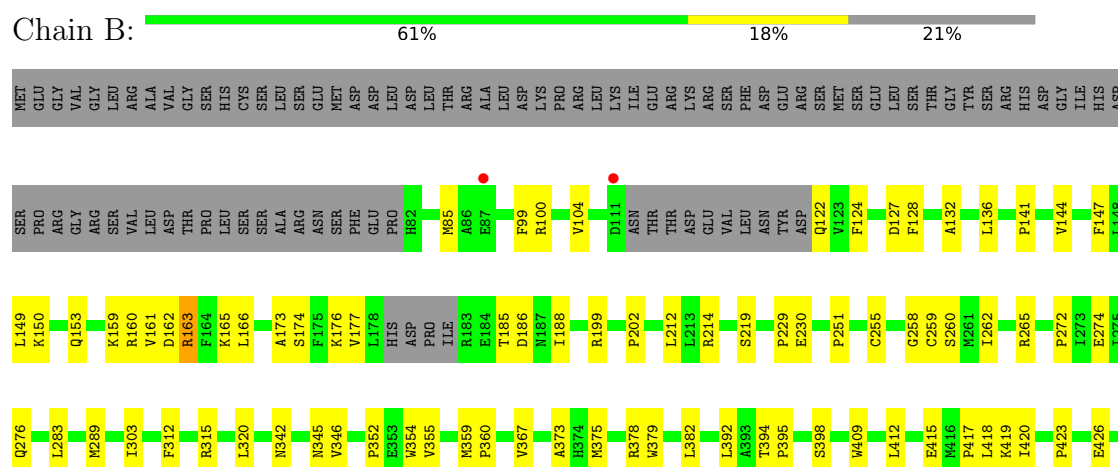
3 Residue-property plots

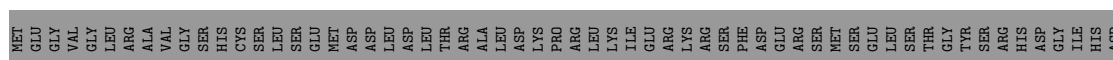
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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

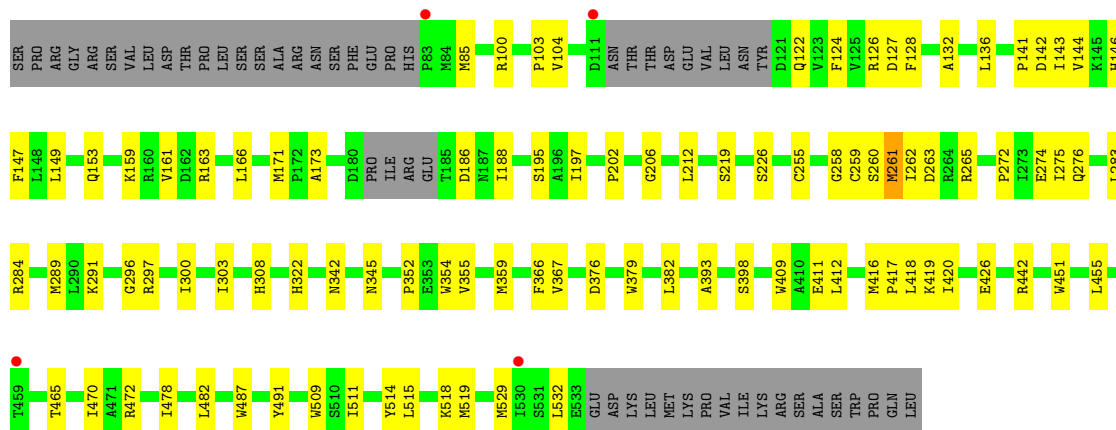
- Molecule 1: Alkaline/neutral invertase CINV1



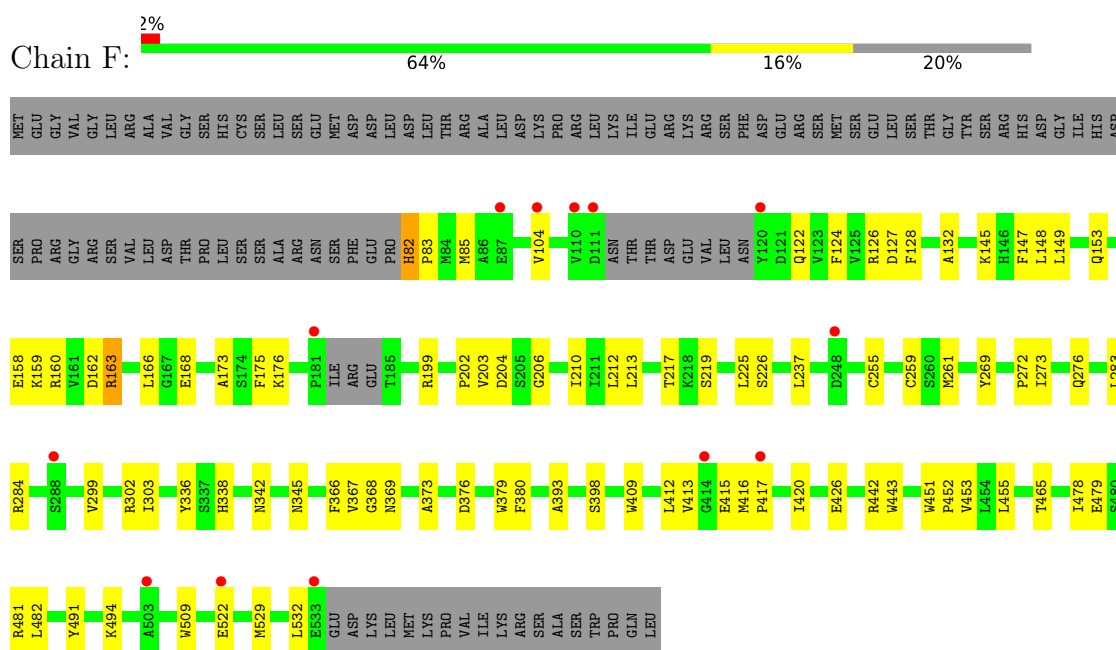
- Molecule 1: Alkaline/neutral invertase CINV1



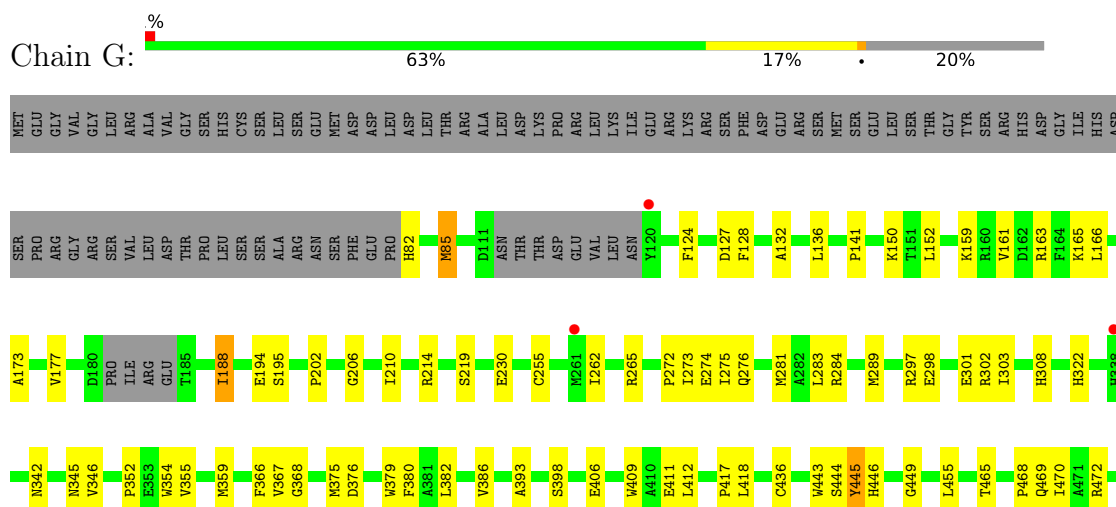


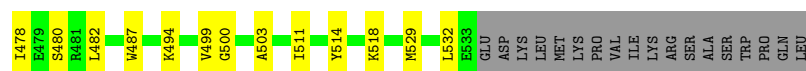


• Molecule 1: Alkaline/neutral invertase CINV1

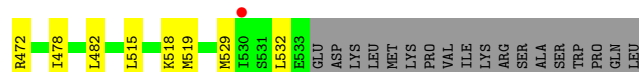
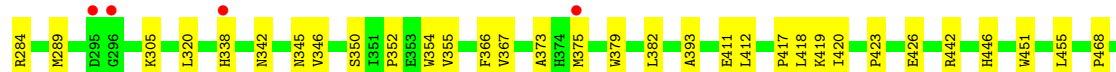
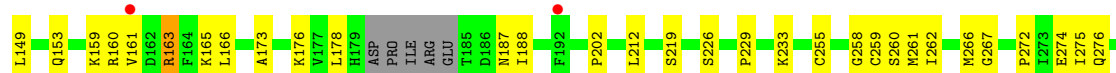
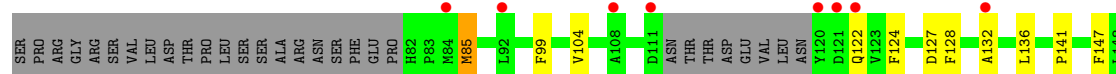
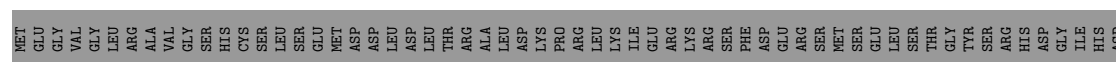


• Molecule 1: Alkaline/neutral invertase CINV1

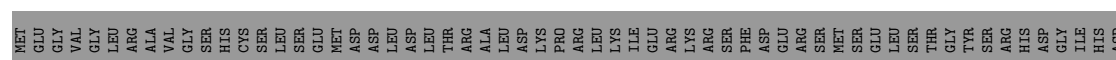




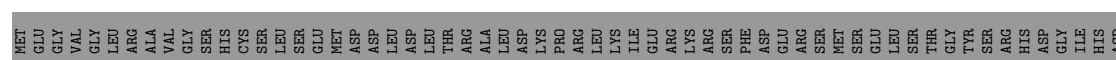
• Molecule 1: Alkaline/neutral invertase CINV1

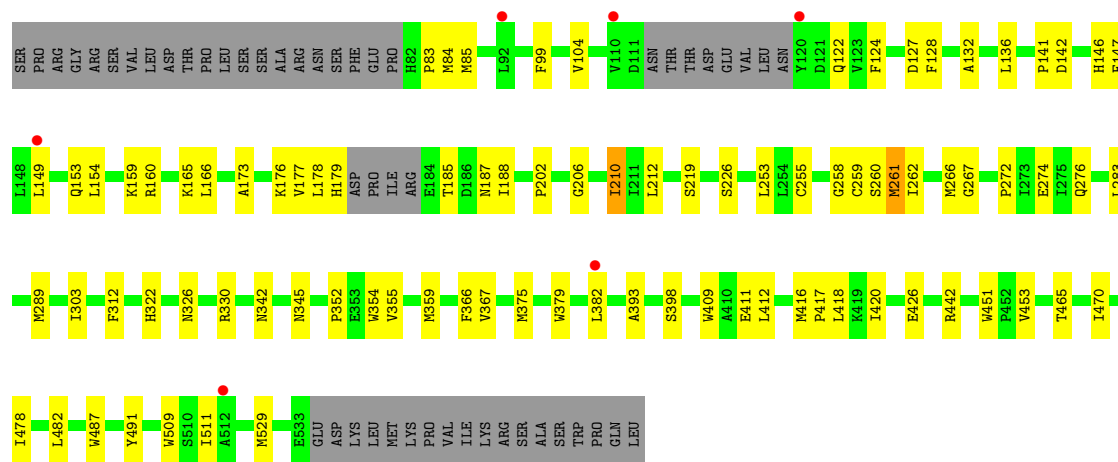


• Molecule 1: Alkaline/neutral invertase CINV1

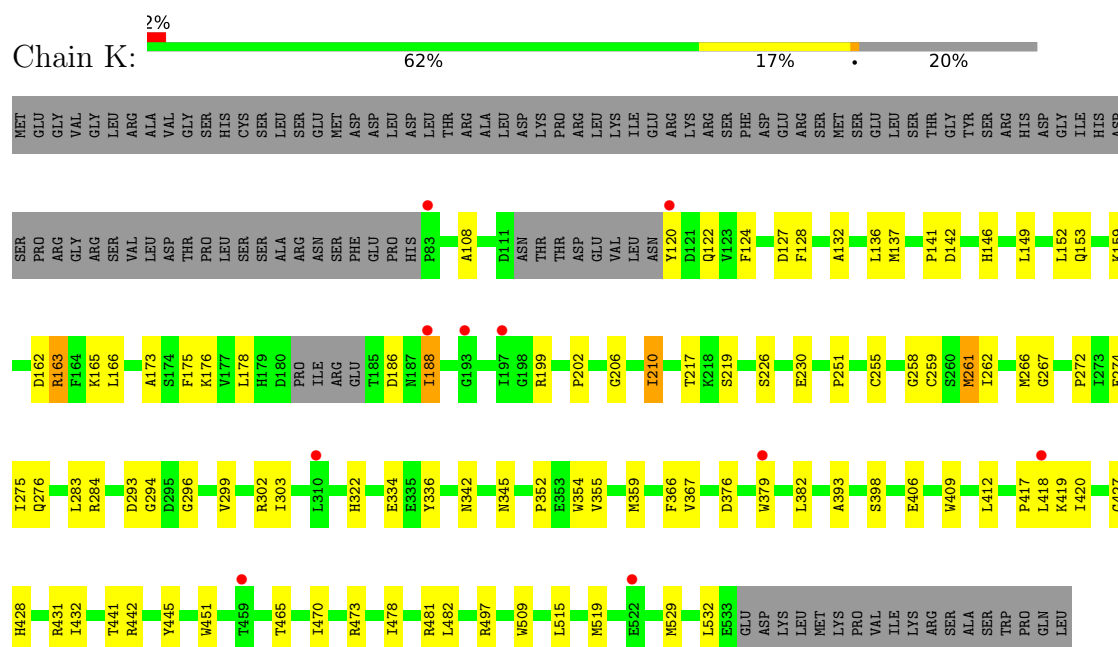


• Molecule 1: Alkaline/neutral invertase CINV1

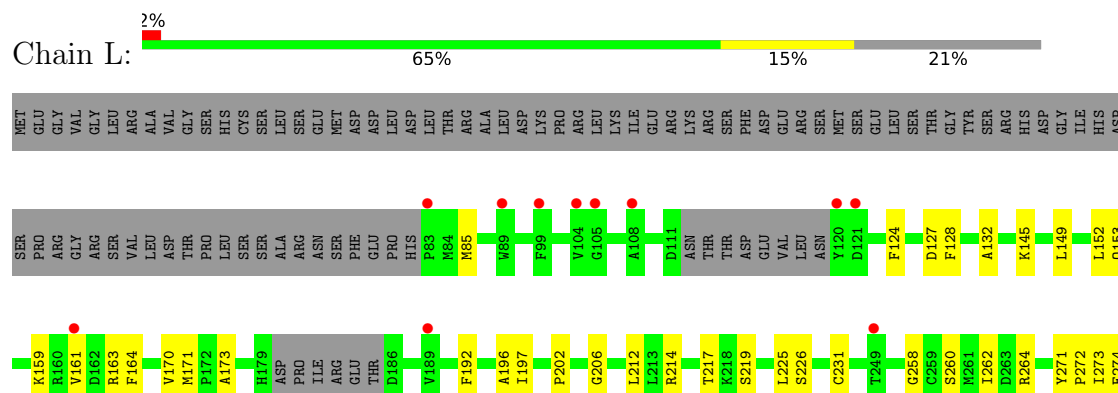


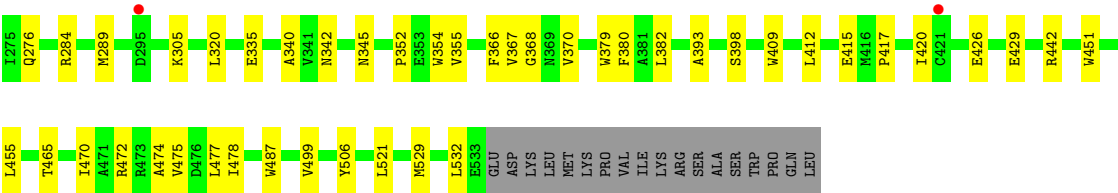


• Molecule 1: Alkaline/neutral invertase CINV1



• Molecule 1: Alkaline/neutral invertase CINV1





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	107.45Å 186.71Å 152.20Å 90.00° 104.86° 90.00°	Depositor
Resolution (Å)	46.65 – 3.39 46.65 – 3.39	Depositor EDS
% Data completeness (in resolution range)	98.7 (46.65-3.39) 98.8 (46.65-3.39)	Depositor EDS
R_{merge}	0.41	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.58 (at 3.40Å)	Xtriage
Refinement program	PHENIX 1.15.2_3472	Depositor
R, R_{free}	0.252 , 0.284 0.253 , 0.284	Depositor DCC
R_{free} test set	3969 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	62.3	Xtriage
Anisotropy	0.115	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 32.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.85	EDS
Total number of atoms	42174	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.86% 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	0.18	0/3559	0.56	3/4822 (0.1%)
1	B	0.17	0/3595	0.53	2/4871 (0.0%)
1	C	0.17	0/3634	0.52	2/4923 (0.0%)
1	D	0.16	0/3624	0.52	2/4910 (0.0%)
1	E	0.17	0/3607	0.53	2/4885 (0.0%)
1	F	0.17	0/3639	0.56	3/4931 (0.1%)
1	G	0.18	0/3634	0.56	4/4922 (0.1%)
1	H	0.19	0/3623	0.54	3/4908 (0.1%)
1	I	0.17	0/3574	0.55	0/4840
1	J	0.16	0/3632	0.55	0/4920
1	K	0.17	0/3620	0.54	0/4903
1	L	0.16	0/3599	0.52	2/4875 (0.0%)
All	All	0.17	0/43340	0.54	23/58710 (0.0%)

There are no bond length outliers.

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	161	VAL	CA-C-N	5.73	132.49	121.54
1	A	161	VAL	C-N-CA	5.73	132.49	121.54
1	G	85	MET	N-CA-C	-5.72	105.41	112.90
1	H	85	MET	N-CA-C	-5.53	105.66	112.90
1	A	84	MET	N-CA-C	5.43	126.19	111.00
1	E	161	VAL	CA-C-N	5.32	131.69	121.54
1	E	161	VAL	C-N-CA	5.32	131.69	121.54
1	D	161	VAL	CA-C-N	5.26	131.59	121.54
1	D	161	VAL	C-N-CA	5.26	131.59	121.54
1	H	161	VAL	CA-C-N	5.26	131.59	121.54
1	H	161	VAL	C-N-CA	5.26	131.59	121.54
1	G	161	VAL	CA-C-N	5.21	131.50	121.54
1	G	161	VAL	C-N-CA	5.21	131.50	121.54
1	G	188	ILE	CA-CB-CG1	5.07	119.02	110.40
1	B	161	VAL	CA-C-N	5.06	131.21	121.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	161	VAL	C-N-CA	5.06	131.21	121.54
1	F	82	HIS	C-N-CD	-5.05	109.49	120.60
1	C	161	VAL	CA-C-N	5.04	131.17	121.54
1	C	161	VAL	C-N-CA	5.04	131.17	121.54
1	F	82	HIS	CA-C-N	5.03	139.08	127.00
1	F	82	HIS	C-N-CA	5.03	139.08	127.00
1	L	161	VAL	CA-C-N	5.00	131.10	121.54
1	L	161	VAL	C-N-CA	5.00	131.10	121.54

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3464	0	3391	70	0
1	B	3499	0	3440	65	0
1	C	3536	0	3471	58	0
1	D	3526	0	3468	55	0
1	E	3511	0	3456	60	0
1	F	3540	0	3478	63	0
1	G	3536	0	3478	63	0
1	H	3525	0	3467	52	0
1	I	3478	0	3428	60	0
1	J	3534	0	3473	56	0
1	K	3523	0	3465	64	0
1	L	3502	0	3443	50	0
All	All	42174	0	41458	676	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 8.

All (676) 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:F:206:GLY:O	1:F:210:ILE:HD12	1.61	0.99
1:B:188:ILE:O	1:B:188:ILE:HD12	1.62	0.98
1:B:99:PHE:CE1	1:B:188:ILE:HG23	1.99	0.97
1:B:162:ASP:OD2	1:B:199:ARG:NH1	2.08	0.86
1:F:416:MET:HG3	1:F:417:PRO:HD2	1.57	0.86
1:K:342:ASN:HD21	1:K:345:ASN:HA	1.43	0.83
1:H:266:MET:H	1:H:267:GLY:HA2	1.42	0.82
1:A:152:LEU:HD11	1:A:209:TRP:CZ2	2.16	0.79
1:J:266:MET:H	1:J:267:GLY:HA2	1.48	0.78
1:I:266:MET:H	1:I:267:GLY:HA2	1.48	0.77
1:H:85:MET:HE2	1:H:518:LYS:HD3	1.64	0.77
1:K:266:MET:H	1:K:267:GLY:HA2	1.49	0.77
1:L:417:PRO:HD3	1:L:477:LEU:HD22	1.66	0.76
1:C:162:ASP:OD2	1:C:199:ARG:NH2	2.16	0.76
1:F:122:GLN:HG2	1:F:176:LYS:HB2	1.68	0.75
1:E:124:PHE:HB2	1:E:127:ASP:HB2	1.66	0.74
1:A:83:PRO:HA	1:A:85:MET:H	1.53	0.72
1:A:173:ALA:HB2	1:A:202:PRO:HG3	1.70	0.72
1:L:159:LYS:NZ	1:L:170:VAL:O	2.23	0.71
1:G:173:ALA:HB2	1:G:202:PRO:HG3	1.72	0.71
1:F:173:ALA:HB2	1:F:202:PRO:HG3	1.71	0.71
1:E:173:ALA:HB2	1:E:202:PRO:HG3	1.73	0.70
1:E:426:GLU:HG2	1:E:442:ARG:HG2	1.74	0.70
1:H:266:MET:N	1:H:267:GLY:HA2	2.07	0.70
1:K:334:GLU:OE2	1:L:264:ARG:NH2	2.19	0.70
1:H:426:GLU:HG2	1:H:442:ARG:HG2	1.73	0.69
1:E:195:SER:OG	1:F:338:HIS:CD2	2.45	0.69
1:C:334:GLU:OE2	1:D:197:ILE:HD13	1.93	0.69
1:F:168:GLU:HG2	1:F:237:LEU:HD22	1.74	0.69
1:C:221:GLY:O	1:E:472:ARG:NH1	2.26	0.69
1:G:346:VAL:HG21	1:G:375:MET:HE3	1.73	0.69
1:F:426:GLU:HG2	1:F:442:ARG:HG2	1.75	0.68
1:J:154:LEU:HD21	1:J:188:ILE:HD13	1.73	0.68
1:D:346:VAL:HG21	1:D:375:MET:HE3	1.74	0.68
1:C:173:ALA:HB2	1:C:202:PRO:HG3	1.75	0.68
1:K:342:ASN:ND2	1:K:345:ASN:HA	2.08	0.68
1:D:173:ALA:HB2	1:D:202:PRO:HG3	1.75	0.67
1:I:99:PHE:HB3	1:I:104:VAL:HG21	1.76	0.67
1:D:426:GLU:HG2	1:D:442:ARG:HG2	1.77	0.67
1:K:441:THR:HG22	1:K:442:ARG:H	1.59	0.67
1:H:173:ALA:HB2	1:H:202:PRO:HG3	1.77	0.67
1:C:102:GLN:HB3	1:C:146:HIS:HD2	1.60	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:426:GLU:HG2	1:B:442:ARG:HG2	1.77	0.66
1:H:346:VAL:HG11	1:H:375:MET:HE3	1.77	0.66
1:A:178:LEU:O	1:A:186:ASP:HA	1.95	0.66
1:I:426:GLU:HG2	1:I:442:ARG:HG2	1.78	0.65
1:I:346:VAL:HG21	1:I:375:MET:HE3	1.77	0.65
1:I:266:MET:N	1:I:267:GLY:HA2	2.12	0.65
1:J:83:PRO:HA	1:J:85:MET:N	2.12	0.65
1:H:515:LEU:O	1:H:519:MET:HG2	1.97	0.65
1:J:426:GLU:HG2	1:J:442:ARG:HG2	1.80	0.64
1:A:230:GLU:HA	1:A:233:LYS:HE3	1.80	0.64
1:H:124:PHE:HB2	1:H:127:ASP:HB2	1.80	0.64
1:F:126:ARG:NH1	1:F:204:ASP:OD2	2.31	0.63
1:L:455:LEU:HD22	1:L:478:ILE:HD13	1.81	0.63
1:E:308:HIS:C	1:F:163:ARG:HH12	2.06	0.63
1:A:214:ARG:HA	1:A:289:MET:HE1	1.81	0.62
1:F:417:PRO:HG3	1:F:478:ILE:HD11	1.80	0.62
1:J:99:PHE:CE2	1:J:188:ILE:HG23	2.34	0.62
1:J:124:PHE:HB2	1:J:127:ASP:HB2	1.81	0.62
1:J:409:TRP:HZ3	1:J:470:ILE:HG23	1.65	0.62
1:K:173:ALA:HB2	1:K:202:PRO:HG3	1.80	0.62
1:K:266:MET:N	1:K:267:GLY:HA2	2.15	0.62
1:A:82:HIS:N	1:A:85:MET:HG3	2.15	0.62
1:I:359:MET:HE1	1:I:364:GLY:HA2	1.81	0.62
1:B:415:GLU:HB3	1:B:481:ARG:HH12	1.64	0.62
1:C:124:PHE:HB2	1:C:127:ASP:HB2	1.82	0.62
1:B:359:MET:HE3	1:B:360:PRO:HD2	1.81	0.62
1:L:426:GLU:HG2	1:L:442:ARG:HG2	1.82	0.62
1:B:99:PHE:CD1	1:B:188:ILE:HG23	2.34	0.61
1:F:158:GLU:OE1	1:F:160:ARG:NH1	2.33	0.61
1:D:124:PHE:HB2	1:D:127:ASP:HB2	1.80	0.61
1:A:308:HIS:HB3	1:B:163:ARG:NH1	2.15	0.61
1:A:375:MET:SD	1:A:377:PHE:CZ	2.93	0.61
1:J:266:MET:N	1:J:267:GLY:HA2	2.14	0.61
1:D:342:ASN:ND2	1:D:345:ASN:HA	2.15	0.61
1:E:159:LYS:HB2	1:E:166:LEU:HB2	1.82	0.61
1:K:274:GLU:HG3	1:K:382:LEU:HD22	1.83	0.61
1:I:214:ARG:HA	1:I:289:MET:HE1	1.82	0.61
1:C:426:GLU:HG2	1:C:442:ARG:HG2	1.81	0.61
1:G:500:GLY:O	1:J:330:ARG:NH1	2.34	0.60
1:C:160:ARG:HD3	1:C:165:LYS:HG3	1.83	0.60
1:D:342:ASN:HD21	1:D:345:ASN:HA	1.67	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:260:SER:C	1:L:262:ILE:H	2.10	0.60
1:I:124:PHE:HB2	1:I:127:ASP:HB2	1.81	0.60
1:B:124:PHE:HB2	1:B:127:ASP:HB2	1.83	0.60
1:K:481:ARG:HH22	1:K:497:ARG:HA	1.67	0.60
1:K:142:ASP:O	1:K:146:HIS:ND1	2.35	0.59
1:B:159:LYS:HB2	1:B:166:LEU:HB2	1.84	0.59
1:J:274:GLU:HG3	1:J:382:LEU:HD22	1.84	0.59
1:I:162:ASP:OD1	1:I:199:ARG:NH1	2.36	0.59
1:I:158:GLU:OE1	1:I:160:ARG:NH1	2.35	0.59
1:F:162:ASP:O	1:F:163:ARG:HG3	2.03	0.59
1:I:108:ALA:HB1	1:I:120:TYR:HD1	1.67	0.59
1:L:124:PHE:HB2	1:L:127:ASP:HB2	1.83	0.59
1:K:178:LEU:O	1:K:186:ASP:HA	2.03	0.58
1:A:217:THR:HG21	1:A:226:SER:HB3	1.83	0.58
1:G:274:GLU:HG3	1:G:382:LEU:HD22	1.85	0.58
1:E:416:MET:HE2	1:E:491:TYR:CD1	2.38	0.58
1:I:173:ALA:HB2	1:I:202:PRO:HG3	1.83	0.58
1:A:426:GLU:HG2	1:A:442:ARG:HG2	1.84	0.58
1:K:124:PHE:HB2	1:K:127:ASP:HB2	1.84	0.58
1:B:173:ALA:HB2	1:B:202:PRO:HG3	1.85	0.58
1:A:308:HIS:HB3	1:B:163:ARG:HH12	1.67	0.58
1:C:478:ILE:HG23	1:C:482:LEU:HD12	1.86	0.58
1:A:162:ASP:OD2	1:A:199:ARG:NH2	2.32	0.58
1:A:335:GLU:OE2	1:A:340:ALA:HB1	2.04	0.58
1:K:188:ILE:HD12	1:K:188:ILE:O	2.04	0.58
1:B:346:VAL:HG21	1:B:375:MET:HE3	1.86	0.58
1:H:159:LYS:HB2	1:H:166:LEU:HB2	1.86	0.58
1:F:124:PHE:HB2	1:F:127:ASP:HB2	1.85	0.57
1:G:124:PHE:HB2	1:G:127:ASP:HB2	1.86	0.57
1:A:124:PHE:HB2	1:A:127:ASP:HB2	1.86	0.57
1:L:214:ARG:HA	1:L:289:MET:HE1	1.84	0.57
1:L:173:ALA:HB2	1:L:202:PRO:HG3	1.86	0.57
1:J:122:GLN:HE21	1:J:176:LYS:HG2	1.69	0.57
1:J:417:PRO:HD2	1:J:478:ILE:HG13	1.85	0.57
1:L:274:GLU:HG3	1:L:382:LEU:HD22	1.87	0.57
1:A:293:ASP:OD1	1:A:294:GLY:N	2.38	0.56
1:C:82:HIS:CE1	1:C:84:MET:HB2	2.40	0.56
1:B:188:ILE:HD12	1:B:188:ILE:C	2.29	0.56
1:B:100:ARG:NH2	1:B:186:ASP:O	2.38	0.56
1:C:154:LEU:HD21	1:C:188:ILE:HG21	1.87	0.56
1:D:255:CYS:HB3	1:D:259:CYS:SG	2.46	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:83:PRO:HA	1:F:85:MET:H	1.70	0.56
1:G:159:LYS:HB2	1:G:166:LEU:HB2	1.86	0.56
1:H:160:ARG:HH11	1:H:165:LYS:HE2	1.70	0.56
1:J:173:ALA:HB2	1:J:202:PRO:HG3	1.88	0.56
1:D:284:ARG:HB3	1:D:532:LEU:HD22	1.87	0.56
1:F:269:TYR:O	1:F:369:ASN:ND2	2.36	0.56
1:F:122:GLN:HB3	1:F:175:PHE:O	2.06	0.56
1:A:159:LYS:HB2	1:A:166:LEU:HB2	1.87	0.56
1:K:122:GLN:HG2	1:K:176:LYS:HB3	1.88	0.56
1:A:416:MET:HE2	1:A:491:TYR:CD1	2.41	0.56
1:B:188:ILE:O	1:B:188:ILE:CD1	2.48	0.56
1:F:455:LEU:HD22	1:F:478:ILE:HD12	1.88	0.56
1:J:367:VAL:HG21	1:J:375:MET:HE2	1.87	0.55
1:K:283:LEU:HD22	1:K:303:ILE:HG23	1.88	0.55
1:E:197:ILE:HG13	1:F:336:TYR:HD1	1.71	0.55
1:K:284:ARG:HH21	1:K:532:LEU:HB3	1.72	0.55
1:H:284:ARG:HH21	1:H:532:LEU:HB3	1.72	0.55
1:I:409:TRP:HZ3	1:I:470:ILE:HG23	1.72	0.55
1:K:165:LYS:HE3	1:L:305:LYS:HD3	1.88	0.55
1:I:417:PRO:HD2	1:I:478:ILE:HG13	1.89	0.54
1:K:162:ASP:OD2	1:K:199:ARG:NH2	2.37	0.54
1:L:258:GLY:HA2	1:L:262:ILE:O	2.07	0.54
1:F:284:ARG:HH21	1:F:532:LEU:HB3	1.73	0.54
1:F:416:MET:CG	1:F:417:PRO:HD2	2.31	0.54
1:K:137:MET:HB3	1:K:519:MET:HE2	1.89	0.54
1:K:137:MET:CB	1:K:519:MET:HE2	2.37	0.54
1:K:427:GLY:H	1:K:442:ARG:HH21	1.55	0.54
1:D:108:ALA:HB1	1:D:120:TYR:HD1	1.71	0.54
1:A:99:PHE:HB3	1:A:104:VAL:HG21	1.90	0.54
1:I:274:GLU:HG3	1:I:382:LEU:HD22	1.89	0.54
1:L:284:ARG:HH21	1:L:532:LEU:HB3	1.73	0.54
1:L:472:ARG:HG2	1:L:521:LEU:HD11	1.90	0.54
1:A:229:PRO:O	1:A:233:LYS:HG2	2.07	0.54
1:A:299:VAL:HG22	1:A:302:ARG:HH21	1.73	0.54
1:K:406:GLU:HG2	1:K:473:ARG:HH12	1.73	0.54
1:I:162:ASP:HB3	1:J:312:PHE:HE2	1.73	0.54
1:G:352:PRO:HG2	1:G:355:VAL:HG23	1.90	0.53
1:H:149:LEU:O	1:H:153:GLN:HG3	2.09	0.53
1:H:136:LEU:HD23	1:H:141:PRO:HB3	1.90	0.53
1:A:326:ASN:HD21	1:L:499:VAL:H	1.56	0.53
1:I:487:TRP:CE3	1:I:511:ILE:HG12	2.43	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:473:ARG:HH12	1:G:472:ARG:HD2	1.72	0.53
1:C:487:TRP:CE3	1:C:511:ILE:HG12	2.44	0.53
1:J:84:MET:SD	1:J:84:MET:N	2.81	0.53
1:K:149:LEU:O	1:K:153:GLN:HG3	2.08	0.53
1:G:284:ARG:HH21	1:G:532:LEU:HB3	1.73	0.53
1:I:283:LEU:HD22	1:I:303:ILE:HG23	1.90	0.53
1:C:274:GLU:HG3	1:C:382:LEU:HD22	1.90	0.53
1:G:85:MET:HE2	1:G:518:LYS:HD3	1.89	0.53
1:H:178:LEU:HD12	1:H:187:ASN:HB2	1.91	0.53
1:K:293:ASP:OD1	1:K:294:GLY:N	2.36	0.53
1:F:417:PRO:HG2	1:F:452:PRO:HA	1.91	0.53
1:H:274:GLU:HG3	1:H:382:LEU:HD22	1.90	0.53
1:B:122:GLN:HB3	1:B:176:LYS:HB3	1.90	0.53
1:D:419:LYS:NZ	1:D:494:LYS:O	2.41	0.53
1:B:274:GLU:HG3	1:B:382:LEU:HD23	1.91	0.53
1:G:409:TRP:HZ3	1:G:470:ILE:HG23	1.74	0.53
1:K:159:LYS:HB2	1:K:166:LEU:HB2	1.91	0.52
1:L:149:LEU:O	1:L:153:GLN:HG3	2.09	0.52
1:F:83:PRO:HA	1:F:85:MET:N	2.25	0.52
1:K:352:PRO:HG2	1:K:355:VAL:HG23	1.91	0.52
1:C:352:PRO:HG2	1:C:355:VAL:HG23	1.90	0.52
1:E:297:ARG:O	1:E:300:ILE:HG22	2.09	0.52
1:D:352:PRO:HG2	1:D:355:VAL:HG23	1.91	0.52
1:I:284:ARG:HH21	1:I:532:LEU:HB3	1.74	0.52
1:A:206:GLY:O	1:A:210:ILE:HD12	2.09	0.52
1:C:84:MET:SD	1:C:84:MET:N	2.82	0.52
1:C:160:ARG:HG2	1:C:165:LYS:HA	1.91	0.52
1:E:274:GLU:HG3	1:E:382:LEU:HD22	1.90	0.52
1:I:149:LEU:O	1:I:153:GLN:HG3	2.08	0.52
1:B:419:LYS:HA	1:B:450:SER:HA	1.91	0.52
1:I:145:LYS:HG3	1:I:225:LEU:HD22	1.91	0.52
1:A:260:SER:OG	1:A:274:GLU:OE1	2.25	0.52
1:F:273:ILE:HG12	1:F:368:GLY:HA2	1.91	0.52
1:G:82:HIS:CD2	1:G:85:MET:HB2	2.45	0.52
1:E:149:LEU:O	1:E:153:GLN:HG3	2.10	0.52
1:C:272:PRO:O	1:C:276:GLN:HG2	2.10	0.52
1:E:352:PRO:HG2	1:E:355:VAL:HG23	1.92	0.52
1:H:188:ILE:HD12	1:H:188:ILE:O	2.09	0.52
1:I:161:VAL:HG23	1:I:166:LEU:HD11	1.91	0.52
1:I:352:PRO:HG2	1:I:355:VAL:HG23	1.91	0.51
1:J:159:LYS:HB2	1:J:166:LEU:HB2	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:283:LEU:HD22	1:G:303:ILE:HG23	1.91	0.51
1:H:258:GLY:HA2	1:H:262:ILE:O	2.10	0.51
1:K:136:LEU:HD23	1:K:141:PRO:HB3	1.93	0.51
1:E:263:ASP:OD1	1:E:263:ASP:N	2.42	0.51
1:C:159:LYS:HB2	1:C:166:LEU:HB2	1.92	0.51
1:C:417:PRO:HD2	1:C:478:ILE:HG13	1.93	0.51
1:D:149:LEU:O	1:D:153:GLN:HG3	2.10	0.51
1:J:352:PRO:HG2	1:J:355:VAL:HG23	1.93	0.51
1:A:265:ARG:HB3	1:B:373:ALA:HB3	1.93	0.51
1:B:283:LEU:HD22	1:B:303:ILE:HG23	1.92	0.51
1:D:104:VAL:O	1:D:147:PHE:HB2	2.11	0.51
1:D:409:TRP:HZ3	1:D:470:ILE:HG23	1.76	0.51
1:E:283:LEU:HD22	1:E:303:ILE:HG23	1.93	0.51
1:I:255:CYS:HB3	1:I:259:CYS:SG	2.51	0.51
1:F:126:ARG:HH12	1:F:204:ASP:CG	2.19	0.50
1:F:272:PRO:O	1:F:276:GLN:HG2	2.10	0.50
1:C:255:CYS:HB3	1:C:259:CYS:SG	2.51	0.50
1:G:352:PRO:HB2	1:G:354:TRP:CD1	2.46	0.50
1:J:142:ASP:O	1:J:146:HIS:ND1	2.44	0.50
1:K:366:PHE:HE2	1:K:393:ALA:HB2	1.76	0.50
1:D:148:LEU:HD13	1:D:213:LEU:HD12	1.92	0.50
1:E:126:ARG:H	1:E:126:ARG:HD2	1.76	0.50
1:K:417:PRO:HD2	1:K:478:ILE:HG13	1.93	0.50
1:L:475:VAL:HA	1:L:478:ILE:HG22	1.92	0.50
1:B:99:PHE:CE2	1:B:150:LYS:HE3	2.47	0.50
1:B:523:ASP:OD1	1:B:525:SER:OG	2.26	0.50
1:D:272:PRO:O	1:D:276:GLN:HG2	2.12	0.50
1:E:195:SER:OG	1:F:338:HIS:NE2	2.44	0.50
1:G:262:ILE:HG23	1:G:446:HIS:CG	2.46	0.50
1:G:281:MET:HE3	1:G:386:VAL:HG13	1.92	0.50
1:L:219:SER:HB2	1:L:529:MET:HE2	1.93	0.50
1:B:455:LEU:HD22	1:B:478:ILE:HD12	1.93	0.50
1:C:148:LEU:HD13	1:C:213:LEU:HD12	1.92	0.50
1:A:272:PRO:O	1:A:276:GLN:HG2	2.10	0.50
1:B:415:GLU:HB3	1:B:481:ARG:NH1	2.26	0.50
1:G:298:GLU:HG2	1:G:302:ARG:HH21	1.75	0.50
1:J:272:PRO:O	1:J:276:GLN:HG2	2.11	0.50
1:A:352:PRO:HB2	1:A:354:TRP:CD1	2.47	0.50
1:E:100:ARG:HG3	1:E:186:ASP:HB2	1.93	0.50
1:G:272:PRO:O	1:G:276:GLN:HG2	2.11	0.50
1:K:272:PRO:O	1:K:276:GLN:HG2	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:99:PHE:CZ	1:C:188:ILE:HG12	2.46	0.50
1:J:255:CYS:HB3	1:J:259:CYS:SG	2.52	0.49
1:D:159:LYS:HB2	1:D:166:LEU:HB2	1.93	0.49
1:E:255:CYS:HB3	1:E:259:CYS:SG	2.51	0.49
1:A:82:HIS:CD2	1:A:83:PRO:O	2.65	0.49
1:B:272:PRO:O	1:B:276:GLN:HG2	2.12	0.49
1:C:366:PHE:HE2	1:C:393:ALA:HB2	1.77	0.49
1:D:487:TRP:CE3	1:D:511:ILE:HG12	2.47	0.49
1:E:142:ASP:O	1:E:146:HIS:ND1	2.44	0.49
1:F:82:HIS:HD2	1:F:479:GLU:CD	2.20	0.49
1:G:366:PHE:HE2	1:G:393:ALA:HB2	1.77	0.49
1:F:85:MET:HE1	1:F:522:GLU:OE2	2.12	0.49
1:I:136:LEU:HD23	1:I:141:PRO:HB3	1.93	0.49
1:I:515:LEU:O	1:I:519:MET:HG3	2.11	0.49
1:K:293:ASP:H	1:K:296:GLY:HA3	1.77	0.49
1:A:274:GLU:HG3	1:A:382:LEU:HD22	1.93	0.49
1:G:265:ARG:HB2	1:H:373:ALA:HB3	1.94	0.49
1:I:272:PRO:O	1:I:276:GLN:HG2	2.11	0.49
1:I:366:PHE:HE2	1:I:393:ALA:HB2	1.78	0.49
1:K:219:SER:HB2	1:K:529:MET:HE2	1.94	0.49
1:K:293:ASP:O	1:K:294:GLY:C	2.55	0.49
1:L:352:PRO:HG2	1:L:355:VAL:HG23	1.94	0.49
1:A:219:SER:HB2	1:A:529:MET:HE2	1.94	0.49
1:H:352:PRO:HG2	1:H:355:VAL:HG23	1.95	0.49
1:J:352:PRO:HB2	1:J:354:TRP:CD1	2.48	0.49
1:D:352:PRO:HB2	1:D:354:TRP:CD1	2.48	0.49
1:E:258:GLY:HA2	1:E:262:ILE:O	2.13	0.49
1:E:420:ILE:HB	1:E:451:TRP:CD1	2.47	0.49
1:F:82:HIS:HB2	1:F:83:PRO:O	2.12	0.49
1:J:226:SER:HB3	1:J:289:MET:HE2	1.94	0.49
1:L:417:PRO:HD3	1:L:477:LEU:CD2	2.41	0.49
1:D:418:LEU:HD23	1:D:419:LYS:N	2.28	0.49
1:E:272:PRO:O	1:E:276:GLN:HG2	2.13	0.49
1:I:379:TRP:CE3	1:I:412:LEU:HD22	2.48	0.49
1:A:455:LEU:HD22	1:A:478:ILE:HD12	1.95	0.48
1:G:165:LYS:NZ	1:H:305:LYS:HD3	2.28	0.48
1:G:298:GLU:OE1	1:G:302:ARG:NH2	2.45	0.48
1:K:417:PRO:HA	1:K:418:LEU:HA	1.62	0.48
1:L:217:THR:HG21	1:L:226:SER:HB3	1.94	0.48
1:A:417:PRO:HA	1:A:418:LEU:HA	1.61	0.48
1:C:482:LEU:HD22	1:C:487:TRP:CE3	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:416:MET:HE2	1:F:491:TYR:CE1	2.48	0.48
1:I:352:PRO:HB2	1:I:354:TRP:CD1	2.48	0.48
1:K:428:HIS:HA	1:K:431:ARG:HD2	1.95	0.48
1:B:409:TRP:HZ3	1:B:470:ILE:HG23	1.78	0.48
1:C:211:ILE:HG12	1:C:281:MET:HG2	1.94	0.48
1:C:283:LEU:HD22	1:C:303:ILE:HG23	1.95	0.48
1:D:366:PHE:HE2	1:D:393:ALA:HB2	1.78	0.48
1:K:108:ALA:HB1	1:K:120:TYR:HD1	1.78	0.48
1:L:352:PRO:HB2	1:L:354:TRP:CD1	2.48	0.48
1:D:136:LEU:HD23	1:D:141:PRO:HB3	1.94	0.48
1:G:273:ILE:HD11	1:G:380:PHE:HB2	1.95	0.48
1:K:352:PRO:HB2	1:K:354:TRP:CD1	2.48	0.48
1:B:149:LEU:O	1:B:153:GLN:HG3	2.12	0.48
1:D:473:ARG:O	1:D:476:ASP:OD1	2.31	0.48
1:E:85:MET:SD	1:E:518:LYS:HD3	2.54	0.48
1:E:367:VAL:HG22	1:E:376:ASP:O	2.13	0.48
1:E:342:ASN:ND2	1:E:345:ASN:HA	2.28	0.48
1:F:162:ASP:OD2	1:F:199:ARG:NH1	2.31	0.48
1:H:366:PHE:HE2	1:H:393:ALA:HB2	1.78	0.48
1:I:467:ARG:HB3	1:I:470:ILE:HD13	1.96	0.48
1:L:272:PRO:O	1:L:276:GLN:HG2	2.13	0.48
1:C:99:PHE:CE2	1:C:177:VAL:HG13	2.48	0.48
1:C:409:TRP:HZ3	1:C:470:ILE:HG23	1.79	0.48
1:D:518:LYS:O	1:D:522:GLU:HG2	2.14	0.48
1:F:149:LEU:O	1:F:153:GLN:HG3	2.13	0.48
1:D:367:VAL:HG13	1:D:376:ASP:O	2.14	0.48
1:E:136:LEU:HD23	1:E:141:PRO:HB3	1.94	0.48
1:J:149:LEU:O	1:J:153:GLN:HG3	2.13	0.48
1:L:379:TRP:CE3	1:L:412:LEU:HD22	2.49	0.48
1:J:261:MET:HE2	1:J:420:ILE:HD12	1.96	0.48
1:K:255:CYS:HB3	1:K:259:CYS:SG	2.54	0.48
1:L:487:TRP:HB2	1:L:506:TYR:HA	1.95	0.48
1:J:219:SER:HB2	1:J:529:MET:HE2	1.96	0.47
1:C:406:GLU:OE2	1:G:469:GLN:NE2	2.47	0.47
1:I:481:ARG:NH2	1:I:496:GLY:O	2.45	0.47
1:J:136:LEU:HD23	1:J:141:PRO:HB3	1.97	0.47
1:B:352:PRO:HG2	1:B:355:VAL:HG23	1.94	0.47
1:G:367:VAL:HG13	1:G:376:ASP:O	2.15	0.47
1:J:379:TRP:CE3	1:J:412:LEU:HD22	2.49	0.47
1:C:195:SER:HG	1:D:338:HIS:CE1	2.32	0.47
1:C:262:ILE:HG23	1:C:446:HIS:CG	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:417:PRO:HA	1:C:418:LEU:HA	1.64	0.47
1:E:352:PRO:HB2	1:E:354:TRP:CD1	2.49	0.47
1:G:455:LEU:HD22	1:G:478:ILE:HD13	1.97	0.47
1:J:177:VAL:HA	1:J:187:ASN:O	2.15	0.47
1:J:366:PHE:HE2	1:J:393:ALA:HB2	1.79	0.47
1:B:136:LEU:HG	1:B:144:VAL:HG21	1.97	0.47
1:D:188:ILE:HD12	1:D:188:ILE:O	2.13	0.47
1:G:411:GLU:N	1:G:411:GLU:OE1	2.47	0.47
1:L:409:TRP:HZ3	1:L:470:ILE:HG23	1.80	0.47
1:B:219:SER:HB2	1:B:529:MET:HE2	1.97	0.47
1:B:315:ARG:NE	1:B:392:LEU:O	2.26	0.47
1:C:128:PHE:CE2	1:C:132:ALA:HB2	2.50	0.47
1:E:261:MET:HE3	1:E:509:TRP:CH2	2.49	0.47
1:B:352:PRO:HB2	1:B:354:TRP:CD1	2.49	0.47
1:E:455:LEU:HD22	1:E:478:ILE:HD12	1.97	0.47
1:F:162:ASP:CG	1:F:199:ARG:HH12	2.22	0.47
1:F:299:VAL:HG22	1:F:302:ARG:HH21	1.80	0.47
1:I:188:ILE:HD12	1:I:188:ILE:O	2.14	0.47
1:G:436:CYS:HB2	1:H:350:SER:HB3	1.95	0.47
1:H:272:PRO:O	1:H:276:GLN:HG2	2.15	0.47
1:L:415:GLU:C	1:L:477:LEU:HD21	2.40	0.47
1:A:352:PRO:HG2	1:A:355:VAL:HG23	1.95	0.47
1:A:373:ALA:HB3	1:B:265:ARG:HB2	1.97	0.47
1:K:478:ILE:HG23	1:K:482:LEU:HD12	1.97	0.47
1:A:83:PRO:HA	1:A:85:MET:N	2.26	0.47
1:A:262:ILE:HG23	1:A:446:HIS:CG	2.50	0.47
1:C:195:SER:OG	1:D:338:HIS:CE1	2.68	0.47
1:C:482:LEU:HD23	1:C:487:TRP:HA	1.97	0.47
1:A:478:ILE:HG23	1:A:482:LEU:HD12	1.97	0.46
1:B:258:GLY:HA2	1:B:262:ILE:O	2.15	0.46
1:D:274:GLU:HG3	1:D:382:LEU:HD22	1.98	0.46
1:H:352:PRO:HB2	1:H:354:TRP:CD1	2.50	0.46
1:I:207:PHE:CD2	1:I:279:PHE:HA	2.49	0.46
1:L:474:ALA:O	1:L:478:ILE:HG22	2.15	0.46
1:B:85:MET:SD	1:B:518:LYS:HD3	2.56	0.46
1:F:145:LYS:HG3	1:F:225:LEU:HD22	1.98	0.46
1:H:417:PRO:HA	1:H:418:LEU:HA	1.61	0.46
1:K:515:LEU:O	1:K:519:MET:HG2	2.15	0.46
1:L:85:MET:HE3	1:L:85:MET:HB3	1.79	0.46
1:E:308:HIS:O	1:F:163:ARG:NH1	2.45	0.46
1:G:165:LYS:HZ2	1:H:305:LYS:HD3	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:342:ASN:ND2	1:I:345:ASN:HA	2.30	0.46
1:B:188:ILE:C	1:B:188:ILE:CD1	2.88	0.46
1:I:159:LYS:HB2	1:I:166:LEU:HB2	1.98	0.46
1:J:179:HIS:HA	1:J:185:THR:HA	1.96	0.46
1:E:366:PHE:HE2	1:E:393:ALA:HB2	1.81	0.46
1:F:255:CYS:HB3	1:F:259:CYS:SG	2.56	0.46
1:G:308:HIS:HB3	1:H:163:ARG:NH1	2.31	0.46
1:H:354:TRP:CZ2	1:H:423:PRO:HD3	2.51	0.46
1:B:128:PHE:CE2	1:B:132:ALA:HB2	2.50	0.46
1:E:265:ARG:HB2	1:F:373:ALA:HB3	1.97	0.46
1:G:478:ILE:C	1:G:480:SER:H	2.24	0.46
1:K:122:GLN:HB3	1:K:175:PHE:O	2.16	0.46
1:H:99:PHE:HB3	1:H:104:VAL:HG21	1.97	0.46
1:H:99:PHE:CE1	1:H:188:ILE:HG23	2.51	0.46
1:I:207:PHE:CE2	1:I:279:PHE:HA	2.51	0.46
1:A:283:LEU:HD22	1:A:303:ILE:HG23	1.98	0.46
1:F:126:ARG:NH1	1:F:126:ARG:HB3	2.31	0.46
1:H:455:LEU:HD22	1:H:478:ILE:HD13	1.98	0.46
1:L:260:SER:O	1:L:262:ILE:HD12	2.15	0.46
1:A:501:LYS:NZ	1:C:349:ASP:OD2	2.49	0.46
1:B:260:SER:C	1:B:262:ILE:N	2.73	0.46
1:G:194:GLU:HG3	1:H:338:HIS:CD2	2.51	0.46
1:G:398:SER:HB3	1:G:465:THR:HG22	1.98	0.46
1:E:487:TRP:CD2	1:E:511:ILE:HG12	2.50	0.45
1:F:159:LYS:HB2	1:F:166:LEU:HB2	1.96	0.45
1:G:219:SER:HB2	1:G:529:MET:HE2	1.99	0.45
1:G:379:TRP:CE3	1:G:412:LEU:HD22	2.51	0.45
1:I:122:GLN:HG2	1:I:176:LYS:HD3	1.97	0.45
1:I:289:MET:HE2	1:I:289:MET:HB3	1.70	0.45
1:B:415:GLU:O	1:B:481:ARG:NH1	2.49	0.45
1:D:145:LYS:HG3	1:D:225:LEU:HD22	1.98	0.45
1:E:260:SER:C	1:E:262:ILE:N	2.74	0.45
1:G:194:GLU:HG2	1:G:195:SER:N	2.30	0.45
1:A:145:LYS:HG3	1:A:225:LEU:HD22	1.99	0.45
1:F:398:SER:HB3	1:F:465:THR:HG22	1.98	0.45
1:I:128:PHE:CE2	1:I:132:ALA:HB2	2.52	0.45
1:A:379:TRP:CE3	1:A:412:LEU:HD22	2.51	0.45
1:E:411:GLU:N	1:E:411:GLU:OE1	2.50	0.45
1:F:203:VAL:HG23	1:F:259:CYS:HB2	1.97	0.45
1:K:206:GLY:O	1:K:210:ILE:HD12	2.17	0.45
1:B:229:PRO:HG2	1:B:230:GLU:OE2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:487:TRP:CD2	1:C:511:ILE:HG12	2.51	0.45
1:F:379:TRP:CE3	1:F:412:LEU:HD22	2.50	0.45
1:H:255:CYS:HB3	1:H:259:CYS:SG	2.57	0.45
1:I:108:ALA:HB3	1:I:120:TYR:O	2.16	0.45
1:J:128:PHE:CE2	1:J:132:ALA:HB2	2.52	0.45
1:L:273:ILE:HG12	1:L:368:GLY:HA2	1.99	0.45
1:F:415:GLU:O	1:F:481:ARG:NH1	2.50	0.45
1:G:128:PHE:CE2	1:G:132:ALA:HB2	2.52	0.45
1:J:482:LEU:HD23	1:J:482:LEU:HA	1.82	0.45
1:D:283:LEU:HD22	1:D:303:ILE:HG23	1.99	0.45
1:D:417:PRO:HA	1:D:418:LEU:HA	1.63	0.45
1:F:283:LEU:HD22	1:F:303:ILE:HG23	1.98	0.45
1:G:322:HIS:N	1:G:359:MET:HE2	2.32	0.45
1:G:443:TRP:CD1	1:G:494:LYS:HG2	2.52	0.45
1:H:104:VAL:O	1:H:147:PHE:HB2	2.17	0.45
1:L:366:PHE:HE2	1:L:393:ALA:HB2	1.81	0.45
1:A:85:MET:HE3	1:A:85:MET:HB2	1.89	0.45
1:A:217:THR:HG21	1:A:226:SER:CB	2.47	0.45
1:G:468:PRO:O	1:G:472:ARG:HG3	2.17	0.45
1:I:171:MET:HG3	1:I:206:GLY:HA2	1.97	0.45
1:J:160:ARG:HD3	1:J:165:LYS:HG3	1.98	0.45
1:A:487:TRP:CD2	1:A:511:ILE:HG12	2.52	0.45
1:E:417:PRO:HD2	1:E:478:ILE:HG13	1.99	0.45
1:F:482:LEU:HA	1:F:482:LEU:HD23	1.80	0.45
1:H:482:LEU:HD23	1:H:482:LEU:HA	1.83	0.45
1:I:207:PHE:HE2	1:I:279:PHE:CD1	2.35	0.45
1:J:416:MET:HE2	1:J:491:TYR:CD1	2.52	0.45
1:D:132:ALA:HB3	1:D:212:LEU:HD11	1.99	0.44
1:D:420:ILE:HB	1:D:451:TRP:CD1	2.51	0.44
1:K:152:LEU:HD11	1:K:230:GLU:O	2.17	0.44
1:A:162:ASP:CG	1:A:199:ARG:HH12	2.25	0.44
1:A:260:SER:OG	1:A:261:MET:N	2.49	0.44
1:B:177:VAL:HG22	1:B:188:ILE:HG22	1.99	0.44
1:F:128:PHE:CE2	1:F:132:ALA:HB2	2.52	0.44
1:J:260:SER:C	1:J:262:ILE:N	2.75	0.44
1:K:255:CYS:SG	1:K:275:ILE:HG21	2.56	0.44
1:A:315:ARG:HH21	1:A:394:THR:HG22	1.83	0.44
1:A:482:LEU:HD23	1:A:482:LEU:HA	1.78	0.44
1:B:255:CYS:HB3	1:B:259:CYS:SG	2.57	0.44
1:C:157:TRP:HH2	1:C:188:ILE:HB	1.82	0.44
1:C:469:GLN:NE2	1:G:406:GLU:OE2	2.50	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:84:MET:HE1	1:D:479:GLU:HA	1.99	0.44
1:I:162:ASP:HB3	1:J:312:PHE:CE2	2.52	0.44
1:J:178:LEU:HB3	1:J:187:ASN:HB2	1.99	0.44
1:J:487:TRP:CD2	1:J:511:ILE:HG12	2.53	0.44
1:C:102:GLN:HB3	1:C:146:HIS:CD2	2.47	0.44
1:K:379:TRP:CE3	1:K:412:LEU:HD22	2.53	0.44
1:L:398:SER:HB3	1:L:465:THR:HG22	1.99	0.44
1:C:352:PRO:HB2	1:C:354:TRP:CD1	2.52	0.44
1:J:258:GLY:HA2	1:J:262:ILE:O	2.18	0.44
1:G:82:HIS:CD2	1:G:514:TYR:OH	2.71	0.44
1:I:320:LEU:HD13	1:I:359:MET:HE3	1.98	0.44
1:C:482:LEU:CD2	1:C:487:TRP:HA	2.47	0.44
1:G:152:LEU:HD11	1:G:230:GLU:O	2.17	0.44
1:J:342:ASN:ND2	1:J:345:ASN:HA	2.33	0.44
1:K:398:SER:HB3	1:K:465:THR:HG22	2.00	0.44
1:K:482:LEU:HD23	1:K:482:LEU:HA	1.79	0.44
1:L:271:TYR:HD2	1:L:370:VAL:O	2.01	0.44
1:C:367:VAL:HG13	1:C:376:ASP:O	2.17	0.44
1:H:379:TRP:CE3	1:H:412:LEU:HD22	2.53	0.44
1:C:85:MET:SD	1:C:518:LYS:HD3	2.58	0.44
1:D:290:LEU:HD23	1:D:300:ILE:HG13	1.99	0.44
1:K:261:MET:HE2	1:K:509:TRP:CH2	2.53	0.44
1:A:251:PRO:HB2	1:A:310:LEU:HD12	1.98	0.43
1:A:409:TRP:HZ3	1:A:470:ILE:HG23	1.83	0.43
1:C:99:PHE:CD2	1:C:177:VAL:HG13	2.53	0.43
1:E:322:HIS:N	1:E:359:MET:HE2	2.33	0.43
1:F:273:ILE:HD11	1:F:380:PHE:HB2	2.00	0.43
1:F:366:PHE:HE2	1:F:393:ALA:HB2	1.83	0.43
1:L:145:LYS:HG3	1:L:225:LEU:HD22	2.00	0.43
1:A:128:PHE:CE2	1:A:132:ALA:HB2	2.52	0.43
1:G:82:HIS:CD2	1:G:514:TYR:HH	2.36	0.43
1:I:229:PRO:HG2	1:I:230:GLU:OE2	2.18	0.43
1:I:370:VAL:HG22	1:I:375:MET:HG3	2.01	0.43
1:I:418:LEU:HD23	1:I:419:LYS:N	2.33	0.43
1:J:398:SER:HB3	1:J:465:THR:HG22	2.00	0.43
1:L:152:LEU:CD1	1:L:231:CYS:HA	2.48	0.43
1:B:160:ARG:HG2	1:B:165:LYS:HG3	2.00	0.43
1:B:320:LEU:HB2	1:B:367:VAL:HG23	1.99	0.43
1:D:455:LEU:HD22	1:D:478:ILE:HD12	2.01	0.43
1:B:214:ARG:HA	1:B:289:MET:HE1	2.00	0.43
1:D:411:GLU:OE1	1:D:411:GLU:N	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:342:ASN:ND2	1:H:345:ASN:HA	2.33	0.43
1:K:418:LEU:HD23	1:K:419:LYS:N	2.34	0.43
1:L:273:ILE:HD11	1:L:380:PHE:HB2	2.01	0.43
1:L:335:GLU:HB3	1:L:340:ALA:HB1	1.99	0.43
1:F:453:VAL:HG21	1:F:509:TRP:CD2	2.53	0.43
1:B:482:LEU:HD23	1:B:482:LEU:HA	1.76	0.43
1:C:322:HIS:N	1:C:359:MET:HE2	2.33	0.43
1:C:379:TRP:CE3	1:C:412:LEU:HD22	2.54	0.43
1:G:482:LEU:HD23	1:G:482:LEU:HA	1.80	0.43
1:C:97:VAL:HG22	1:C:107:LEU:HG	2.01	0.43
1:H:260:SER:C	1:H:262:ILE:N	2.76	0.43
1:I:162:ASP:CG	1:I:199:ARG:HH12	2.26	0.43
1:L:342:ASN:ND2	1:L:345:ASN:HA	2.34	0.43
1:E:103:PRO:HB2	1:E:143:ILE:HD13	1.99	0.43
1:E:171:MET:HG3	1:E:206:GLY:HA2	2.01	0.43
1:F:342:ASN:ND2	1:F:345:ASN:HA	2.34	0.43
1:G:342:ASN:ND2	1:G:345:ASN:HA	2.34	0.43
1:H:226:SER:HB3	1:H:289:MET:HE2	2.01	0.43
1:H:468:PRO:HB2	1:H:472:ARG:NH1	2.34	0.43
1:I:379:TRP:O	1:I:421:CYS:HA	2.19	0.43
1:K:336:TYR:CE1	1:L:196:ALA:HA	2.53	0.43
1:B:136:LEU:HD23	1:B:141:PRO:HB3	2.01	0.43
1:J:132:ALA:HB3	1:J:212:LEU:HD11	2.01	0.43
1:L:128:PHE:CE2	1:L:132:ALA:HB2	2.54	0.43
1:L:132:ALA:HB3	1:L:212:LEU:HD11	2.01	0.43
1:A:164:PHE:HB2	1:B:251:PRO:HG3	2.00	0.43
1:F:443:TRP:HE1	1:F:494:LYS:HE3	1.82	0.43
1:D:180:ASP:H	1:D:185:THR:HB	1.83	0.42
1:E:284:ARG:HH21	1:E:532:LEU:HB3	1.82	0.42
1:J:104:VAL:O	1:J:147:PHE:HB2	2.19	0.42
1:K:128:PHE:CE2	1:K:132:ALA:HB2	2.54	0.42
1:A:136:LEU:HD23	1:A:141:PRO:HB3	2.01	0.42
1:A:209:TRP:HE3	1:A:210:ILE:HG13	1.84	0.42
1:B:417:PRO:HA	1:B:418:LEU:HA	1.59	0.42
1:D:159:LYS:NZ	1:D:168:GLU:O	2.51	0.42
1:E:122:GLN:HG3	1:E:124:PHE:CE1	2.54	0.42
1:E:260:SER:O	1:E:262:ILE:N	2.52	0.42
1:K:367:VAL:HG13	1:K:376:ASP:O	2.19	0.42
1:C:203:VAL:HG12	1:C:207:PHE:HE2	1.83	0.42
1:D:450:SER:HB2	1:D:491:TYR:HB2	2.00	0.42
1:F:217:THR:HG21	1:F:226:SER:HB3	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:177:VAL:HG12	1:G:188:ILE:HG22	2.01	0.42
1:G:417:PRO:HA	1:G:418:LEU:HA	1.63	0.42
1:K:409:TRP:HZ3	1:K:470:ILE:HG23	1.83	0.42
1:A:366:PHE:HE2	1:A:393:ALA:HB2	1.83	0.42
1:G:487:TRP:CD2	1:G:511:ILE:HG12	2.54	0.42
1:A:312:PHE:C	1:A:312:PHE:CD1	2.96	0.42
1:C:432:ILE:HD12	1:D:432:ILE:HD12	2.01	0.42
1:D:255:CYS:SG	1:D:275:ILE:HG21	2.59	0.42
1:E:188:ILE:O	1:E:188:ILE:HD12	2.19	0.42
1:E:379:TRP:CE3	1:E:412:LEU:HD22	2.55	0.42
1:F:478:ILE:HG23	1:F:482:LEU:HD12	2.00	0.42
1:G:499:VAL:HG23	1:J:326:ASN:OD1	2.20	0.42
1:I:398:SER:HB3	1:I:465:THR:HG22	2.01	0.42
1:I:487:TRP:CD2	1:I:511:ILE:HG12	2.55	0.42
1:J:411:GLU:N	1:J:411:GLU:OE1	2.52	0.42
1:J:417:PRO:HA	1:J:418:LEU:HA	1.64	0.42
1:J:453:VAL:HG21	1:J:509:TRP:CD2	2.54	0.42
1:A:284:ARG:NH1	1:A:391:SER:HB2	2.35	0.42
1:B:487:TRP:CD2	1:B:511:ILE:HG12	2.54	0.42
1:E:291:LYS:O	1:E:296:GLY:HA3	2.19	0.42
1:E:515:LEU:O	1:E:519:MET:HG3	2.18	0.42
1:G:297:ARG:O	1:G:301:GLU:HG3	2.20	0.42
1:I:207:PHE:N	1:I:207:PHE:CD1	2.87	0.42
1:E:219:SER:HB2	1:E:529:MET:HE2	2.00	0.42
1:E:417:PRO:HA	1:E:418:LEU:HA	1.60	0.42
1:H:229:PRO:HB2	1:H:233:LYS:HE3	2.02	0.42
1:J:322:HIS:N	1:J:359:MET:HE2	2.34	0.42
1:K:162:ASP:O	1:K:163:ARG:HG3	2.20	0.42
1:K:432:ILE:HD13	1:L:429:GLU:HA	2.02	0.42
1:A:261:MET:HE2	1:A:509:TRP:CZ3	2.54	0.42
1:A:420:ILE:HB	1:A:451:TRP:CD1	2.55	0.42
1:B:132:ALA:HB3	1:B:212:LEU:HD11	2.01	0.42
1:B:394:THR:OG1	1:B:395:PRO:HD2	2.20	0.42
1:B:398:SER:HB3	1:B:465:THR:HG22	2.00	0.42
1:A:132:ALA:HB3	1:A:212:LEU:HD11	2.00	0.42
1:A:429:GLU:HA	1:B:432:ILE:HD13	2.02	0.42
1:B:478:ILE:HG23	1:B:482:LEU:HD12	2.01	0.42
1:D:260:SER:OG	1:D:261:MET:N	2.51	0.42
1:D:487:TRP:HB2	1:D:506:TYR:HA	2.01	0.42
1:E:398:SER:HB3	1:E:465:THR:HG22	2.02	0.42
1:F:132:ALA:HB3	1:F:212:LEU:HD11	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:409:TRP:CZ2	1:F:413:VAL:HG11	2.54	0.42
1:G:136:LEU:HD23	1:G:141:PRO:HB3	2.02	0.42
1:K:188:ILE:H	1:K:188:ILE:HG13	1.64	0.42
1:B:418:LEU:CD1	1:B:458:LEU:HD12	2.50	0.42
1:H:262:ILE:HD12	1:H:446:HIS:HB3	2.02	0.42
1:H:420:ILE:HB	1:H:451:TRP:CD1	2.55	0.42
1:J:283:LEU:HD22	1:J:303:ILE:HG23	2.00	0.42
1:D:151:THR:HB	1:D:171:MET:HE1	2.02	0.41
1:F:168:GLU:HG2	1:F:237:LEU:CD2	2.44	0.41
1:F:455:LEU:HD22	1:F:478:ILE:CD1	2.49	0.41
1:H:122:GLN:HG2	1:H:176:LYS:HB3	2.02	0.41
1:H:132:ALA:HB3	1:H:212:LEU:HD11	2.02	0.41
1:K:258:GLY:HA2	1:K:262:ILE:O	2.20	0.41
1:E:136:LEU:HG	1:E:144:VAL:HG21	2.01	0.41
1:G:255:CYS:SG	1:G:275:ILE:HG21	2.59	0.41
1:G:443:TRP:HE1	1:G:494:LYS:HE2	1.83	0.41
1:A:188:ILE:HD12	1:A:188:ILE:O	2.20	0.41
1:B:420:ILE:HB	1:B:451:TRP:CD1	2.55	0.41
1:C:411:GLU:N	1:C:411:GLU:OE1	2.53	0.41
1:F:148:LEU:HD13	1:F:213:LEU:HD12	2.01	0.41
1:F:219:SER:HB2	1:F:529:MET:HE2	2.02	0.41
1:H:219:SER:HB2	1:H:529:MET:HE2	2.02	0.41
1:K:322:HIS:N	1:K:359:MET:HE2	2.35	0.41
1:L:171:MET:HG3	1:L:206:GLY:HA2	2.02	0.41
1:B:378:ARG:HG2	1:B:423:PRO:HD2	2.01	0.41
1:D:260:SER:C	1:D:262:ILE:N	2.77	0.41
1:E:226:SER:HB3	1:E:289:MET:HE2	2.02	0.41
1:B:379:TRP:CE3	1:B:412:LEU:HD22	2.55	0.41
1:D:128:PHE:CE2	1:D:132:ALA:HB2	2.55	0.41
1:F:104:VAL:O	1:F:147:PHE:HB2	2.20	0.41
1:G:206:GLY:O	1:G:210:ILE:HD12	2.20	0.41
1:G:444:SER:OG	1:G:445:TYR:N	2.52	0.41
1:H:255:CYS:SG	1:H:275:ILE:HG21	2.60	0.41
1:I:257:ASP:OD1	1:I:265:ARG:HA	2.21	0.41
1:J:122:GLN:HG2	1:J:176:LYS:HD3	2.03	0.41
1:A:83:PRO:HB2	1:A:84:MET:SD	2.60	0.41
1:A:163:ARG:HD2	1:B:312:PHE:CB	2.51	0.41
1:A:260:SER:C	1:A:262:ILE:N	2.77	0.41
1:B:453:VAL:HG21	1:B:509:TRP:CE3	2.55	0.41
1:D:453:VAL:HG21	1:D:509:TRP:CD2	2.56	0.41
1:E:255:CYS:SG	1:E:275:ILE:HG21	2.60	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:99:PHE:CE2	1:I:150:LYS:HE3	2.55	0.41
1:K:217:THR:HG21	1:K:226:SER:HB3	2.03	0.41
1:A:258:GLY:HA2	1:A:262:ILE:O	2.20	0.41
1:B:104:VAL:O	1:B:147:PHE:HB2	2.21	0.41
1:E:418:LEU:HD23	1:E:419:LYS:N	2.35	0.41
1:G:150:LYS:HD2	1:G:150:LYS:HA	1.81	0.41
1:G:503:ALA:HB3	1:J:330:ARG:HH12	1.86	0.41
1:H:320:LEU:HB2	1:H:367:VAL:HG23	2.02	0.41
1:I:469:GLN:H	1:I:469:GLN:CD	2.28	0.41
1:B:124:PHE:CD2	1:B:174:SER:HB3	2.55	0.41
1:D:322:HIS:N	1:D:359:MET:HE2	2.36	0.41
1:F:420:ILE:HB	1:F:451:TRP:CD1	2.56	0.41
1:H:418:LEU:HD23	1:H:419:LYS:N	2.35	0.41
1:K:251:PRO:HG3	1:L:164:PHE:HB2	2.02	0.41
1:K:420:ILE:HB	1:K:451:TRP:CD1	2.56	0.41
1:A:209:TRP:CE3	1:A:210:ILE:HG13	2.56	0.41
1:B:342:ASN:ND2	1:B:345:ASN:HA	2.36	0.41
1:C:398:SER:HB3	1:C:465:THR:HG22	2.02	0.41
1:E:132:ALA:HB3	1:E:212:LEU:HD11	2.03	0.41
1:F:367:VAL:HG13	1:F:376:ASP:O	2.21	0.41
1:I:417:PRO:HA	1:I:418:LEU:HA	1.68	0.41
1:J:83:PRO:HA	1:J:84:MET:C	2.46	0.41
1:J:85:MET:HE2	1:J:85:MET:HB3	1.84	0.41
1:L:159:LYS:HE2	1:L:192:PHE:CE2	2.56	0.41
1:D:99:PHE:HB3	1:D:104:VAL:CG2	2.51	0.41
1:G:214:ARG:HA	1:G:289:MET:HE1	2.03	0.41
1:L:320:LEU:HB2	1:L:367:VAL:HG23	2.02	0.41
1:C:281:MET:HB2	1:C:390:SER:HB3	2.03	0.40
1:D:99:PHE:HB3	1:D:104:VAL:HG21	2.02	0.40
1:D:298:GLU:H	1:D:298:GLU:HG2	1.72	0.40
1:E:85:MET:HG2	1:E:514:TYR:HE2	1.85	0.40
1:H:519:MET:HE2	1:H:519:MET:HB3	1.88	0.40
1:I:160:ARG:HG3	1:I:165:LYS:HG3	2.03	0.40
1:I:406:GLU:OE2	1:I:470:ILE:HD11	2.21	0.40
1:I:468:PRO:HB2	1:I:472:ARG:NH1	2.36	0.40
1:A:151:THR:HB	1:A:171:MET:HE1	2.03	0.40
1:A:422:TYR:HA	1:A:423:PRO:HA	1.90	0.40
1:C:162:ASP:CG	1:C:199:ARG:HH12	2.29	0.40
1:E:104:VAL:O	1:E:147:PHE:HB2	2.21	0.40
1:G:444:SER:O	1:G:449:GLY:N	2.55	0.40
1:G:455:LEU:HD22	1:G:478:ILE:CD1	2.51	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:299:VAL:HG22	1:K:302:ARG:HH21	1.86	0.40
1:A:169:GLY:C	1:A:241:LEU:HD11	2.45	0.40
1:E:478:ILE:HG23	1:E:482:LEU:HD12	2.03	0.40
1:H:411:GLU:OE1	1:H:411:GLU:N	2.54	0.40
1:J:206:GLY:O	1:J:210:ILE:HD12	2.21	0.40
1:J:420:ILE:HB	1:J:451:TRP:CD1	2.56	0.40
1:K:137:MET:HB2	1:K:519:MET:HE2	2.02	0.40
1:L:289:MET:HE2	1:L:289:MET:HB3	1.74	0.40
1:L:420:ILE:HB	1:L:451:TRP:CD1	2.56	0.40
1:C:160:ARG:HD3	1:C:165:LYS:HE2	2.04	0.40
1:C:260:SER:C	1:C:262:ILE:N	2.77	0.40
1:D:398:SER:HB3	1:D:465:THR:HG22	2.03	0.40
1:E:128:PHE:CE2	1:E:132:ALA:HB2	2.55	0.40
1:E:409:TRP:HZ3	1:E:470:ILE:HG23	1.86	0.40
1:G:478:ILE:HG23	1:G:482:LEU:HD12	2.03	0.40
1:C:103:PRO:HB2	1:C:143:ILE:HD13	2.04	0.40
1:G:273:ILE:HG12	1:G:368:GLY:HA2	2.04	0.40
1:H:128:PHE:CE2	1:H:132:ALA:HB2	2.56	0.40
1:K:336:TYR:HD1	1:L:197:ILE:HG13	1.86	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	426/551 (77%)	406 (95%)	18 (4%)	2 (0%)	24 54
1	B	432/551 (78%)	408 (94%)	21 (5%)	3 (1%)	18 47
1	C	435/551 (79%)	412 (95%)	21 (5%)	2 (0%)	24 54
1	D	434/551 (79%)	413 (95%)	20 (5%)	1 (0%)	43 71
1	E	432/551 (78%)	407 (94%)	23 (5%)	2 (0%)	24 54

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	435/551 (79%)	408 (94%)	25 (6%)	2 (0%)	24	54
1	G	434/551 (79%)	407 (94%)	25 (6%)	2 (0%)	24	54
1	H	433/551 (79%)	411 (95%)	20 (5%)	2 (0%)	24	54
1	I	428/551 (78%)	406 (95%)	20 (5%)	2 (0%)	24	54
1	J	434/551 (79%)	409 (94%)	24 (6%)	1 (0%)	43	71
1	K	433/551 (79%)	411 (95%)	19 (4%)	3 (1%)	18	47
1	L	431/551 (78%)	407 (94%)	23 (5%)	1 (0%)	43	71
All	All	5187/6612 (78%)	4905 (95%)	259 (5%)	23 (0%)	30	59

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	261	MET
1	H	261	MET
1	A	261	MET
1	B	163	ARG
1	B	185	THR
1	C	163	ARG
1	G	163	ARG
1	K	163	ARG
1	K	261	MET
1	A	163	ARG
1	B	445	TYR
1	C	261	MET
1	D	163	ARG
1	E	163	ARG
1	F	163	ARG
1	F	261	MET
1	H	163	ARG
1	I	163	ARG
1	L	163	ARG
1	G	445	TYR
1	I	445	TYR
1	J	261	MET
1	K	445	TYR

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	363/477 (76%)	363 (100%)	0	100	100
1	B	369/477 (77%)	369 (100%)	0	100	100
1	C	373/477 (78%)	373 (100%)	0	100	100
1	D	373/477 (78%)	373 (100%)	0	100	100
1	E	371/477 (78%)	371 (100%)	0	100	100
1	F	375/477 (79%)	375 (100%)	0	100	100
1	G	374/477 (78%)	374 (100%)	0	100	100
1	H	373/477 (78%)	373 (100%)	0	100	100
1	I	366/477 (77%)	366 (100%)	0	100	100
1	J	374/477 (78%)	372 (100%)	2 (0%)	81	81
1	K	373/477 (78%)	371 (100%)	2 (0%)	81	81
1	L	370/477 (78%)	370 (100%)	0	100	100
All	All	4454/5724 (78%)	4450 (100%)	4 (0%)	88	89

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	J	210	ILE
1	J	253	LEU
1	K	188	ILE
1	K	210	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	326	ASN
1	A	338	HIS
1	A	526	HIS
1	B	82	HIS
1	B	428	HIS

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Mol	Chain	Res	Type
1	C	146	HIS
1	C	153	GLN
1	C	507	GLN
1	C	526	HIS
1	D	446	HIS
1	E	187	ASN
1	E	446	HIS
1	F	122	GLN
1	F	179	HIS
1	F	338	HIS
1	F	428	HIS
1	F	446	HIS
1	G	446	HIS
1	G	526	HIS
1	H	338	HIS
1	J	179	HIS
1	J	187	ASN
1	J	526	HIS
1	K	102	GLN
1	K	179	HIS
1	L	428	HIS

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	434/551 (78%)	0.53	7 (1%) 70 56	40, 58, 80, 107	0
1	B	438/551 (79%)	0.50	2 (0%) 87 78	37, 53, 75, 98	0
1	C	441/551 (80%)	0.55	8 (1%) 67 53	39, 56, 81, 102	0
1	D	440/551 (79%)	0.50	8 (1%) 67 53	37, 53, 74, 99	0
1	E	438/551 (79%)	0.53	4 (0%) 81 68	37, 57, 78, 99	0
1	F	441/551 (80%)	0.58	13 (2%) 53 39	37, 53, 78, 95	0
1	G	440/551 (79%)	0.54	3 (0%) 84 73	43, 58, 83, 99	0
1	H	439/551 (79%)	0.64	15 (3%) 48 35	40, 60, 86, 111	0
1	I	434/551 (78%)	0.56	4 (0%) 81 68	39, 58, 82, 101	0
1	J	440/551 (79%)	0.66	6 (1%) 73 59	44, 61, 85, 99	0
1	K	439/551 (79%)	0.67	10 (2%) 61 46	38, 62, 87, 106	0
1	L	437/551 (79%)	0.68	13 (2%) 52 39	43, 67, 92, 103	0
All	All	5261/6612 (79%)	0.58	93 (1%) 67 53	37, 58, 84, 111	0

All (93) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	J	120	TYR	3.9
1	L	120	TYR	3.7
1	I	120	TYR	3.4
1	G	120	TYR	3.3
1	K	83	PRO	3.3
1	F	120	TYR	3.2
1	J	110	VAL	3.2
1	C	120	TYR	3.1
1	H	161	VAL	3.1
1	L	121	ASP	3.0
1	D	181	PRO	2.9

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Mol	Chain	Res	Type	RSRZ
1	F	417	PRO	2.9
1	I	83	PRO	2.9
1	A	111	ASP	2.8
1	D	120	TYR	2.8
1	L	83	PRO	2.8
1	B	111	ASP	2.8
1	E	111	ASP	2.8
1	H	120	TYR	2.8
1	H	122	GLN	2.8
1	L	108	ALA	2.7
1	K	120	TYR	2.7
1	H	121	ASP	2.7
1	B	87	GLU	2.7
1	L	105	GLY	2.7
1	F	181	PRO	2.6
1	J	382	LEU	2.6
1	I	258	GLY	2.6
1	C	188	ILE	2.5
1	C	242	CYS	2.5
1	A	338	HIS	2.5
1	D	336	TYR	2.5
1	K	193	GLY	2.5
1	A	120	TYR	2.5
1	C	110	VAL	2.5
1	F	110	VAL	2.5
1	K	310	LEU	2.5
1	L	104	VAL	2.5
1	J	512	ALA	2.5
1	H	375	MET	2.4
1	H	84	MET	2.4
1	F	248	ASP	2.4
1	I	189	VAL	2.4
1	E	530	ILE	2.4
1	A	188	ILE	2.4
1	K	188	ILE	2.4
1	K	418	LEU	2.4
1	H	530	ILE	2.4
1	E	459	THR	2.3
1	A	231	CYS	2.3
1	A	289	MET	2.3
1	K	379	TRP	2.3
1	H	108	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	99	PHE	2.3
1	L	421	CYS	2.3
1	L	161	VAL	2.3
1	H	338	HIS	2.3
1	A	121	ASP	2.3
1	F	87	GLU	2.2
1	H	192	PHE	2.2
1	L	99	PHE	2.2
1	F	111	ASP	2.2
1	F	288	SER	2.2
1	J	92	LEU	2.2
1	D	334	GLU	2.2
1	G	261	MET	2.2
1	H	296	GLY	2.2
1	H	111	ASP	2.2
1	L	89	TRP	2.2
1	K	459	THR	2.2
1	F	104	VAL	2.2
1	D	382	LEU	2.2
1	D	261	MET	2.2
1	F	503	ALA	2.2
1	K	197	ILE	2.1
1	E	83	PRO	2.1
1	L	249	THR	2.1
1	K	522	GLU	2.1
1	H	295	ASP	2.1
1	F	414	GLY	2.1
1	F	522	GLU	2.1
1	C	161	VAL	2.1
1	D	357	ASP	2.1
1	G	338	HIS	2.1
1	C	183	ARG	2.1
1	H	92	LEU	2.1
1	L	295	ASP	2.1
1	F	533	GLU	2.1
1	C	530	ILE	2.0
1	J	149	LEU	2.0
1	L	189	VAL	2.0
1	H	132	ALA	2.0
1	D	111	ASP	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.