



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 12:43 PM UTC

PDB ID : 4FQV / pdb_00004fqv
Title : Crystal structure of broadly neutralizing antibody CR9114 bound to H7 influenza hemagglutinin
Authors : Ekiert, D.C.; Dreyfus, C.; Wilson, I.A.
Deposited on : 2012-06-25
Resolution : 5.75 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

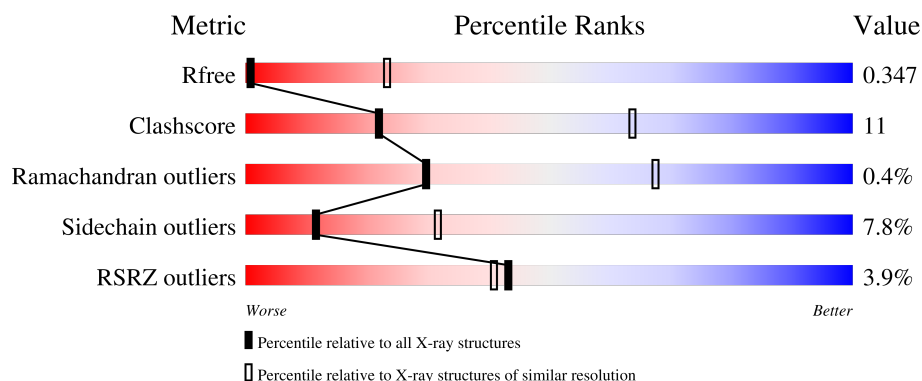
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 5.75 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	1119 (7.50-4.00)
Clashscore	190562	1182 (7.50-4.00)
Ramachandran outliers	187476	1014 (7.50-4.00)
Sidechain outliers	187428	1006 (7.54-3.98)
RSRZ outliers	180081	1112 (7.50-4.00)

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	327	4% 74% 17% 5% • •
1	C	327	3% 73% 18% 5% •
1	E	327	4% 77% 16% • •
2	B	176	7% 64% 26% 7% •
2	D	176	8% 63% 29% 5% •

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Mol	Chain	Length	Quality of chain
2	F	176	10% 65% 26% 5%
3	H	224	% 75% 18%
3	I	224	2% 84% 11%
3	J	224	6% 87% 8%
4	L	216	76% 19%
4	M	216	3% 85% 10%
4	N	216	80% 16%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 20938 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 Hemagglutinin HA1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	316	Total	C	N	O	S	0	0	0
			2423	1503	435	471	14			
1	C	316	Total	C	N	O	S	0	0	0
			2423	1503	435	471	14			
1	E	317	Total	C	N	O	S	0	0	0
			2427	1505	436	472	14			

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	7	ALA	-	expression tag	UNP Q6VMK1
A	8	ASP	-	expression tag	UNP Q6VMK1
A	9	PRO	-	expression tag	UNP Q6VMK1
A	10	GLY	-	expression tag	UNP Q6VMK1
A	254	LEU	PRO	conflict	UNP Q6VMK1
C	7	ALA	-	expression tag	UNP Q6VMK1
C	8	ASP	-	expression tag	UNP Q6VMK1
C	9	PRO	-	expression tag	UNP Q6VMK1
C	10	GLY	-	expression tag	UNP Q6VMK1
C	254	LEU	PRO	conflict	UNP Q6VMK1
E	7	ALA	-	expression tag	UNP Q6VMK1
E	8	ASP	-	expression tag	UNP Q6VMK1
E	9	PRO	-	expression tag	UNP Q6VMK1
E	10	GLY	-	expression tag	UNP Q6VMK1
E	254	LEU	PRO	conflict	UNP Q6VMK1

- Molecule 2 is a protein called Hemagglutinin HA2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	170	Total	C	N	O	S	0	0	0
			1380	851	243	279	7			

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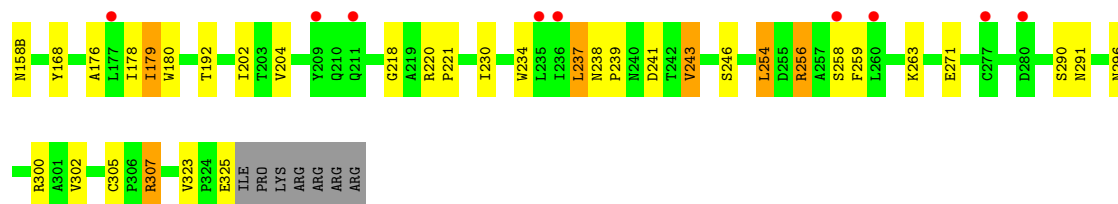
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	171	Total	C	N	O	S	0	0	0
			1388	857	244	280	7			
2	F	169	Total	C	N	O	S	0	0	0
			1369	845	239	278	7			

- Molecule 3 is a protein called Antibody CR9114 heavy chain.

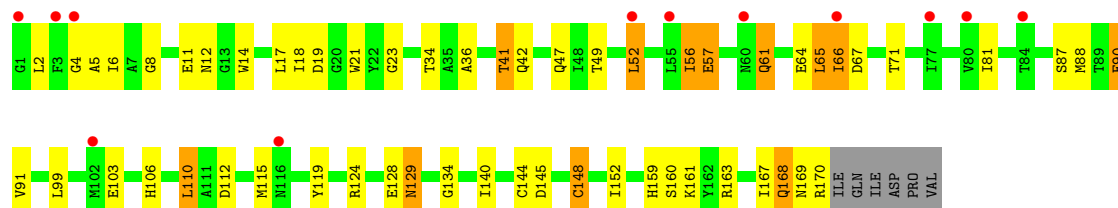
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	215	Total	C	N	O	S	0	2	0
			1608	1013	268	320	7			
3	I	215	Total	C	N	O	S	0	2	0
			1608	1013	268	320	7			
3	J	215	Total	C	N	O	S	0	2	0
			1608	1013	268	320	7			

- Molecule 4 is a protein called Antibody CR9114 light chain.

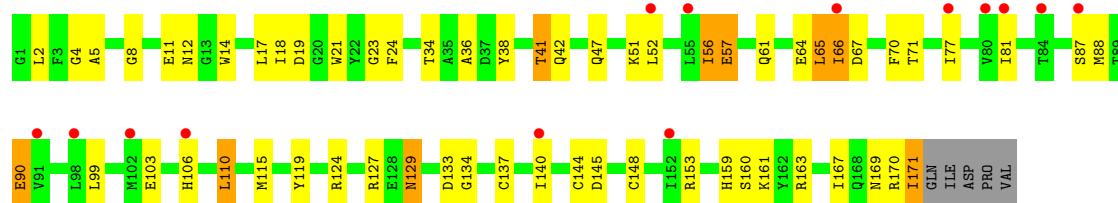
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	L	211	Total	C	N	O	S	0	0	0
			1568	978	266	320	4			
4	M	211	Total	C	N	O	S	0	0	0
			1568	978	266	320	4			
4	N	211	Total	C	N	O	S	0	0	0
			1568	978	266	320	4			



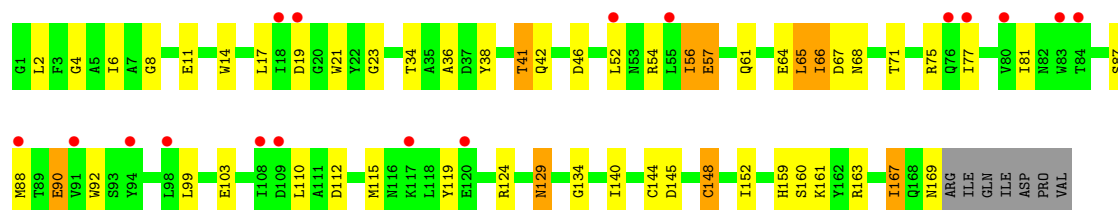
• Molecule 2: Hemagglutinin HA2



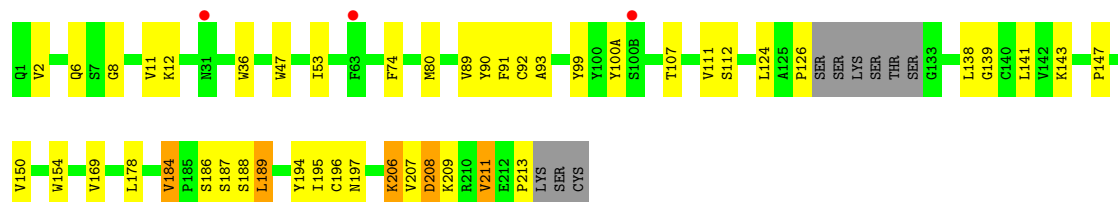
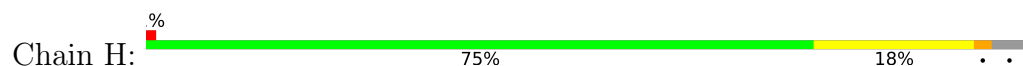
• Molecule 2: Hemagglutinin HA2



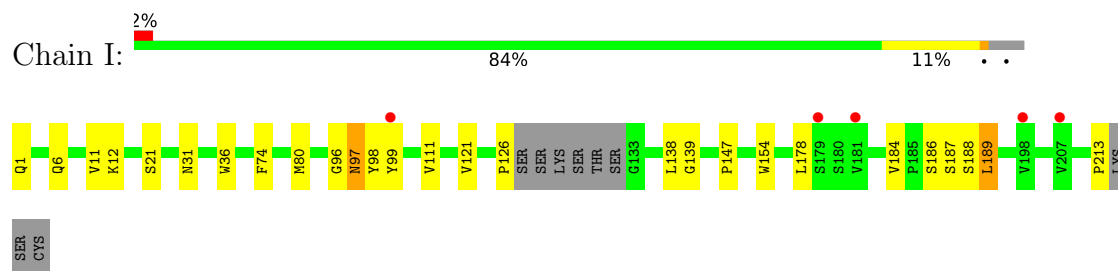
• Molecule 2: Hemagglutinin HA2



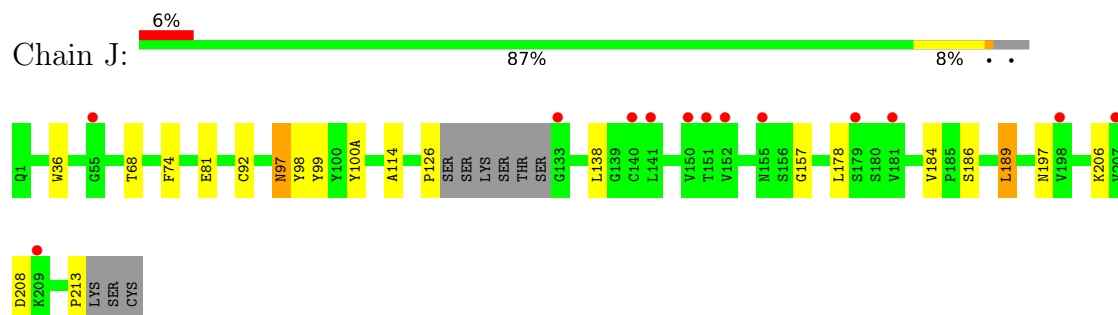
• Molecule 3: Antibody CR9114 heavy chain



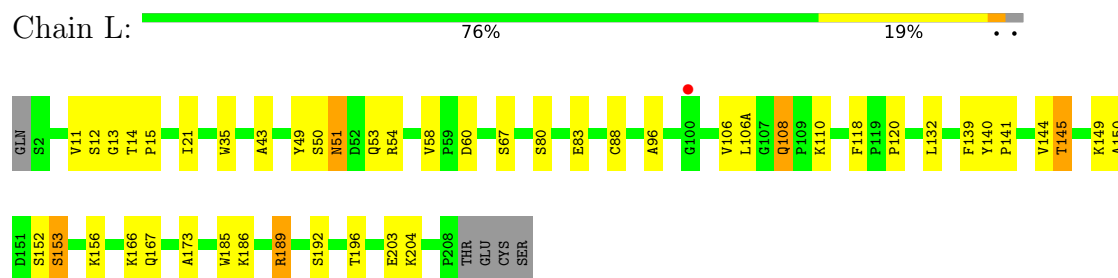
- Molecule 3: Antibody CR9114 heavy chain



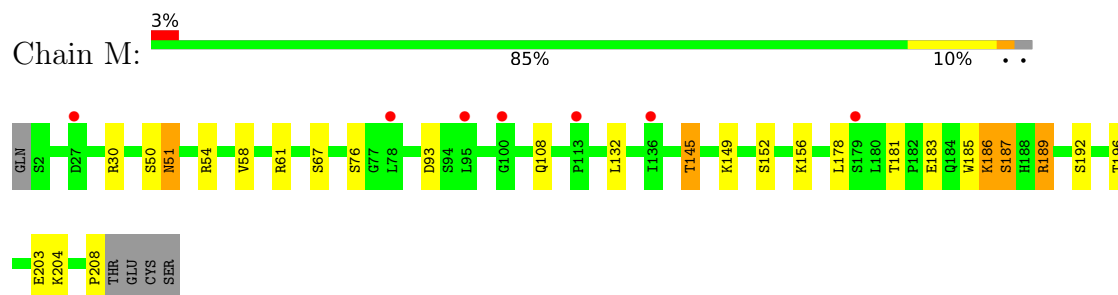
- Molecule 3: Antibody CR9114 heavy chain



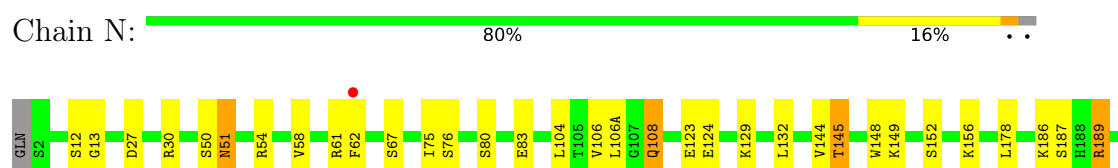
- Molecule 4: Antibody CR9114 light chain



- Molecule 4: Antibody CR9114 light chain



- Molecule 4: Antibody CR9114 light chain



S192					
T196					
H197					
E203					
K204					
P208					
THR					
GLU					
CYS					
SER					

4 Data and refinement statistics

Property	Value	Source
Space group	P 42 21 2	Depositor
Cell constants a, b, c, α , β , γ	197.57Å 197.57Å 223.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.35 – 5.75 49.35 – 5.75	Depositor EDS
% Data completeness (in resolution range)	99.4 (49.35-5.75) 98.7 (49.35-5.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.35 (at 5.73Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8_1069)	Depositor
R, R_{free}	0.314 , 0.339 0.323 , 0.347	Depositor DCC
R_{free} test set	642 reflections (4.89%)	wwPDB-VP
Wilson B-factor (Å ²)	319.3	Xtriage
Anisotropy	0.293	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 390.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.81	EDS
Total number of atoms	20938	wwPDB-VP
Average B, all atoms (Å ²)	244.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.52% 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	0.44	0/2468	0.82	3/3336 (0.1%)
1	C	0.45	0/2468	0.85	2/3336 (0.1%)
1	E	0.44	0/2472	0.83	3/3341 (0.1%)
2	B	0.41	0/1404	0.76	1/1892 (0.1%)
2	D	0.47	0/1412	0.80	0/1903
2	F	0.43	0/1393	0.79	2/1878 (0.1%)
3	H	0.49	0/1647	1.12	12/2245 (0.5%)
3	I	0.43	0/1647	0.82	4/2245 (0.2%)
3	J	0.49	0/1647	0.84	2/2245 (0.1%)
4	L	0.45	0/1606	0.90	5/2193 (0.2%)
4	M	0.48	0/1606	0.83	1/2193 (0.0%)
4	N	0.43	0/1606	0.87	3/2193 (0.1%)
All	All	0.45	0/21376	0.86	38/29000 (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	1
1	C	0	1
1	E	0	1
2	B	0	1
2	D	0	1
2	F	0	1
4	M	0	1
All	All	0	7

There are no bond length outliers.

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	L	153	SER	CA-C-N	-11.41	105.58	119.84
4	L	153	SER	C-N-CA	-11.41	105.58	119.84
3	H	92	CYS	N-CA-C	-10.22	93.14	109.59
3	I	97	ASN	N-CA-C	9.60	120.54	108.45
4	N	108	GLN	CA-C-N	9.09	128.84	119.76
4	N	108	GLN	C-N-CA	9.09	128.84	119.76
4	N	108	GLN	N-CA-C	8.15	119.92	109.65
2	F	167	ILE	N-CA-C	7.78	119.83	111.58
3	H	207	VAL	N-CA-C	7.22	118.54	107.78
3	H	207	VAL	CA-C-N	-7.06	113.25	123.00
3	H	207	VAL	C-N-CA	-7.06	113.25	123.00
3	H	187	SER	N-CA-C	-6.62	104.15	111.36
3	H	89	VAL	N-CA-C	-6.27	98.84	107.99
3	H	112	SER	N-CA-C	6.16	117.59	108.60
3	H	206	LYS	N-CA-C	-6.09	98.46	108.76
3	H	208	ASP	N-CA-C	-6.05	99.32	109.07
4	M	187	SER	N-CA-C	5.86	119.69	112.54
1	C	132	THR	N-CA-C	-5.68	106.00	114.64
3	I	96	GLY	N-CA-C	-5.63	108.47	114.67
1	E	305	CYS	CA-C-N	-5.62	113.99	119.78
1	E	305	CYS	C-N-CA	-5.62	113.99	119.78
4	L	108	GLN	CA-C-N	5.62	125.55	119.76
4	L	108	GLN	C-N-CA	5.62	125.55	119.76
3	I	187	SER	N-CA-C	-5.45	105.42	111.36
3	J	114	ALA	N-CA-C	5.39	118.10	110.50
3	H	93	ALA	N-CA-C	5.36	117.55	109.23
3	H	2	VAL	N-CA-C	5.25	116.54	108.71
1	C	243	VAL	CB-CA-C	-5.16	104.15	111.21
2	B	61	GLN	N-CA-C	5.15	117.12	108.73
1	E	84	ALA	N-CA-C	5.14	117.06	108.99
2	F	68	ASN	N-CA-C	5.11	116.50	107.61
3	J	97	ASN	N-CA-C	5.09	114.86	108.45
3	I	121	VAL	N-CA-C	5.08	115.28	108.17
4	L	150	ALA	N-CA-C	-5.06	100.27	108.52
3	H	6	GLN	N-CA-C	5.04	116.61	110.41
1	A	132	THR	N-CA-C	-5.02	107.01	114.64
1	A	305	CYS	CA-C-N	-5.01	114.62	119.78
1	A	305	CYS	C-N-CA	-5.01	114.62	119.78

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	29	LEU	Peptide
2	B	56	ILE	Peptide
1	C	29	LEU	Peptide
2	D	56	ILE	Peptide
1	E	29	LEU	Peptide
2	F	56	ILE	Peptide
4	M	108	GLN	Sidechain

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	2423	0	2373	65	0
1	C	2423	0	2373	57	0
1	E	2427	0	2376	53	0
2	B	1380	0	1284	53	0
2	D	1388	0	1295	58	0
2	F	1369	0	1271	49	0
3	H	1608	0	1556	63	0
3	I	1608	0	1556	35	0
3	J	1608	0	1556	36	2
4	L	1568	0	1517	37	2
4	M	1568	0	1517	15	1
4	N	1568	0	1517	30	0
All	All	20938	0	20191	439	3

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 11.

All (439) 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)
2:B:19:ASP:OD2	3:H:100(A):TYR:OH	1.65	1.11
1:C:123:MET:SD	1:C:254:LEU:HD13	2.00	1.01
3:H:11:VAL:CG2	3:H:147:PRO:HG3	1.94	0.97
1:C:123:MET:HE3	1:C:168:TYR:CG	2.01	0.96
1:E:123:MET:SD	1:E:254:LEU:CD1	2.55	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:38:TYR:HE1	3:I:97:ASN:HB2	1.30	0.94
1:E:123:MET:HE3	1:E:168:TYR:CG	2.01	0.94
2:D:56:ILE:HG12	3:J:74:PHE:CE1	2.03	0.94
1:A:123:MET:SD	1:A:254:LEU:HD13	2.07	0.93
2:D:56:ILE:HG12	3:J:74:PHE:CD1	2.03	0.93
1:E:123:MET:HE3	1:E:168:TYR:CD2	2.03	0.92
1:A:123:MET:SD	1:A:254:LEU:CD1	2.58	0.91
1:A:291:ASN:HD22	3:H:74:PHE:HB2	1.34	0.91
1:C:123:MET:HE3	1:C:168:TYR:CD2	2.06	0.91
4:L:15:PRO:HD3	4:L:106(A):LEU:O	1.70	0.90
1:C:70:ILE:HG21	1:C:179:ILE:HD13	1.53	0.89
1:C:123:MET:SD	1:C:254:LEU:CD1	2.60	0.89
1:E:123:MET:SD	1:E:254:LEU:HD13	2.12	0.87
1:A:123:MET:HE3	1:A:168:TYR:CD2	2.09	0.86
2:D:19:ASP:OD1	3:J:99:TYR:CD2	2.28	0.86
2:F:56:ILE:HG12	3:I:74:PHE:CD1	2.10	0.86
2:F:38:TYR:CE1	3:I:97:ASN:HB2	2.12	0.84
2:D:19:ASP:O	3:J:98:TYR:OH	1.95	0.84
1:E:123:MET:SD	1:E:254:LEU:HD11	2.17	0.84
2:D:56:ILE:CG1	3:J:74:PHE:CD1	2.62	0.82
1:A:123:MET:HE3	1:A:168:TYR:CG	2.14	0.82
3:H:11:VAL:HG21	3:H:147:PRO:HG3	1.62	0.81
3:H:206:LYS:NZ	3:H:208:ASP:OD1	2.16	0.79
1:C:291:ASN:HD22	3:J:74:PHE:HB2	1.47	0.77
4:M:187:SER:O	4:M:189:ARG:NH2	2.17	0.77
2:B:56:ILE:HG13	3:H:74:PHE:CD1	2.20	0.76
2:F:38:TYR:HE1	3:I:97:ASN:CB	1.97	0.76
2:B:19:ASP:OD1	3:H:99:TYR:CE2	2.38	0.75
4:L:13:GLY:CA	4:L:106(A):LEU:HD12	2.17	0.74
2:D:56:ILE:CG1	3:J:74:PHE:CE1	2.71	0.74
2:D:19:ASP:OD2	3:J:100(A):TYR:CZ	2.40	0.74
2:B:56:ILE:HG13	3:H:74:PHE:CE1	2.25	0.72
1:A:123:MET:SD	1:A:254:LEU:HD11	2.28	0.72
2:B:56:ILE:CG1	3:H:74:PHE:CE1	2.73	0.72
2:D:19:ASP:OD2	3:J:100(A):TYR:CE2	2.43	0.71
2:B:56:ILE:CG1	3:H:74:PHE:CD1	2.74	0.71
3:H:138:LEU:HD13	3:H:211:VAL:HG11	1.73	0.70
4:L:13:GLY:HA2	4:L:106(A):LEU:HD12	1.73	0.70
2:F:38:TYR:CE1	3:I:97:ASN:CB	2.74	0.70
4:M:181:THR:HB	4:M:183:GLU:OE1	1.92	0.70
2:F:56:ILE:HG12	3:I:74:PHE:HD1	1.55	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:123:MET:CE	1:E:168:TYR:CG	2.74	0.69
2:D:19:ASP:OD2	3:J:100(A):TYR:OH	2.09	0.69
1:E:178:ILE:HG23	1:E:254:LEU:CD2	2.21	0.69
2:D:124:ARG:HD3	2:F:134:GLY:HA2	1.74	0.69
1:E:179:ILE:O	1:E:254:LEU:HD23	1.93	0.69
1:A:158(A):ASP:O	1:A:158(B):ASN:HB2	1.92	0.69
1:A:179:ILE:O	1:A:254:LEU:HD23	1.93	0.69
1:E:158(A):ASP:O	1:E:158(B):ASN:HB2	1.93	0.69
1:E:178:ILE:HG23	1:E:254:LEU:HD21	1.74	0.68
1:A:291:ASN:ND2	3:H:74:PHE:HB2	2.07	0.68
2:F:19:ASP:OD1	3:I:99:TYR:CZ	2.47	0.68
2:B:124:ARG:HD3	2:D:134:GLY:HA2	1.74	0.68
4:N:13:GLY:HA2	4:N:106(A):LEU:CD1	2.23	0.67
1:C:158(A):ASP:O	1:C:158(B):ASN:HB2	1.94	0.67
1:C:123:MET:CE	1:C:168:TYR:CG	2.76	0.67
3:J:189:LEU:HD21	3:J:213:PRO:CG	2.25	0.67
2:F:56:ILE:HG12	3:I:74:PHE:CE1	2.28	0.67
1:A:15:LEU:H	1:A:15:LEU:HD22	1.59	0.67
1:A:291:ASN:CB	3:H:74:PHE:HB3	2.25	0.66
4:L:83:GLU:HG2	4:L:166:LYS:NZ	2.09	0.66
2:B:134:GLY:HA2	2:F:124:ARG:HD3	1.76	0.66
1:A:178:ILE:HG23	1:A:254:LEU:CD2	2.26	0.66
4:L:106:VAL:HG12	4:L:108:GLN:HG3	1.76	0.66
4:N:12:SER:OG	4:N:106(A):LEU:HD21	1.95	0.65
4:N:13:GLY:N	4:N:106(A):LEU:HG	2.10	0.65
1:C:178:ILE:HG23	1:C:254:LEU:CD2	2.26	0.65
3:H:36:TRP:CE2	3:H:80:MET:HB2	2.31	0.65
2:B:19:ASP:CG	3:H:100(A):TYR:HH	1.91	0.64
3:I:36:TRP:CE2	3:I:80:MET:HB2	2.33	0.64
2:F:56:ILE:CG1	3:I:74:PHE:CD1	2.81	0.64
4:N:54:ARG:NH1	4:N:58:VAL:O	2.31	0.64
2:D:56:ILE:HG21	3:J:74:PHE:HD1	1.63	0.63
2:D:19:ASP:OD1	3:J:99:TYR:CG	2.51	0.63
1:E:302:VAL:HG11	2:F:65:LEU:HD13	1.80	0.63
1:C:254:LEU:HD12	1:C:254:LEU:O	1.99	0.63
2:B:66:ILE:HD13	2:B:81:ILE:HG21	1.81	0.62
3:I:189:LEU:HD21	3:I:213:PRO:CG	2.29	0.62
4:L:83:GLU:HB2	4:L:106:VAL:HG23	1.79	0.62
2:B:128:GLU:O	2:B:170:ARG:NH1	2.33	0.61
3:H:189:LEU:HD21	3:H:213:PRO:CG	2.30	0.61
4:L:83:GLU:HG2	4:L:166:LYS:HZ1	1.63	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:L:145:THR:HG22	4:L:196:THR:OG1	2.00	0.61
1:C:179:ILE:O	1:C:254:LEU:HD23	2.01	0.61
1:E:123:MET:CE	1:E:168:TYR:CD2	2.82	0.61
1:A:70:ILE:CG2	1:A:179:ILE:HG13	2.31	0.61
1:C:300:ARG:HH21	2:D:67:ASP:HB3	1.65	0.61
4:L:15:PRO:CD	4:L:106(A):LEU:O	2.46	0.61
1:A:254:LEU:O	1:A:254:LEU:HD12	2.01	0.60
1:E:15:LEU:HD22	1:E:15:LEU:H	1.64	0.60
1:E:254:LEU:HD12	1:E:254:LEU:O	2.01	0.60
4:L:106:VAL:HB	4:L:108:GLN:OE1	2.02	0.60
1:C:123:MET:SD	1:C:254:LEU:HD11	2.41	0.60
1:A:70:ILE:HG21	1:A:179:ILE:CD1	2.32	0.60
1:A:15:LEU:H	1:A:15:LEU:CD2	2.14	0.60
2:B:56:ILE:HG12	3:H:74:PHE:CE1	2.36	0.60
2:D:17:LEU:HD11	2:D:36:ALA:HB2	1.82	0.59
3:H:169:VAL:O	3:H:169:VAL:HG13	2.00	0.59
1:C:291:ASN:ND2	3:J:74:PHE:HB2	2.18	0.59
1:A:123:MET:CE	1:A:168:TYR:CG	2.85	0.59
3:J:126:PRO:HG3	3:J:138:LEU:HB3	1.83	0.59
1:A:302:VAL:HG11	2:B:65:LEU:HD13	1.84	0.59
2:F:41:THR:HG21	3:I:98:TYR:CZ	2.37	0.58
4:N:13:GLY:HA2	4:N:106(A):LEU:HD12	1.85	0.58
4:M:61:ARG:HB2	4:M:76:SER:O	2.02	0.58
1:C:17:HIS:HA	2:D:21:TRP:O	2.03	0.58
4:N:50:SER:O	4:N:51:ASN:HB2	2.03	0.58
2:B:56:ILE:CG2	3:H:74:PHE:HD1	2.16	0.58
2:D:99:LEU:O	2:D:103:GLU:HG2	2.04	0.58
2:D:88:MET:HE2	2:F:87:SER:HB3	1.85	0.58
1:C:15:LEU:H	1:C:15:LEU:HD22	1.68	0.58
1:C:44:GLU:OE1	1:C:290:SER:HB2	2.03	0.58
3:H:126:PRO:HG3	3:H:138:LEU:HB3	1.86	0.57
1:C:120:LYS:HE3	1:C:258:SER:OG	2.04	0.57
1:C:302:VAL:HG11	2:D:65:LEU:HD13	1.85	0.57
1:E:70:ILE:HG21	1:E:179:ILE:CD1	2.35	0.57
1:E:15:LEU:H	1:E:15:LEU:CD2	2.17	0.57
2:F:17:LEU:HD11	2:F:36:ALA:HB2	1.86	0.57
1:C:123:MET:CE	1:C:168:TYR:CD2	2.85	0.57
2:D:56:ILE:CG2	3:J:74:PHE:HD1	2.18	0.57
2:D:66:ILE:HD13	2:D:81:ILE:HG21	1.85	0.57
2:D:56:ILE:CG1	3:J:74:PHE:HD1	2.16	0.57
1:E:70:ILE:CG2	1:E:179:ILE:HG13	2.34	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:M:50:SER:O	4:M:51:ASN:HB2	2.04	0.57
2:B:17:LEU:HD11	2:B:36:ALA:HB2	1.86	0.56
4:M:145:THR:HG22	4:M:196:THR:OG1	2.04	0.56
2:B:4:GLY:O	2:B:8:GLY:HA3	2.04	0.56
1:A:291:ASN:HD22	3:H:74:PHE:CB	2.15	0.56
2:D:4:GLY:O	2:D:8:GLY:HA3	2.06	0.56
2:B:87:SER:HB3	2:F:88:MET:HE2	1.87	0.55
2:F:163:ARG:O	2:F:167:ILE:HG13	2.06	0.55
2:D:56:ILE:HG12	3:J:74:PHE:HE1	1.65	0.55
2:B:56:ILE:CG1	3:H:74:PHE:HD1	2.19	0.55
2:D:163:ARG:O	2:D:167:ILE:HG13	2.07	0.55
1:E:100:GLY:HA3	1:E:230:ILE:O	2.07	0.55
1:A:123:MET:CE	1:A:168:TYR:CD2	2.85	0.55
1:A:172:ARG:HD3	1:A:259:PHE:CZ	2.42	0.55
2:B:19:ASP:OD2	3:H:100(A):TYR:CZ	2.58	0.55
2:B:61:GLN:OE1	2:D:90:GLU:HG2	2.06	0.55
1:A:180:TRP:HZ3	1:A:235:LEU:HD23	1.71	0.54
3:J:178:LEU:HD12	3:J:178:LEU:C	2.32	0.54
2:D:56:ILE:HG13	3:J:74:PHE:CD1	2.43	0.54
1:C:15:LEU:H	1:C:15:LEU:CD2	2.21	0.54
2:B:52:LEU:HD23	3:H:53:ILE:CD1	2.37	0.54
1:C:237:LEU:HG	1:C:243:VAL:HG22	1.90	0.54
4:L:50:SER:O	4:L:51:ASN:HB2	2.08	0.54
1:A:178:ILE:HG23	1:A:254:LEU:HD21	1.90	0.54
1:A:220:ARG:HB3	1:A:221:PRO:HD2	1.90	0.54
3:I:126:PRO:HG3	3:I:138:LEU:HB3	1.88	0.54
2:D:38:TYR:HE1	3:J:97:ASN:CB	2.20	0.54
4:L:167:GLN:OE1	4:L:173:ALA:HB2	2.07	0.53
3:I:11:VAL:HG21	3:I:147:PRO:HG3	1.91	0.53
3:I:178:LEU:C	3:I:178:LEU:HD12	2.34	0.53
1:C:172:ARG:HD3	1:C:259:PHE:CZ	2.44	0.53
1:E:325:GLU:N	1:E:325:GLU:OE1	2.42	0.53
4:M:183:GLU:O	4:M:186:LYS:HG2	2.09	0.53
1:A:291:ASN:HB3	3:H:74:PHE:HB3	1.89	0.52
1:E:120:LYS:HE3	1:E:258:SER:OG	2.09	0.52
1:A:291:ASN:ND2	3:H:74:PHE:CB	2.72	0.52
1:C:150:GLU:OE1	1:C:256:ARG:HD3	2.09	0.52
1:A:292:LEU:HD21	3:H:74:PHE:CZ	2.45	0.52
2:D:56:ILE:HG12	3:J:74:PHE:HD1	1.63	0.52
4:N:62:PHE:CD1	4:N:75:ILE:HG12	2.44	0.52
2:B:56:ILE:CG1	3:H:74:PHE:HE1	2.18	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:129:ASN:N	2:B:129:ASN:OD1	2.42	0.52
1:C:201:LEU:HD12	1:E:218:GLY:HA3	1.92	0.52
4:N:13:GLY:CA	4:N:106(A):LEU:HG	2.40	0.52
1:A:121:GLU:OE2	1:A:172:ARG:HD2	2.10	0.52
2:D:47:GLN:OE1	2:D:110:LEU:HD11	2.10	0.52
4:N:80:SER:HA	4:N:106:VAL:HG21	1.92	0.52
1:A:73:PRO:HB2	1:A:75:GLN:OE1	2.09	0.52
2:F:38:TYR:CE1	3:I:97:ASN:ND2	2.77	0.52
2:B:145:ASP:OD1	2:B:145:ASP:N	2.44	0.51
1:C:31:GLU:HG3	1:C:32:ARG:N	2.25	0.51
4:L:13:GLY:C	4:L:106(A):LEU:HB2	2.35	0.51
1:A:237:LEU:HG	1:A:243:VAL:HG22	1.92	0.51
2:B:128:GLU:OE1	2:D:170:ARG:NH2	2.44	0.51
1:C:100:GLY:HA3	1:C:230:ILE:O	2.10	0.51
1:E:237:LEU:HG	1:E:243:VAL:HG22	1.93	0.51
2:F:99:LEU:O	2:F:103:GLU:HG2	2.11	0.51
4:N:13:GLY:CA	4:N:106(A):LEU:CD1	2.87	0.51
2:B:168:GLN:C	2:B:170:ARG:H	2.18	0.51
1:C:220:ARG:HB3	1:C:221:PRO:HD2	1.92	0.51
2:D:129:ASN:OD1	2:D:129:ASN:N	2.43	0.51
3:H:11:VAL:HG21	3:H:147:PRO:CG	2.38	0.51
3:I:186:SER:O	3:I:189:LEU:HB2	2.10	0.51
1:A:150:GLU:OE1	1:A:256:ARG:HD3	2.10	0.51
2:B:163:ARG:O	2:B:167:ILE:HG13	2.11	0.51
1:A:291:ASN:CB	3:H:74:PHE:CB	2.89	0.50
2:B:56:ILE:HG12	3:H:74:PHE:CD1	2.46	0.50
2:D:145:ASP:OD1	2:D:145:ASP:N	2.43	0.50
1:E:220:ARG:HB3	1:E:221:PRO:HD2	1.93	0.50
4:N:145:THR:HG22	4:N:196:THR:OG1	2.12	0.50
1:A:120:LYS:HE3	1:A:258:SER:OG	2.12	0.50
2:B:41:THR:HG22	2:B:42:GLN:N	2.25	0.50
2:D:56:ILE:CG2	2:D:56:ILE:O	2.59	0.50
1:E:254:LEU:HD12	1:E:254:LEU:C	2.37	0.50
2:F:159:HIS:CG	2:F:160:SER:N	2.79	0.50
1:C:178:ILE:HG23	1:C:254:LEU:HD21	1.93	0.50
2:B:99:LEU:O	2:B:103:GLU:HG2	2.11	0.50
1:E:291:ASN:HD22	3:I:74:PHE:HB2	1.75	0.50
2:F:66:ILE:HD13	2:F:81:ILE:HG21	1.94	0.50
4:M:54:ARG:NH1	4:M:58:VAL:O	2.45	0.50
2:B:6:ILE:N	2:B:112:ASP:OD1	2.41	0.50
2:D:38:TYR:HE1	3:J:97:ASN:HB2	1.77	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:129:ASN:OD1	2:F:129:ASN:N	2.44	0.50
3:H:186:SER:O	3:H:189:LEU:HB2	2.12	0.49
2:B:19:ASP:OD1	3:H:99:TYR:CD2	2.64	0.49
1:A:44:GLU:OE1	1:A:290:SER:HB2	2.12	0.49
2:B:90:GLU:HG2	2:F:61:GLN:OE1	2.11	0.49
2:D:61:GLN:OE1	2:F:90:GLU:HG2	2.13	0.49
2:F:19:ASP:OD1	3:I:99:TYR:CE1	2.65	0.49
1:C:180:TRP:CH2	1:C:204:VAL:HG11	2.48	0.49
4:L:80:SER:HA	4:L:106:VAL:HG21	1.93	0.49
4:L:149:LYS:HB2	4:L:192:SER:HB2	1.94	0.49
2:D:41:THR:HG22	2:D:42:GLN:N	2.27	0.49
2:F:145:ASP:OD1	2:F:145:ASP:N	2.46	0.49
4:M:149:LYS:HB2	4:M:192:SER:HB2	1.93	0.49
1:A:17:HIS:HA	2:B:21:TRP:O	2.13	0.49
4:N:148:TRP:CE3	4:N:178:LEU:HD22	2.48	0.49
1:E:31:GLU:HG3	1:E:32:ARG:N	2.28	0.48
1:A:31:GLU:HG3	1:A:32:ARG:N	2.29	0.48
3:H:36:TRP:CD2	3:H:80:MET:HB2	2.48	0.48
4:L:49:TYR:O	4:L:53:GLN:HB2	2.13	0.48
1:C:80:LEU:O	1:C:120:LYS:NZ	2.46	0.48
2:F:4:GLY:O	2:F:8:GLY:HA3	2.14	0.48
2:F:41:THR:HG22	2:F:42:GLN:N	2.28	0.48
4:L:139:PHE:CE2	4:L:144:VAL:HG13	2.49	0.48
3:J:189:LEU:CD2	3:J:213:PRO:CG	2.92	0.48
1:E:178:ILE:CG2	1:E:254:LEU:HD21	2.42	0.48
2:F:148:CYS:O	2:F:152:ILE:HG13	2.13	0.48
1:E:103:VAL:O	1:E:104:ASN:HB2	2.13	0.48
1:E:180:TRP:CE2	1:E:204:VAL:HG21	2.49	0.48
1:E:70:ILE:HG21	1:E:179:ILE:HD11	1.96	0.48
1:E:300:ARG:HH21	2:F:67:ASP:HB3	1.79	0.48
3:H:188:SER:O	3:H:188:SER:OG	2.29	0.48
4:L:13:GLY:C	4:L:106(A):LEU:HD12	2.39	0.48
4:N:106:VAL:O	4:N:108:GLN:OE1	2.31	0.48
1:C:295:GLN:O	1:C:308:TYR:HA	2.14	0.48
1:E:150:GLU:OE1	1:E:256:ARG:HD3	2.13	0.48
3:H:90:TYR:N	3:H:90:TYR:CD2	2.82	0.48
2:B:23:GLY:HA3	2:B:36:ALA:HA	1.95	0.47
1:C:237:LEU:HD11	1:C:243:VAL:HG23	1.96	0.47
1:C:260:LEU:HD13	1:C:260:LEU:N	2.28	0.47
1:E:237:LEU:HD22	1:E:241:ASP:HB3	1.96	0.47
3:I:188:SER:O	3:I:188:SER:OG	2.31	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:254:LEU:HD12	1:A:254:LEU:C	2.38	0.47
1:C:103:VAL:O	1:C:104:ASN:HB2	2.14	0.47
1:A:292:LEU:CD2	3:H:74:PHE:CZ	2.97	0.47
3:I:126:PRO:HD2	3:I:213:PRO:HA	1.96	0.47
1:C:60:ASP:OD1	1:C:60:ASP:C	2.58	0.47
4:N:203:GLU:HG2	4:N:204:LYS:N	2.29	0.47
1:C:41:GLU:OE1	1:C:313:SER:OG	2.17	0.47
2:F:46:ASP:OD2	3:I:31:ASN:ND2	2.46	0.47
3:H:141:LEU:HG	3:H:143:LYS:HG3	1.97	0.47
3:I:36:TRP:CD2	3:I:80:MET:HB2	2.50	0.47
1:A:16:GLY:HA3	2:B:14:TRP:CH2	2.50	0.47
3:H:126:PRO:HD2	3:H:213:PRO:HA	1.97	0.47
3:H:138:LEU:HD13	3:H:211:VAL:CG1	2.43	0.47
3:J:189:LEU:HD21	3:J:213:PRO:HG3	1.95	0.47
1:C:296:ASN:CG	1:C:296:ASN:O	2.57	0.47
1:C:178:ILE:CG2	1:C:254:LEU:CD2	2.93	0.46
1:C:180:TRP:CE2	1:C:204:VAL:HG21	2.50	0.46
1:E:123:MET:CE	1:E:168:TYR:CB	2.94	0.46
1:A:100:GLY:HA3	1:A:230:ILE:O	2.16	0.46
4:L:106:VAL:CB	4:L:108:GLN:OE1	2.62	0.46
2:F:41:THR:HG21	3:I:98:TYR:CE2	2.51	0.46
1:C:121:GLU:OE2	1:C:172:ARG:HD2	2.16	0.46
1:A:70:ILE:HG21	1:A:179:ILE:HD11	1.97	0.46
2:D:133:ASP:OD1	2:D:137:CYS:O	2.34	0.46
1:E:179:ILE:HD13	1:E:234:TRP:HB3	1.97	0.46
4:L:54:ARG:NH1	4:L:58:VAL:O	2.42	0.46
1:A:201:LEU:CD2	1:C:218:GLY:HA3	2.46	0.46
2:D:115:MET:HE3	2:D:115:MET:HA	1.98	0.46
1:E:237:LEU:HD11	1:E:243:VAL:HG23	1.96	0.46
2:F:23:GLY:HA3	2:F:36:ALA:HA	1.97	0.46
1:A:132:THR:HG23	1:A:152:LYS:HD3	1.98	0.46
2:B:91:VAL:CG2	2:F:88:MET:HE1	2.46	0.46
2:D:5:ALA:HB1	2:D:115:MET:HG2	1.98	0.46
3:J:68:THR:HB	3:J:81:GLU:HB3	1.97	0.46
2:B:19:ASP:OD1	3:H:99:TYR:CZ	2.69	0.46
4:N:156:LYS:HB2	4:N:156:LYS:HE3	1.77	0.46
1:A:18:HIS:CE1	1:A:320:MET:HE3	2.51	0.46
2:D:77:ILE:HD11	2:F:77:ILE:HD12	1.98	0.46
4:L:13:GLY:O	4:L:106(A):LEU:N	2.46	0.46
4:M:203:GLU:HG2	4:M:204:LYS:N	2.31	0.46
3:J:186:SER:O	3:J:189:LEU:HB2	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:47:GLN:OE1	2:B:110:LEU:HD11	2.15	0.45
1:C:123:MET:HE2	1:C:123:MET:HB3	1.91	0.45
4:N:132:LEU:HB2	4:N:178:LEU:HB3	1.97	0.45
2:B:56:ILE:CG2	2:B:56:ILE:O	2.63	0.45
1:C:113:ARG:HB3	1:C:267:GLY:HA3	1.97	0.45
1:E:176:ALA:HB2	1:E:259:PHE:CE2	2.51	0.45
1:A:238:ASN:O	1:A:239:PRO:C	2.59	0.45
1:E:17:HIS:HA	2:F:21:TRP:O	2.17	0.45
4:L:54:ARG:NE	4:L:60:ASP:HA	2.31	0.45
2:D:127:ARG:HG3	2:D:159:HIS:CG	2.52	0.45
1:E:108:LEU:HD22	1:E:234:TRP:CD1	2.51	0.45
1:A:296:ASN:CG	1:A:296:ASN:O	2.59	0.45
4:N:13:GLY:C	4:N:106(A):LEU:HG	2.42	0.45
4:L:110:LYS:HG3	4:L:141:PRO:HD3	1.98	0.45
3:I:189:LEU:CD2	3:I:213:PRO:CG	2.94	0.45
1:C:188:THR:O	1:C:192:THR:HG23	2.17	0.45
2:F:56:ILE:CG1	3:I:74:PHE:HD1	2.22	0.45
4:N:12:SER:CB	4:N:106(A):LEU:HD21	2.46	0.45
2:B:159:HIS:CG	2:B:160:SER:N	2.84	0.45
4:N:187:SER:O	4:N:189:ARG:NH2	2.49	0.45
1:A:291:ASN:HB2	3:H:74:PHE:CG	2.52	0.45
2:D:19:ASP:OD1	3:J:99:TYR:CE2	2.67	0.44
2:D:133:ASP:OD2	2:D:137:CYS:HB2	2.17	0.44
4:N:12:SER:HB2	4:N:106(A):LEU:CD2	2.46	0.44
1:C:107:ALA:HB1	2:F:75:ARG:HB3	1.99	0.44
2:F:41:THR:CG2	3:I:98:TYR:CE2	3.01	0.44
3:H:47:TRP:CG	4:L:96:ALA:HB3	2.52	0.44
3:H:139:GLY:HA2	3:H:154:TRP:CH2	2.53	0.44
3:H:184:VAL:HG11	3:H:194:TYR:OH	2.18	0.44
3:I:6:GLN:HA	3:I:21:SER:O	2.17	0.44
2:D:106:HIS:O	2:D:110:LEU:HB2	2.18	0.44
1:E:307:ARG:HG2	2:F:92:TRP:CE2	2.53	0.44
3:H:196:CYS:O	3:H:196:CYS:SG	2.75	0.44
4:L:156:LYS:HB2	4:L:156:LYS:HE3	1.84	0.44
4:N:149:LYS:HB2	4:N:192:SER:HB2	1.98	0.44
2:F:56:ILE:O	2:F:56:ILE:CG2	2.64	0.44
1:C:82:PHE:CE1	1:C:117:GLY:HA2	2.53	0.44
2:D:56:ILE:HG21	3:J:74:PHE:CD1	2.48	0.44
1:E:180:TRP:CH2	1:E:204:VAL:HG11	2.52	0.44
3:H:189:LEU:CD2	3:H:213:PRO:CG	2.94	0.44
4:M:156:LYS:HB2	4:M:156:LYS:HE3	1.79	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:202:ILE:N	1:C:202:ILE:HD12	2.33	0.44
1:E:70:ILE:CD1	1:E:118:ILE:HD12	2.48	0.44
3:H:11:VAL:HG23	3:H:147:PRO:HG3	1.90	0.44
1:C:108:LEU:HD22	1:C:234:TRP:CD1	2.53	0.44
1:A:201:LEU:HD23	1:C:218:GLY:HA3	2.00	0.43
1:A:325:GLU:HA	2:B:12:ASN:HD22	1.82	0.43
2:B:88:MET:HE2	2:D:87:SER:HB3	2.00	0.43
1:E:178:ILE:CG2	1:E:254:LEU:CD2	2.93	0.43
2:F:6:ILE:N	2:F:112:ASP:OD1	2.45	0.43
1:C:123:MET:CE	1:C:168:TYR:CB	2.96	0.43
1:E:291:ASN:HD22	3:I:74:PHE:CB	2.31	0.43
2:F:56:ILE:HG21	3:I:74:PHE:HA	2.01	0.43
3:H:139:GLY:HA2	3:H:154:TRP:CZ2	2.53	0.43
2:D:24:PHE:CD1	2:D:153:ARG:HG2	2.53	0.43
1:A:237:LEU:HD11	1:A:243:VAL:HG23	2.01	0.43
1:E:307:ARG:HG2	2:F:92:TRP:CD2	2.53	0.43
2:D:23:GLY:HA3	2:D:36:ALA:HA	2.00	0.43
1:E:296:ASN:CG	1:E:296:ASN:O	2.61	0.43
1:A:170:ASN:HB2	1:A:237:LEU:HD13	2.00	0.43
4:N:13:GLY:C	4:N:106(A):LEU:HD12	2.44	0.43
1:A:202:ILE:HD12	1:A:202:ILE:N	2.33	0.43
1:A:300:ARG:HH21	2:B:67:ASP:HB3	1.83	0.43
4:L:120:PRO:HD3	4:L:132:LEU:HG	1.99	0.43
1:A:237:LEU:HD22	1:A:241:ASP:HB3	2.00	0.43
1:E:238:ASN:O	1:E:239:PRO:C	2.61	0.43
1:A:103:VAL:O	1:A:104:ASN:HB2	2.18	0.43
2:B:49:THR:OG1	3:H:53:ILE:HG21	2.18	0.43
4:N:61:ARG:HB2	4:N:76:SER:O	2.18	0.42
2:F:115:MET:HE3	2:F:115:MET:HA	2.01	0.42
3:H:91:PHE:CE1	4:L:43:ALA:HA	2.55	0.42
3:H:196:CYS:SG	3:H:209:LYS:HB3	2.59	0.42
4:M:185:TRP:CE2	4:M:208:PRO:HB3	2.53	0.42
1:A:60:ASP:OD1	1:A:60:ASP:C	2.62	0.42
2:B:159:HIS:C	2:B:161:LYS:H	2.27	0.42
2:B:56:ILE:HG12	3:H:74:PHE:HE1	1.80	0.42
3:H:8:GLY:O	3:H:107:THR:HG23	2.19	0.42
3:I:139:GLY:HA2	3:I:154:TRP:CH2	2.55	0.42
4:M:187:SER:O	4:M:189:ARG:CZ	2.67	0.42
1:A:291:ASN:CG	3:H:74:PHE:HB3	2.45	0.42
4:N:27:ASP:O	4:N:30:ARG:HD3	2.20	0.42
2:D:159:HIS:C	2:D:161:LYS:H	2.28	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:202:ILE:N	1:E:202:ILE:HD12	2.34	0.42
3:H:12:LYS:O	3:H:111:VAL:HA	2.20	0.42
4:L:11:VAL:HG21	4:L:21:ILE:HG12	2.01	0.42
4:L:14:THR:HA	4:L:106(A):LEU:O	2.19	0.42
1:A:123:MET:HE2	1:A:123:MET:HB3	1.98	0.42
1:A:180:TRP:HZ3	1:A:235:LEU:CD2	2.31	0.42
1:E:16:GLY:HA3	2:F:14:TRP:CH2	2.55	0.42
3:J:189:LEU:CD2	3:J:213:PRO:HG2	2.49	0.42
4:N:83:GLU:HG3	4:N:104:LEU:O	2.20	0.42
3:H:169:VAL:O	3:H:169:VAL:CG1	2.67	0.42
4:L:108:GLN:NE2	4:L:140:TYR:CE2	2.88	0.42
4:L:140:TYR:HA	4:L:141:PRO:C	2.44	0.42
3:J:197:ASN:ND2	3:J:208:ASP:OD1	2.48	0.42
3:I:189:LEU:HD21	3:I:213:PRO:HG3	2.02	0.42
3:J:126:PRO:HD2	3:J:213:PRO:HA	2.02	0.42
4:N:13:GLY:CA	4:N:106(A):LEU:HD12	2.47	0.42
1:A:29:LEU:HD12	2:F:54:ARG:CZ	2.50	0.41
3:H:178:LEU:HD12	3:H:178:LEU:C	2.44	0.41
4:L:106:VAL:CG1	4:L:108:GLN:HG3	2.48	0.41
3:I:1:GLN:OE1	3:I:1:GLN:N	2.48	0.41
1:A:28:THR:C	1:A:30:THR:H	2.27	0.41
2:D:159:HIS:CG	2:D:160:SER:N	2.87	0.41
2:D:38:TYR:HE1	3:J:97:ASN:HB3	1.84	0.41
1:A:320:MET:HG3	1:A:321:LYS:O	2.20	0.41
1:C:16:GLY:HA3	2:D:14:TRP:CH2	2.55	0.41
2:D:56:ILE:CG2	3:J:74:PHE:CD1	3.00	0.41
1:A:180:TRP:CZ3	1:A:235:LEU:CD2	3.03	0.41
1:C:41:GLU:HG3	1:C:42:THR:N	2.34	0.41
1:C:70:ILE:HD13	1:C:118:ILE:HD12	2.03	0.41
1:E:123:MET:HE2	1:E:123:MET:HB3	1.92	0.41
4:L:120:PRO:HD2	4:L:185:TRP:CZ2	2.55	0.41
4:M:132:LEU:HB2	4:M:178:LEU:HB3	2.02	0.41
1:E:70:ILE:HD13	1:E:118:ILE:HD12	2.03	0.41
1:A:179:ILE:HD13	1:A:234:TRP:HB3	2.02	0.41
1:A:180:TRP:CE2	1:A:204:VAL:HG21	2.55	0.41
1:E:60:ASP:OD1	1:E:60:ASP:C	2.63	0.41
1:E:132:THR:O	1:E:132:THR:CG2	2.68	0.41
4:L:35:TRP:CH2	4:L:88:CYS:HB3	2.55	0.41
4:M:181:THR:C	4:M:183:GLU:N	2.76	0.41
4:N:13:GLY:HA2	4:N:106(A):LEU:HD11	2.02	0.41
2:B:5:ALA:HB1	2:B:115:MET:HG2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:52:LEU:CD2	3:H:53:ILE:HD12	2.51	0.41
2:B:106:HIS:O	2:B:110:LEU:HB2	2.21	0.41
1:C:254:LEU:HD12	1:C:254:LEU:C	2.45	0.41
1:C:325:GLU:HA	2:D:12:ASN:HD22	1.85	0.41
4:N:123:GLU:OE1	4:N:123:GLU:N	2.41	0.41
4:N:124:GLU:HG2	4:N:129:LYS:O	2.21	0.41
2:D:51:LYS:NZ	2:D:103:GLU:O	2.54	0.41
2:D:170:ARG:O	2:D:171:ILE:HD13	2.21	0.41
3:H:124:LEU:HB3	4:L:118:PHE:CD1	2.56	0.41
4:L:12:SER:HB2	4:L:106(A):LEU:HD21	2.03	0.41
3:I:12:LYS:O	3:I:111:VAL:HA	2.20	0.41
4:N:144:VAL:HG12	4:N:197:HIS:HB2	2.03	0.41
4:L:203:GLU:O	4:L:204:LYS:HG2	2.21	0.40
2:B:66:ILE:HD13	2:B:81:ILE:CG2	2.49	0.40
3:H:195:ILE:HG23	3:H:209:LYS:O	2.22	0.40
4:M:30:ARG:NH1	4:M:93:ASP:OD2	2.53	0.40
3:J:36:TRP:CZ3	3:J:92:CYS:HB3	2.56	0.40
2:F:56:ILE:HG21	3:I:74:PHE:HD1	1.86	0.40
2:F:159:HIS:C	2:F:161:LYS:H	2.28	0.40
2:B:49:THR:HG23	3:H:53:ILE:HG13	2.04	0.40
2:B:148:CYS:O	2:B:152:ILE:HG13	2.20	0.40
1:C:269:GLN:HB3	2:D:70:PHE:CE1	2.57	0.40
3:H:197:ASN:HB3	3:H:206:LYS:NZ	2.36	0.40
3:J:197:ASN:HB3	3:J:206:LYS:NZ	2.36	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:L:189:ARG:NH2	3:J:157:GLY:O[6_554]	1.99	0.21
4:L:189:ARG:NE	3:J:157:GLY:O[6_554]	2.10	0.10
4:M:187:SER:OG	4:M:187:SER:OG[2_465]	2.14	0.06

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	314/327 (96%)	295 (94%)	19 (6%)	0	100	100
1	C	314/327 (96%)	297 (95%)	17 (5%)	0	100	100
1	E	315/327 (96%)	296 (94%)	19 (6%)	0	100	100
2	B	168/176 (96%)	151 (90%)	14 (8%)	3 (2%)	6	34
2	D	169/176 (96%)	152 (90%)	14 (8%)	3 (2%)	6	34
2	F	167/176 (95%)	151 (90%)	14 (8%)	2 (1%)	10	43
3	H	213/224 (95%)	210 (99%)	3 (1%)	0	100	100
3	I	213/224 (95%)	212 (100%)	1 (0%)	0	100	100
3	J	213/224 (95%)	212 (100%)	1 (0%)	0	100	100
4	L	209/216 (97%)	204 (98%)	4 (2%)	1 (0%)	24	63
4	M	209/216 (97%)	206 (99%)	2 (1%)	1 (0%)	24	63
4	N	209/216 (97%)	205 (98%)	3 (1%)	1 (0%)	24	63
All	All	2713/2829 (96%)	2591 (96%)	111 (4%)	11 (0%)	30	67

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	57	GLU
2	D	57	GLU
2	F	57	GLU
2	B	119	TYR
2	D	119	TYR
2	F	119	TYR
2	B	169	ASN
2	D	169	ASN
4	L	51	ASN
4	M	51	ASN
4	N	51	ASN

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	268/277 (97%)	232 (87%)	36 (13%)	4	14
1	C	268/277 (97%)	233 (87%)	35 (13%)	4	15
1	E	268/277 (97%)	236 (88%)	32 (12%)	5	17
2	B	145/151 (96%)	127 (88%)	18 (12%)	4	16
2	D	146/151 (97%)	128 (88%)	18 (12%)	4	17
2	F	144/151 (95%)	127 (88%)	17 (12%)	5	18
3	H	180/187 (96%)	176 (98%)	4 (2%)	45	64
3	I	180/187 (96%)	178 (99%)	2 (1%)	65	75
3	J	180/187 (96%)	178 (99%)	2 (1%)	65	75
4	L	175/180 (97%)	169 (97%)	6 (3%)	32	54
4	M	175/180 (97%)	170 (97%)	5 (3%)	37	57
4	N	175/180 (97%)	170 (97%)	5 (3%)	37	57
All	All	2304/2385 (97%)	2124 (92%)	180 (8%)	11	32

All (180) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	ASP
1	A	15	LEU
1	A	18	HIS
1	A	29	LEU
1	A	30	THR
1	A	31	GLU
1	A	34	VAL
1	A	42	THR
1	A	70	ILE
1	A	75	GLN
1	A	96	VAL
1	A	104	ASN
1	A	118	ILE
1	A	123	MET
1	A	131	ARG
1	A	132	THR
1	A	136	THR
1	A	140	ARG
1	A	146	SER
1	A	152	LYS

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Mol	Chain	Res	Type
1	A	155	LEU
1	A	173	LYS
1	A	179	ILE
1	A	192	THR
1	A	235	LEU
1	A	237	LEU
1	A	243	VAL
1	A	246	SER
1	A	254	LEU
1	A	256	ARG
1	A	263	LYS
1	A	271	GLU
1	A	290	SER
1	A	307	ARG
1	A	321	LYS
1	A	323	VAL
2	B	2	LEU
2	B	11	GLU
2	B	18	ILE
2	B	34	THR
2	B	41	THR
2	B	52	LEU
2	B	57	GLU
2	B	64	GLU
2	B	65	LEU
2	B	66	ILE
2	B	71	THR
2	B	90	GLU
2	B	110	LEU
2	B	129	ASN
2	B	140	ILE
2	B	144	CYS
2	B	148	CYS
2	B	168	GLN
1	C	15	LEU
1	C	18	HIS
1	C	29	LEU
1	C	30	THR
1	C	31	GLU
1	C	34	VAL
1	C	42	THR
1	C	70	ILE

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Mol	Chain	Res	Type
1	C	75	GLN
1	C	96	VAL
1	C	104	ASN
1	C	118	ILE
1	C	123	MET
1	C	131	ARG
1	C	132	THR
1	C	136	THR
1	C	140	ARG
1	C	146	SER
1	C	152	LYS
1	C	155	LEU
1	C	173	LYS
1	C	179	ILE
1	C	192	THR
1	C	201	LEU
1	C	237	LEU
1	C	243	VAL
1	C	246	SER
1	C	254	LEU
1	C	256	ARG
1	C	260	LEU
1	C	263	LYS
1	C	271	GLU
1	C	307	ARG
1	C	323	VAL
1	C	325	GLU
2	D	2	LEU
2	D	11	GLU
2	D	18	ILE
2	D	34	THR
2	D	41	THR
2	D	52	LEU
2	D	57	GLU
2	D	64	GLU
2	D	65	LEU
2	D	66	ILE
2	D	71	THR
2	D	90	GLU
2	D	110	LEU
2	D	129	ASN
2	D	140	ILE

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Mol	Chain	Res	Type
2	D	144	CYS
2	D	148	CYS
2	D	171	ILE
1	E	11	ASP
1	E	15	LEU
1	E	18	HIS
1	E	29	LEU
1	E	30	THR
1	E	31	GLU
1	E	34	VAL
1	E	42	THR
1	E	70	ILE
1	E	96	VAL
1	E	104	ASN
1	E	118	ILE
1	E	123	MET
1	E	131	ARG
1	E	132	THR
1	E	136	THR
1	E	140	ARG
1	E	146	SER
1	E	152	LYS
1	E	155	LEU
1	E	179	ILE
1	E	192	THR
1	E	237	LEU
1	E	243	VAL
1	E	246	SER
1	E	254	LEU
1	E	256	ARG
1	E	263	LYS
1	E	271	GLU
1	E	290	SER
1	E	307	ARG
1	E	323	VAL
2	F	2	LEU
2	F	11	GLU
2	F	34	THR
2	F	41	THR
2	F	52	LEU
2	F	57	GLU
2	F	64	GLU

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Mol	Chain	Res	Type
2	F	65	LEU
2	F	66	ILE
2	F	71	THR
2	F	90	GLU
2	F	110	LEU
2	F	129	ASN
2	F	140	ILE
2	F	144	CYS
2	F	148	CYS
2	F	169	ASN
3	H	150	VAL
3	H	184	VAL
3	H	189	LEU
3	H	211	VAL
4	L	67	SER
4	L	145	THR
4	L	152	SER
4	L	153	SER
4	L	186	LYS
4	L	189	ARG
3	I	184	VAL
3	I	189	LEU
4	M	67	SER
4	M	145	THR
4	M	152	SER
4	M	186	LYS
4	M	189	ARG
3	J	184	VAL
3	J	189	LEU
4	N	67	SER
4	N	145	THR
4	N	152	SER
4	N	186	LYS
4	N	189	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	158(B)	ASN
1	A	191	GLN
1	C	158(B)	ASN
1	E	158(B)	ASN

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Mol	Chain	Res	Type
1	E	191	GLN
2	F	60	ASN
3	H	164	HIS
3	H	192	GLN
3	H	204	ASN
3	I	192	GLN
3	I	204	ASN
4	M	108	GLN
4	M	197	HIS
3	J	192	GLN
4	N	37	GLN
4	N	108	GLN

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	316/327 (96%)	0.08	12 (3%) 44 41	178, 369, 432, 457	0
1	C	316/327 (96%)	0.07	9 (2%) 55 47	145, 357, 410, 426	0
1	E	317/327 (96%)	0.09	14 (4%) 39 38	166, 309, 394, 402	0
2	B	170/176 (96%)	0.22	12 (7%) 22 27	175, 222, 260, 276	0
2	D	171/176 (97%)	0.45	14 (8%) 17 24	168, 211, 267, 300	0
2	F	169/176 (96%)	0.38	17 (10%) 12 19	166, 217, 262, 274	0
3	H	215/224 (95%)	-0.07	3 (1%) 73 62	110, 233, 322, 341	2 (0%)
3	I	215/224 (95%)	-0.01	5 (2%) 61 52	65, 159, 190, 206	2 (0%)
3	J	215/224 (95%)	0.09	13 (6%) 27 32	52, 143, 198, 232	2 (0%)
4	L	211/216 (97%)	-0.30	1 (0%) 87 77	218, 272, 365, 402	0
4	M	211/216 (97%)	-0.05	7 (3%) 49 44	159, 188, 242, 274	0
4	N	211/216 (97%)	-0.33	1 (0%) 87 77	116, 159, 194, 204	0
All	All	2737/2829 (96%)	0.04	108 (3%) 43 41	52, 223, 397, 457	6 (0%)

All (108) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	212	SER	7.4
2	D	98	LEU	7.4
1	E	177	LEU	7.2
2	D	80	VAL	7.1
1	C	317	ALA	6.7
2	F	77	ILE	6.1
1	A	103	VAL	5.8
2	F	80	VAL	5.8
1	C	13	ILE	5.7
3	I	181	VAL	5.7
2	F	98	LEU	5.5

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Mol	Chain	Res	Type	RSRZ
4	M	136	ILE	5.5
2	D	84	THR	5.5
2	F	52	LEU	5.4
1	A	233	HIS	5.2
1	C	294	PHE	5.0
2	F	55	LEU	4.9
1	E	70	ILE	4.9
2	B	55	LEU	4.8
2	D	77	ILE	4.8
3	J	179	SER	4.7
2	F	19	ASP	4.7
1	E	260	LEU	4.5
3	J	198	VAL	4.5
1	E	277	CYS	4.5
2	F	108	ILE	4.3
1	A	213	PHE	4.1
1	E	258	SER	4.0
3	J	207	VAL	4.0
2	D	87	SER	4.0
2	D	102	MET	4.0
2	D	52	LEU	3.9
1	A	211	GLN	3.9
1	C	316	LEU	3.9
1	C	42	THR	3.9
3	J	150	VAL	3.9
2	F	76	GLN	3.7
2	F	84	THR	3.6
3	I	198	VAL	3.6
2	B	3	PHE	3.5
2	B	116	ASN	3.5
3	J	155	ASN	3.4
3	J	152	VAL	3.4
2	D	55	LEU	3.3
1	A	294	PHE	3.2
1	C	177	LEU	3.2
2	B	77	ILE	3.2
1	C	319	GLY	3.1
2	B	52	LEU	3.1
1	A	106	GLU	3.0
3	I	207	VAL	3.0
4	M	27	ASP	3.0
3	J	181	VAL	3.0

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Mol	Chain	Res	Type	RSRZ
2	B	60	ASN	3.0
3	J	209	LYS	2.9
3	I	99	TYR	2.9
1	E	29	LEU	2.9
2	B	4	GLY	2.8
2	D	140	ILE	2.8
2	D	106	HIS	2.7
1	A	158(B)	ASN	2.6
1	E	236	ILE	2.6
1	E	235	LEU	2.6
1	E	52	CYS	2.6
3	J	55	GLY	2.6
1	C	15	LEU	2.5
4	N	62	PHE	2.5
2	F	109	ASP	2.5
3	J	141	LEU	2.5
2	F	91	VAL	2.5
2	B	102	MET	2.5
2	B	80	VAL	2.4
4	M	113	PRO	2.4
1	E	209	TYR	2.4
1	A	174	ASP	2.4
4	M	179	SER	2.4
2	D	91	VAL	2.3
3	H	31	ASN	2.3
3	H	100(B)	SER	2.3
4	M	78	LEU	2.3
3	J	140	CYS	2.3
2	B	66	ILE	2.3
1	A	105	GLU	2.3
1	A	101	LYS	2.3
2	D	66	ILE	2.2
1	E	10	GLY	2.2
2	F	117	LYS	2.2
2	F	94	TYR	2.2
3	J	151	THR	2.2
3	H	63	PHE	2.2
2	B	84	THR	2.2
2	D	81	ILE	2.2
3	I	179	SER	2.2
1	E	71	THR	2.2
2	D	152	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	158(A)	ASP	2.1
4	M	95	LEU	2.1
2	F	120	GLU	2.1
2	F	18	ILE	2.1
2	F	88	MET	2.1
1	E	211	GLN	2.1
4	L	100	GLY	2.1
2	B	1	GLY	2.1
1	C	129	GLY	2.0
4	M	100	GLY	2.0
3	J	133	GLY	2.0
1	E	280	ASP	2.0
2	F	83	TRP	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.