



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:16 PM UTC

PDB ID : 8ZUA / pdb_00008zua
Title : The complex structure of MPXV M1R and neutralizing antibody A138
Authors : Ge, J.W.
Deposited on : 2024-06-08
Resolution : 3.49 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

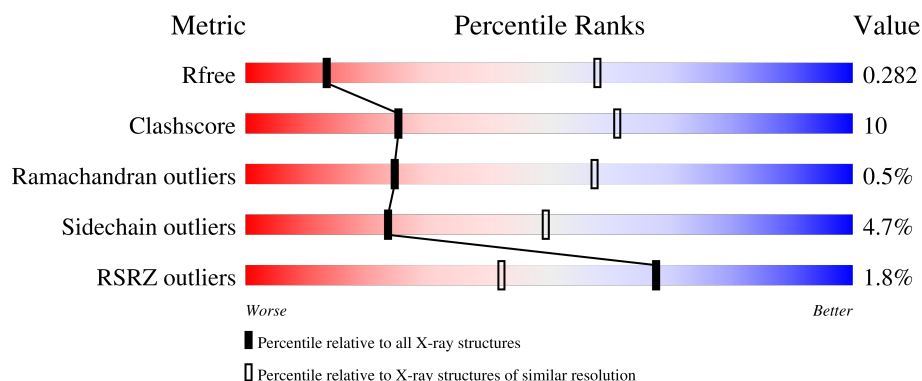
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.49 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	1085 (3.54-3.46)
Clashscore	190562	1140 (3.54-3.46)
Ramachandran outliers	187476	1113 (3.54-3.46)
Sidechain outliers	187428	1114 (3.54-3.46)
RSRZ outliers	180081	1084 (3.54-3.46)

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	218	72% 27%
1	B	218	77% 18%
1	E	218	83% 15%
1	J	218	78% 19%
2	C	187	68% 19% 11%

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Mol	Chain	Length	Quality of chain
2	F	187	% 61% 27% • 11%
2	H	187	% 63% 26% • 11%
2	K	187	2% 65% 22% • 11%
3	D	213	% 76% 17% 7%
3	G	213	2% 73% 19% • 7%
3	I	213	2% 73% 19% • 7%
3	L	213	4% 63% 26% • 7%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 17537 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 A138 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	218	Total	C	N	O	S	0	0	0
			1634	1030	275	323	6			
1	B	218	Total	C	N	O	S	0	0	0
			1618	1022	268	322	6			
1	E	218	Total	C	N	O	S	0	0	0
			1634	1030	275	323	6			
1	J	217	Total	C	N	O	S	0	0	0
			1621	1022	271	322	6			

- Molecule 2 is a protein called Entry-fusion complex associated protein OPG095.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	F	167	Total	C	N	O	S	0	0	0
			1247	766	210	261	10			
2	C	167	Total	C	N	O	S	0	0	0
			1253	769	213	261	10			
2	H	167	Total	C	N	O	S	0	0	0
			1249	767	213	259	10			
2	K	167	Total	C	N	O	S	0	0	0
			1243	764	210	259	10			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	146	MET	LEU	conflict	UNP M1LBP0
F	182	HIS	-	expression tag	UNP M1LBP0
F	183	HIS	-	expression tag	UNP M1LBP0
F	184	HIS	-	expression tag	UNP M1LBP0
F	185	HIS	-	expression tag	UNP M1LBP0
F	186	HIS	-	expression tag	UNP M1LBP0
F	187	HIS	-	expression tag	UNP M1LBP0
C	146	MET	LEU	conflict	UNP M1LBP0

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Chain	Residue	Modelled	Actual	Comment	Reference
C	182	HIS	-	expression tag	UNP M1LBP0
C	183	HIS	-	expression tag	UNP M1LBP0
C	184	HIS	-	expression tag	UNP M1LBP0
C	185	HIS	-	expression tag	UNP M1LBP0
C	186	HIS	-	expression tag	UNP M1LBP0
C	187	HIS	-	expression tag	UNP M1LBP0
H	146	MET	LEU	conflict	UNP M1LBP0
H	182	HIS	-	expression tag	UNP M1LBP0
H	183	HIS	-	expression tag	UNP M1LBP0
H	184	HIS	-	expression tag	UNP M1LBP0
H	185	HIS	-	expression tag	UNP M1LBP0
H	186	HIS	-	expression tag	UNP M1LBP0
H	187	HIS	-	expression tag	UNP M1LBP0
K	146	MET	LEU	conflict	UNP M1LBP0
K	182	HIS	-	expression tag	UNP M1LBP0
K	183	HIS	-	expression tag	UNP M1LBP0
K	184	HIS	-	expression tag	UNP M1LBP0
K	185	HIS	-	expression tag	UNP M1LBP0
K	186	HIS	-	expression tag	UNP M1LBP0
K	187	HIS	-	expression tag	UNP M1LBP0

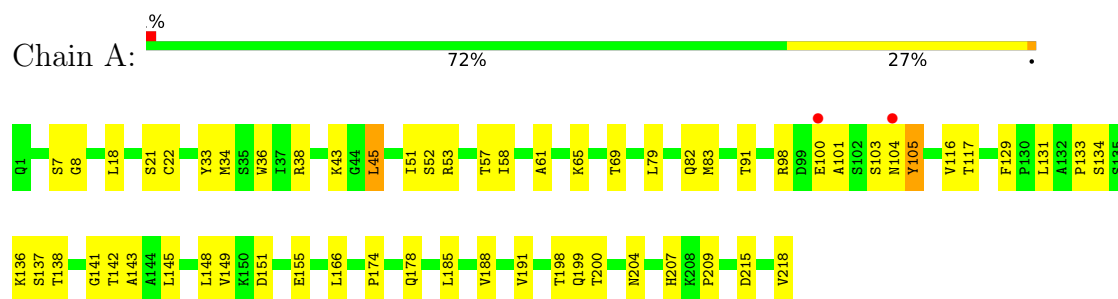
- Molecule 3 is a protein called A138 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	G	198	Total	C	N	O	S	0	0	0
			1508	944	253	307	4			
3	D	199	Total	C	N	O	S	0	0	0
			1514	947	254	309	4			
3	I	199	Total	C	N	O	S	0	0	0
			1508	944	251	309	4			
3	L	199	Total	C	N	O	S	0	0	0
			1508	944	251	309	4			

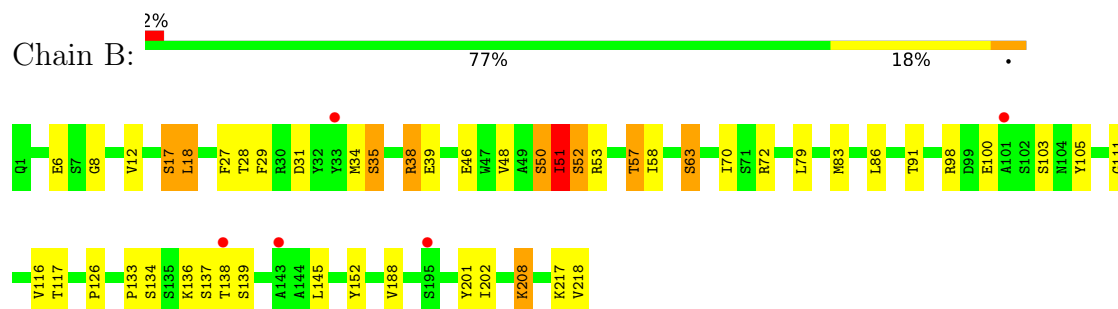
3 Residue-property plots [i](#)

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.

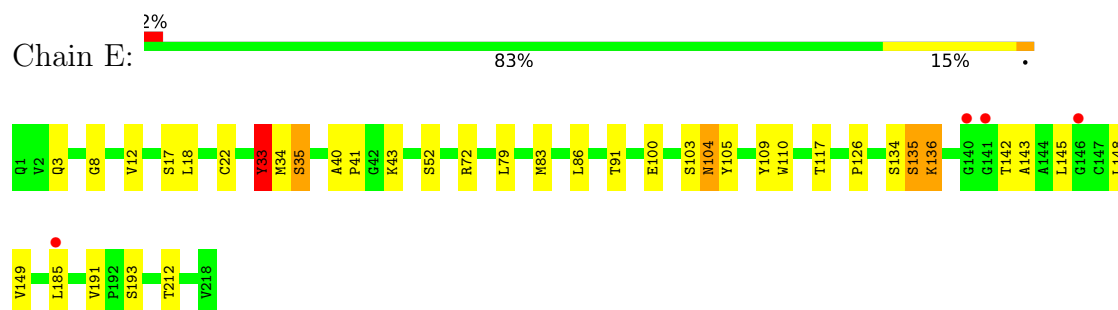
- Molecule 1: A138 heavy chain



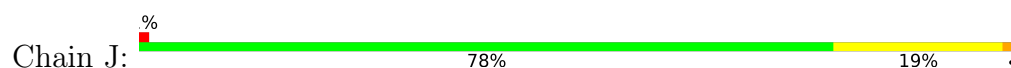
- Molecule 1: A138 heavy chain



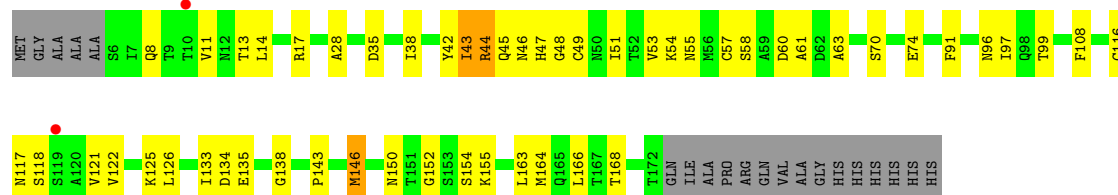
- Molecule 1: A138 heavy chain



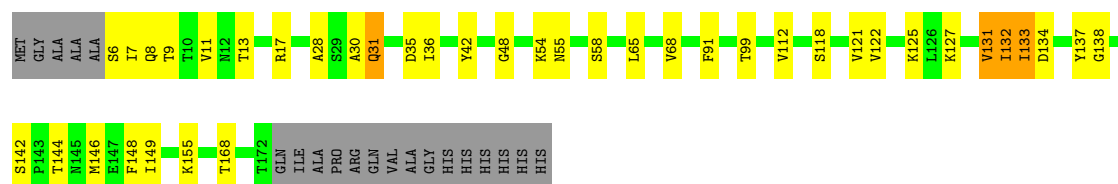
- Molecule 1: A138 heavy chain



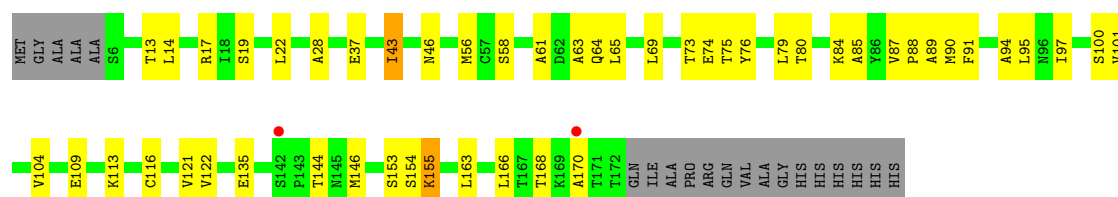
- Molecule 2: Entry-fusion complex associated protein OPG095



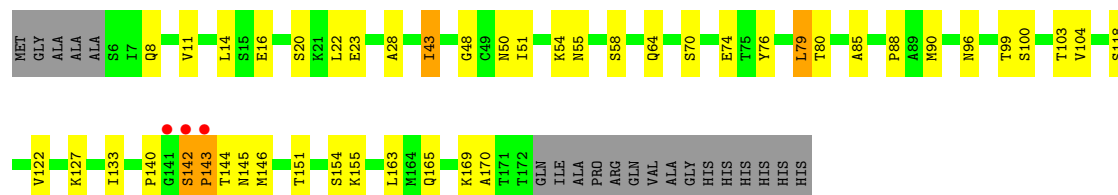
- Molecule 2: Entry-fusion complex associated protein OPG095



- Molecule 2: Entry-fusion complex associated protein OPG095

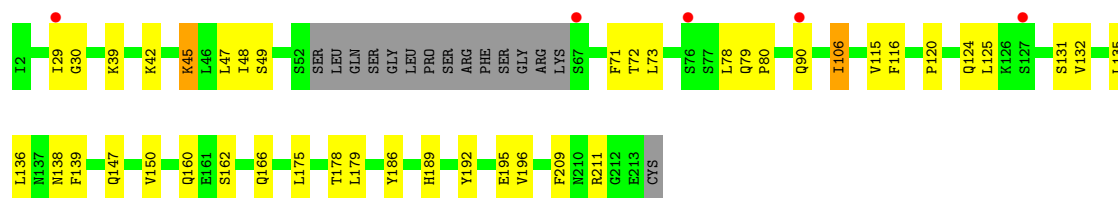


- Molecule 2: Entry-fusion complex associated protein OPG095



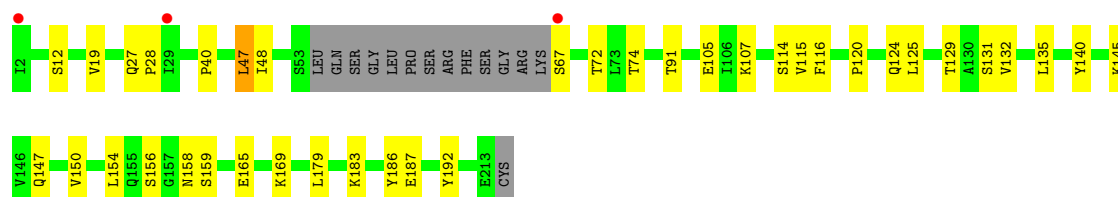
- Molecule 3: A138 light chain

Chain G:  73% 19% 7%



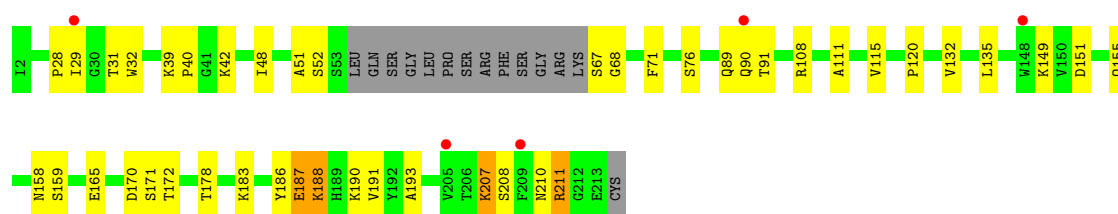
• Molecule 3: A138 light chain

Chain D:  76% 17% 7%



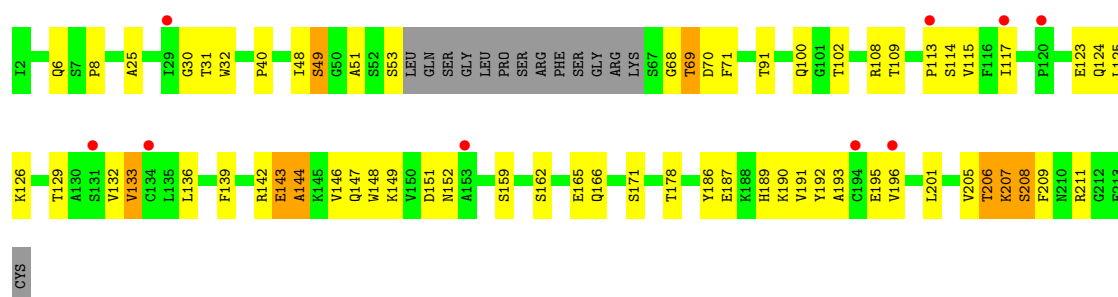
• Molecule 3: A138 light chain

Chain I:  73% 19% 7%



• Molecule 3: A138 light chain

Chain L:  63% 26% 7%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	82.39Å 136.68Å 123.77Å 90.00° 96.55° 90.00°	Depositor
Resolution (Å)	46.67 – 3.49 46.67 – 3.49	Depositor EDS
% Data completeness (in resolution range)	97.1 (46.67-3.49) 97.3 (46.67-3.49)	Depositor EDS
R_{merge}	0.25	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.18 (at 3.48Å)	Xtriage
Refinement program	PHENIX (1.18.2_3874: ???)	Depositor
R, R_{free}	0.237 , 0.286 0.238 , 0.282	Depositor DCC
R_{free} test set	1566 reflections (4.49%)	wwPDB-VP
Wilson B-factor (Å ²)	64.3	Xtriage
Anisotropy	0.626	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 54.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	17537	wwPDB-VP
Average B, all atoms (Å ²)	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.09% 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	0/1673	0.81	3/2277 (0.1%)
1	B	0.48	0/1657	0.81	2/2258 (0.1%)
1	E	0.46	1/1673 (0.1%)	0.76	2/2277 (0.1%)
1	J	0.45	0/1660	0.78	1/2260 (0.0%)
2	C	0.43	0/1266	0.84	0/1718
2	F	0.39	0/1260	0.72	0/1711
2	H	0.49	0/1262	0.90	0/1713
2	K	0.46	0/1256	0.82	1/1706 (0.1%)
3	D	0.40	0/1545	0.75	0/2101
3	G	0.42	0/1539	0.77	1/2093 (0.0%)
3	I	0.40	0/1539	0.82	1/2094 (0.0%)
3	L	0.47	1/1539 (0.1%)	0.87	4/2094 (0.2%)
All	All	0.44	2/17869 (0.0%)	0.81	15/24302 (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	B	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	L	49	SER	CA-CB	-5.62	1.46	1.53
1	E	35	SER	CA-CB	-5.20	1.47	1.53

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	139	PHE	CA-CB-CG	7.09	120.89	113.80
1	B	39	GLU	CB-CA-C	-5.74	101.43	110.79

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	105	TYR	CB-CA-C	5.55	116.63	109.80
3	L	132	VAL	CA-C-N	-5.38	116.12	123.12
3	L	132	VAL	C-N-CA	-5.38	116.12	123.12
1	J	44	GLY	CA-C-O	-5.37	118.53	122.23
1	A	131	LEU	N-CA-C	-5.34	100.39	108.67
3	L	113	PRO	CB-CA-C	-5.33	102.77	111.56
2	K	16	GLU	N-CA-C	-5.32	107.16	113.97
1	E	33	TYR	CB-CA-C	-5.26	103.54	112.00
1	B	63	SER	N-CA-C	-5.24	107.44	113.88
3	I	187	GLU	N-CA-C	-5.18	106.98	113.15
1	A	155	GLU	CA-C-O	-5.13	113.88	121.27
3	L	139	PHE	CA-CB-CG	5.11	118.91	113.80
1	E	104	ASN	N-CA-C	-5.04	104.68	111.54

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	38	ARG	Sidechain

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	1634	0	1596	39	0
1	B	1618	0	1568	37	0
1	E	1634	0	1596	24	0
1	J	1621	0	1576	30	1
2	C	1253	0	1233	23	0
2	F	1247	0	1222	32	1
2	H	1249	0	1229	32	0
2	K	1243	0	1218	23	0
3	D	1514	0	1476	27	0
3	G	1508	0	1471	27	0
3	I	1508	0	1465	28	0
3	L	1508	0	1465	41	0
All	All	17537	0	17115	336	1

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 10.

All (336) 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)
3:I:28:PRO:HB3	3:I:67:SER:HA	1.48	0.93
1:A:34:MET:HB3	1:A:79:LEU:HD22	1.70	0.72
3:I:151:ASP:HA	3:I:191:VAL:HB	1.72	0.71
1:A:151:ASP:OD1	1:A:178:GLN:NE2	2.24	0.71
2:K:58:SER:OG	2:K:64:GLN:NE2	2.25	0.70
2:C:8:GLN:HA	2:C:11:VAL:HG22	1.72	0.70
2:H:13:THR:O	2:H:17:ARG:HG3	1.92	0.70
2:H:17:ARG:NH2	2:H:74:GLU:OE1	2.25	0.69
3:L:187:GLU:HG2	3:L:211:ARG:NH1	2.08	0.69
2:C:133:ILE:HG21	2:C:168:THR:HG21	1.74	0.68
3:D:145:LYS:HE3	3:D:147:GLN:HG3	1.76	0.68
2:H:100:SER:O	2:H:104:VAL:HG13	1.94	0.68
3:L:195:GLU:HB2	3:L:206:THR:HG23	1.74	0.68
2:C:42:TYR:HB2	2:C:132:ILE:HG22	1.76	0.68
1:J:105:TYR:O	3:L:91:THR:HB	1.94	0.68
1:B:38:ARG:NH1	1:B:46:GLU:OE1	2.27	0.67
1:B:83:MET:HE1	1:B:116:VAL:HG11	1.76	0.66
1:E:149:VAL:HB	1:E:185:LEU:HB3	1.76	0.66
3:D:125:LEU:HD21	3:D:186:TYR:HD2	1.60	0.66
2:C:28:ALA:HB2	2:C:58:SER:HB3	1.77	0.66
3:D:120:PRO:HD3	3:D:132:VAL:HG22	1.75	0.66
3:I:31:THR:H	3:I:52:SER:HB3	1.60	0.65
2:F:8:GLN:HA	2:F:11:VAL:HG22	1.77	0.65
1:A:8:GLY:O	1:A:18:LEU:HD12	1.98	0.64
2:K:50:ASN:HB3	2:K:145:ASN:OD1	1.98	0.64
3:I:32:TRP:HB3	3:I:91:THR:HG22	1.80	0.63
1:E:22:CYS:HB3	1:E:79:LEU:HB3	1.80	0.63
1:E:104:ASN:HD22	3:I:51:ALA:HB2	1.63	0.63
1:J:57:THR:HG21	2:K:55:ASN:ND2	2.14	0.63
1:B:100:GLU:HB2	1:B:103:SER:HB3	1.80	0.63
2:F:43:ILE:O	2:F:44:ARG:C	2.41	0.63
1:A:83:MET:HE1	1:A:116:VAL:HG11	1.80	0.63
3:D:124:GLN:HG2	3:D:129:THR:O	1.99	0.62
2:K:165:GLN:O	2:K:169:LYS:HG2	2.00	0.62
2:H:122:VAL:HG22	2:H:155:LYS:HG2	1.82	0.62
2:K:14:LEU:HD21	2:K:163:LEU:HD21	1.80	0.62
1:J:40:ALA:HB1	1:J:41:PRO:HD2	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:K:28:ALA:HB2	2:K:58:SER:HB3	1.83	0.61
2:F:28:ALA:HB2	2:F:58:SER:HB3	1.83	0.61
3:D:116:PHE:HD2	3:D:135:LEU:HD23	1.66	0.61
3:L:187:GLU:HG2	3:L:211:ARG:HH12	1.63	0.61
1:J:98:ARG:HE	1:J:108:ASP:HB2	1.67	0.60
3:I:193:ALA:HA	3:I:208:SER:HA	1.82	0.60
3:G:189:HIS:O	3:G:211:ARG:HD3	2.02	0.59
2:H:116:CYS:HA	2:H:121:VAL:HG21	1.83	0.59
3:L:125:LEU:HD21	3:L:186:TYR:CD2	2.38	0.59
2:H:90:MET:SD	2:H:170:ALA:HB2	2.43	0.59
3:G:79:GLN:HG3	3:G:80:PRO:HD2	1.85	0.59
3:I:120:PRO:HD3	3:I:132:VAL:HG22	1.83	0.58
1:E:135:SER:O	1:E:136:LYS:C	2.46	0.58
3:L:193:ALA:HB1	3:L:206:THR:HG22	1.85	0.58
1:B:27:PHE:CE2	1:B:98:ARG:HD3	2.37	0.58
3:I:108:ARG:NH1	3:I:111:ALA:HB2	2.19	0.58
3:G:73:LEU:HD11	3:I:48:ILE:HD11	1.86	0.57
1:A:129:PHE:CD2	3:G:124:GLN:HG3	2.39	0.57
2:H:122:VAL:HG13	2:H:155:LYS:HE3	1.87	0.57
1:A:61:ALA:O	1:A:65:LYS:HB2	2.05	0.57
1:B:51:ILE:HD11	1:B:70:ILE:HG12	1.87	0.57
3:I:149:LYS:HB2	3:I:193:ALA:HB3	1.87	0.57
1:A:7:SER:OG	1:A:21:SER:OG	2.23	0.57
2:F:13:THR:O	2:F:17:ARG:HG2	2.05	0.56
1:A:207:HIS:HD2	1:A:209:PRO:HD2	1.70	0.56
1:A:207:HIS:CD2	1:A:209:PRO:HD2	2.39	0.56
1:A:104:ASN:O	1:A:105:TYR:C	2.48	0.56
1:B:51:ILE:HG23	1:B:58:ILE:HG12	1.87	0.56
2:H:14:LEU:HD21	2:H:163:LEU:HD11	1.88	0.56
2:K:8:GLN:HA	2:K:11:VAL:HG22	1.87	0.56
1:A:143:ALA:HB3	1:A:191:VAL:HG23	1.88	0.56
3:G:45:LYS:HE2	3:I:76:SER:HB2	1.88	0.56
2:H:19:SER:HB3	2:H:56:MET:HE3	1.87	0.55
1:B:51:ILE:CG2	1:B:58:ILE:HG12	2.37	0.55
3:L:124:GLN:HG2	3:L:129:THR:O	2.07	0.55
2:K:48:GLY:O	2:K:143:PRO:HB2	2.06	0.55
3:L:69:THR:OG1	3:L:70:ASP:N	2.39	0.55
2:H:28:ALA:HB2	2:H:58:SER:OG	2.06	0.55
2:F:35:ASP:HB2	2:F:152:GLY:O	2.07	0.55
1:B:52:SER:O	1:B:72:ARG:NH1	2.40	0.55
3:L:147:GLN:HB2	3:L:195:GLU:O	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:69:THR:OG1	1:A:82:GLN:HB3	2.08	0.54
3:L:148:TRP:CD1	3:L:159:SER:HG	2.25	0.54
1:A:91:THR:HG23	1:A:117:THR:HA	1.89	0.54
2:F:116:CYS:O	2:F:155:LYS:HB2	2.08	0.54
1:J:40:ALA:O	1:J:42:GLY:N	2.41	0.54
2:C:13:THR:O	2:C:17:ARG:HG2	2.09	0.54
1:J:98:ARG:O	1:J:107:PHE:HA	2.08	0.53
2:F:14:LEU:HD21	2:F:163:LEU:HD21	1.90	0.53
2:C:65:LEU:HD11	2:C:112:VAL:HG12	1.89	0.53
2:K:100:SER:O	2:K:104:VAL:HG23	2.08	0.53
1:A:101:ALA:HA	1:A:105:TYR:CE1	2.43	0.53
1:B:137:SER:HA	3:D:116:PHE:HD1	1.74	0.53
2:C:131:VAL:HG11	2:C:148:PHE:CE2	2.43	0.53
1:J:103:SER:OG	1:J:108:ASP:OD1	2.21	0.53
1:B:138:THR:OG1	3:D:115:VAL:O	2.24	0.53
3:D:125:LEU:HD21	3:D:186:TYR:CD2	2.43	0.53
2:H:14:LEU:HD13	2:H:75:THR:OG1	2.09	0.53
2:K:22:LEU:HD11	2:K:151:THR:HG22	1.91	0.53
1:A:133:PRO:HD3	1:A:218:VAL:HG12	1.90	0.53
3:L:6:GLN:H	3:L:100:GLN:HG2	1.74	0.53
2:C:30:ALA:HA	2:C:55:ASN:O	2.09	0.52
3:I:183:LYS:HG2	3:I:187:GLU:HG3	1.89	0.52
1:J:143:ALA:HB3	1:J:191:VAL:O	2.09	0.52
1:B:6:GLU:OE2	1:B:111:GLY:HA3	2.10	0.52
1:J:134:SER:C	1:J:136:LYS:H	2.17	0.52
2:F:51:ILE:HG12	2:F:146:MET:HB2	1.91	0.52
3:I:29:ILE:HB	3:I:90:GLN:HE21	1.75	0.51
1:J:40:ALA:HB3	1:J:43:LYS:HB2	1.93	0.51
1:B:34:MET:HB3	1:B:79:LEU:HD22	1.91	0.51
1:B:51:ILE:HG12	1:B:70:ILE:HD13	1.92	0.51
1:B:105:TYR:O	3:D:91:THR:HB	2.10	0.51
1:B:51:ILE:CD1	1:B:70:ILE:HG12	2.41	0.51
3:L:192:TYR:HB2	3:L:209:PHE:CE1	2.46	0.51
2:F:45:GLN:HG2	2:F:134:ASP:O	2.11	0.51
2:F:46:ASN:OD1	2:F:47:HIS:N	2.44	0.51
1:J:17:SER:HA	1:J:83:MET:O	2.10	0.51
1:B:133:PRO:HD3	1:B:218:VAL:HG12	1.93	0.50
2:F:47:HIS:O	2:F:138:GLY:N	2.45	0.50
1:E:52:SER:O	1:E:72:ARG:NH1	2.45	0.50
1:J:83:MET:HE1	1:J:116:VAL:HG11	1.92	0.50
3:L:143:GLU:O	3:L:144:ALA:HB2	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:143:ALA:HB2	1:E:193:SER:HA	1.94	0.50
3:G:116:PHE:HD2	3:G:135:LEU:HD23	1.76	0.49
3:G:147:GLN:HG2	3:G:195:GLU:HB3	1.93	0.49
2:H:109:GLU:HB3	2:H:113:LYS:NZ	2.27	0.49
2:K:90:MET:SD	2:K:170:ALA:HB2	2.52	0.49
2:F:164:MET:O	2:F:168:THR:HG23	2.12	0.49
3:G:106:ILE:O	3:G:166:GLN:NE2	2.33	0.49
2:F:91:PHE:CD2	2:F:99:THR:HG21	2.48	0.49
2:C:131:VAL:HG11	2:C:148:PHE:CZ	2.48	0.49
3:I:115:VAL:HA	3:I:135:LEU:O	2.13	0.49
2:K:88:PRO:HA	2:K:99:THR:HG23	1.93	0.49
1:E:17:SER:HA	1:E:83:MET:O	2.13	0.49
2:H:22:LEU:HD21	2:H:64:GLN:HA	1.94	0.49
1:B:137:SER:HA	3:D:116:PHE:CD1	2.48	0.49
2:C:42:TYR:O	2:C:132:ILE:HA	2.12	0.49
3:I:40:PRO:HB3	3:I:165:GLU:HG3	1.95	0.49
1:J:131:LEU:HB2	1:J:146:GLY:C	2.38	0.49
1:E:143:ALA:HB3	1:E:191:VAL:HG23	1.94	0.48
1:B:126:PRO:HB3	1:B:152:TYR:HB3	1.95	0.48
1:J:155:GLU:N	1:J:156:PRO:HD2	2.28	0.48
3:L:149:LYS:HD3	3:L:152:ASN:HA	1.95	0.48
1:A:51:ILE:HG13	1:A:58:ILE:HG12	1.94	0.48
2:H:43:ILE:HB	2:H:46:ASN:HB2	1.94	0.48
3:L:40:PRO:HG3	3:L:165:GLU:OE1	2.14	0.48
1:A:145:LEU:HB2	1:A:218:VAL:HG11	1.94	0.48
2:H:109:GLU:O	2:H:113:LYS:HG3	2.14	0.48
3:L:31:THR:OG1	3:L:51:ALA:C	2.57	0.48
3:D:183:LYS:HE3	3:D:187:GLU:OE2	2.13	0.48
3:I:31:THR:HA	3:I:71:PHE:HZ	1.79	0.48
1:E:34:MET:HB3	1:E:34:MET:HE3	1.70	0.48
1:A:136:LYS:HE2	3:G:209:PHE:HB3	1.96	0.48
1:B:83:MET:HB3	1:B:86:LEU:HD21	1.96	0.48
1:E:12:VAL:HB	1:E:18:LEU:HD21	1.96	0.48
1:J:138:THR:HG21	3:L:115:VAL:H	1.79	0.48
3:L:190:LYS:HA	3:L:211:ARG:HB2	1.94	0.48
2:F:91:PHE:HE1	2:F:166:LEU:HD11	1.79	0.47
2:H:76:TYR:HE1	2:H:87:VAL:HG21	1.79	0.47
1:J:97:ALA:HB1	1:J:107:PHE:HB3	1.96	0.47
1:J:136:LYS:HE2	1:J:136:LYS:HB2	1.55	0.47
2:K:100:SER:N	2:K:103:THR:OG1	2.43	0.47
2:H:144:THR:HG22	2:H:146:MET:HG3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:151:ASP:OD2	3:L:189:HIS:HB3	2.13	0.47
2:F:91:PHE:HZ	2:F:108:PHE:HB2	1.78	0.47
2:H:76:TYR:OH	2:H:101:VAL:HG13	2.14	0.47
3:I:170:ASP:O	3:I:172:THR:HG23	2.15	0.47
1:J:61:ALA:O	1:J:65:LYS:HB2	2.14	0.47
1:J:4:LEU:HB3	1:J:22:CYS:SG	2.55	0.47
3:L:108:ARG:HG2	3:L:109:THR:O	2.15	0.47
1:A:142:THR:HG22	1:A:191:VAL:O	2.15	0.47
1:B:12:VAL:HG11	1:B:86:LEU:HD12	1.97	0.47
1:B:51:ILE:HG12	1:B:70:ILE:CD1	2.45	0.47
1:B:139:SER:HB3	3:D:114:SER:OG	2.15	0.47
1:A:43:LYS:HE2	1:A:43:LYS:HB3	1.61	0.47
1:A:33:TYR:CD1	1:A:52:SER:HA	2.49	0.47
2:F:133:ILE:HG21	2:F:168:THR:HG21	1.96	0.47
3:I:186:TYR:C	3:I:188:LYS:H	2.22	0.47
3:G:160:GLN:O	3:G:178:THR:N	2.48	0.47
1:B:202:ILE:HD13	1:B:217:LYS:HA	1.96	0.47
3:D:150:VAL:HG22	3:D:192:TYR:CD2	2.50	0.46
1:E:83:MET:HB3	1:E:86:LEU:HD21	1.97	0.46
3:L:193:ALA:HA	3:L:208:SER:HB3	1.96	0.46
2:F:46:ASN:ND2	2:F:49:CYS:O	2.48	0.46
3:G:120:PRO:HD3	3:G:132:VAL:HG22	1.97	0.46
1:J:159:VAL:HG21	1:J:185:LEU:HD11	1.96	0.46
1:A:204:ASN:ND2	1:A:215:ASP:OD1	2.35	0.46
1:E:8:GLY:O	1:E:18:LEU:HD12	2.16	0.46
2:K:70:SER:O	2:K:74:GLU:HG3	2.15	0.46
2:F:116:CYS:HA	2:F:121:VAL:HG21	1.96	0.46
2:F:48:GLY:HA2	2:F:143:PRO:HB3	1.97	0.46
3:L:148:TRP:HD1	3:L:159:SER:HG	1.63	0.46
3:D:131:SER:HA	3:D:179:LEU:O	2.16	0.46
1:B:145:LEU:HD21	1:B:201:TYR:CD2	2.51	0.46
2:F:146:MET:HE2	2:F:146:MET:HB3	1.49	0.46
3:G:136:LEU:HB2	3:G:175:LEU:HB3	1.98	0.46
1:B:52:SER:HB3	1:B:57:THR:O	2.16	0.46
3:L:25:ALA:HB3	3:L:69:THR:HA	1.97	0.46
3:L:30:GLY:O	3:L:68:GLY:HA3	2.16	0.46
3:L:123:GLU:HA	3:L:126:LYS:HE2	1.96	0.46
3:L:133:VAL:HG22	3:L:178:THR:HG22	1.97	0.46
1:A:138:THR:HG21	3:G:115:VAL:HB	1.98	0.46
2:F:70:SER:O	2:F:74:GLU:HG3	2.16	0.46
3:I:210:ASN:O	3:I:211:ARG:C	2.58	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:K:8:GLN:O	2:K:11:VAL:HG22	2.15	0.46
1:B:105:TYR:OH	2:C:35:ASP:HA	2.16	0.45
1:J:31:ASP:O	2:K:127:LYS:NZ	2.46	0.45
1:A:36:TRP:HE1	1:A:79:LEU:HG	1.81	0.45
3:G:136:LEU:HD21	3:G:196:VAL:HG13	1.97	0.45
1:E:105:TYR:O	3:I:91:THR:OG1	2.34	0.45
1:E:148:LEU:HD12	1:E:185:LEU:O	2.16	0.45
1:J:103:SER:CB	1:J:108:ASP:OD1	2.64	0.45
2:F:117:ASN:OD1	2:F:155:LYS:HE3	2.17	0.45
2:C:91:PHE:CD2	2:C:99:THR:HG21	2.51	0.45
2:H:69:LEU:O	2:H:73:THR:HG23	2.16	0.45
2:K:51:ILE:HG12	2:K:146:MET:HB2	1.99	0.45
2:C:7:ILE:O	2:C:11:VAL:HG13	2.17	0.45
3:D:125:LEU:O	3:D:183:LYS:HD2	2.17	0.45
1:J:12:VAL:HG11	1:J:86:LEU:HD13	1.99	0.45
3:L:201:LEU:HD13	3:L:205:VAL:HG22	1.98	0.45
2:C:122:VAL:HG13	2:C:155:LYS:HD2	1.99	0.45
3:L:32:TRP:HB3	3:L:91:THR:OG1	2.16	0.45
3:L:207:LYS:HD2	3:L:207:LYS:HA	1.56	0.45
2:H:79:LEU:HD23	2:H:79:LEU:HA	1.77	0.45
2:H:87:VAL:C	2:H:89:ALA:N	2.74	0.45
2:F:61:ALA:HB1	2:F:155:LYS:HD3	1.99	0.44
2:C:54:LYS:HB3	2:C:149:ILE:HD13	1.98	0.44
1:E:3:GLN:HA	1:E:109:TYR:HE2	1.81	0.44
3:L:48:ILE:HG22	3:L:49:SER:H	1.80	0.44
3:G:30:GLY:O	3:G:71:PHE:HZ	2.00	0.44
3:L:108:ARG:HG2	3:L:109:THR:N	2.33	0.44
1:A:53:ARG:HH11	1:A:101:ALA:HB1	1.82	0.44
1:A:137:SER:CB	1:A:141:GLY:HA2	2.47	0.44
2:C:48:GLY:O	2:C:138:GLY:HA3	2.18	0.44
1:J:176:VAL:HG13	3:L:162:SER:OG	2.18	0.44
2:F:118:SER:O	2:F:122:VAL:HG23	2.16	0.44
3:G:72:THR:OG1	3:I:51:ALA:HB3	2.18	0.44
1:B:29:PHE:CE1	1:B:34:MET:HG3	2.53	0.44
3:D:145:LYS:HE3	3:D:147:GLN:CG	2.47	0.44
2:C:31:GLN:H	2:C:31:GLN:HG2	1.40	0.44
2:H:122:VAL:HG22	2:H:155:LYS:CG	2.48	0.44
1:J:22:CYS:HB3	1:J:79:LEU:HB3	2.00	0.44
2:F:38:ILE:HB	2:F:53:VAL:HG11	1.98	0.44
3:G:29:ILE:HG22	3:G:90:GLN:HE21	1.82	0.44
3:G:72:THR:OG1	1:E:104:ASN:ND2	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:40:ALA:HB1	1:J:41:PRO:CD	2.46	0.44
2:K:54:LYS:HA	2:K:54:LYS:HD3	1.85	0.44
2:K:76:TYR:O	2:K:79:LEU:HB2	2.18	0.44
2:K:118:SER:O	2:K:122:VAL:HG23	2.17	0.44
1:A:174:PRO:O	3:G:162:SER:OG	2.33	0.44
3:D:114:SER:HB3	3:D:116:PHE:CZ	2.53	0.44
3:I:207:LYS:HD3	3:I:207:LYS:HA	1.70	0.44
1:E:52:SER:HB2	2:H:37:GLU:OE1	2.18	0.43
3:I:155:GLN:HG3	3:I:158:ASN:OD1	2.17	0.43
1:B:208:LYS:HB2	1:B:208:LYS:HE2	1.45	0.43
2:C:118:SER:O	2:C:122:VAL:HG23	2.18	0.43
3:D:40:PRO:HB3	3:D:165:GLU:HG3	2.00	0.43
1:J:207:HIS:CD2	1:J:209:PRO:HD2	2.53	0.43
3:L:31:THR:HA	3:L:71:PHE:HZ	1.83	0.43
1:B:18:LEU:HD12	1:B:18:LEU:HA	1.77	0.43
3:G:150:VAL:HG22	3:G:192:TYR:CD2	2.54	0.43
2:C:6:SER:OG	2:C:9:THR:HG23	2.18	0.43
2:C:134:ASP:N	2:C:134:ASP:OD1	2.51	0.43
3:I:31:THR:HG22	3:I:31:THR:O	2.19	0.43
2:K:20:SER:O	2:K:23:GLU:HG2	2.18	0.43
3:D:47:LEU:HD23	3:D:47:LEU:HA	1.75	0.43
3:D:74:THR:H	3:L:49:SER:HG	1.67	0.43
2:H:155:LYS:H	2:H:155:LYS:HD2	1.83	0.43
3:I:39:LYS:HE3	3:I:42:LYS:HD3	2.01	0.43
1:J:134:SER:C	1:J:136:LYS:N	2.75	0.43
3:G:125:LEU:HA	3:G:125:LEU:HD23	1.63	0.43
3:D:27:GLN:O	3:D:28:PRO:C	2.59	0.43
2:F:54:LYS:HD3	2:F:54:LYS:HA	1.72	0.42
1:A:199:GLN:HG3	1:A:200:THR:N	2.34	0.42
3:D:107:LYS:HG2	3:D:140:TYR:OH	2.18	0.42
1:J:177:LEU:HB2	1:J:183:TYR:CE1	2.53	0.42
1:A:188:VAL:HG21	3:G:135:LEU:HD22	2.00	0.42
1:B:8:GLY:O	1:B:18:LEU:HG	2.18	0.42
1:E:91:THR:HG23	1:E:117:THR:HA	2.02	0.42
1:E:103:SER:O	1:E:104:ASN:C	2.61	0.42
1:A:7:SER:HG	1:A:21:SER:HG	1.67	0.42
1:A:148:LEU:HD12	1:A:185:LEU:O	2.19	0.42
3:I:89:GLN:HG2	3:I:90:GLN:O	2.20	0.42
3:D:169:LYS:HD2	3:D:169:LYS:HA	1.90	0.42
1:E:110:TRP:CD1	1:E:110:TRP:N	2.88	0.42
1:A:137:SER:HA	3:G:116:PHE:HD1	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:52:SER:OG	1:B:53:ARG:N	2.53	0.42
2:H:22:LEU:HD23	2:H:22:LEU:HA	1.88	0.42
1:A:100:GLU:OE1	1:A:103:SER:HB2	2.20	0.42
1:A:166:LEU:HD12	1:A:166:LEU:HA	1.91	0.42
2:F:91:PHE:HD2	2:F:99:THR:HG21	1.85	0.42
3:D:147:GLN:OE1	3:D:154:LEU:HD11	2.20	0.42
3:L:189:HIS:C	3:L:190:LYS:HD2	2.44	0.42
2:F:60:ASP:HB3	2:F:63:ALA:HB3	2.02	0.42
3:L:136:LEU:HD21	3:L:196:VAL:HG11	2.02	0.42
3:L:190:LYS:HG3	3:L:211:ARG:HB3	2.02	0.42
2:C:65:LEU:O	2:C:68:VAL:HG22	2.20	0.42
3:D:12:SER:HA	3:D:105:GLU:O	2.20	0.42
1:E:40:ALA:HB1	1:E:41:PRO:HD2	2.02	0.42
3:G:47:LEU:HD23	3:G:47:LEU:HA	1.81	0.41
1:B:28:THR:OG1	1:B:31:ASP:HB2	2.20	0.41
2:K:43:ILE:HA	2:K:133:ILE:O	2.20	0.41
3:L:148:TRP:HD1	3:L:159:SER:OG	2.03	0.41
1:A:149:VAL:HB	1:A:185:LEU:HB3	2.02	0.41
2:F:121:VAL:O	2:F:154:SER:OG	2.26	0.41
1:B:17:SER:HA	1:B:83:MET:O	2.20	0.41
1:B:91:THR:HG23	1:B:117:THR:HA	2.02	0.41
1:B:188:VAL:HG21	3:D:135:LEU:HD22	2.02	0.41
2:H:94:ALA:HB3	2:H:166:LEU:HD23	2.01	0.41
1:E:126:PRO:HD2	1:E:212:THR:OG1	2.20	0.41
2:H:91:PHE:CE2	2:H:97:ILE:HB	2.55	0.41
1:A:45:LEU:HD23	1:A:45:LEU:HA	1.85	0.41
2:H:61:ALA:HB1	2:H:155:LYS:HD3	2.02	0.41
2:H:85:ALA:O	2:H:88:PRO:HD2	2.20	0.41
1:A:34:MET:HB3	1:A:34:MET:HE3	1.91	0.41
1:A:57:THR:HG21	2:F:55:ASN:ND2	2.36	0.41
1:J:166:LEU:HD12	1:J:166:LEU:HA	1.87	0.41
3:G:78:LEU:HD23	3:G:78:LEU:HA	1.85	0.41
2:C:65:LEU:HD11	2:C:112:VAL:CG1	2.51	0.41
1:E:43:LYS:HA	1:E:43:LYS:HD3	1.82	0.41
1:A:22:CYS:HB3	1:A:79:LEU:HB3	2.03	0.41
2:K:85:ALA:O	2:K:88:PRO:HD2	2.20	0.41
3:G:131:SER:HA	3:G:179:LEU:O	2.21	0.41
1:B:134:SER:HB3	1:B:136:LYS:HE3	2.03	0.41
1:E:33:TYR:CD1	1:E:52:SER:HA	2.56	0.41
3:I:108:ARG:HG3	3:I:171:SER:HB2	2.02	0.41
3:L:151:ASP:OD1	3:L:191:VAL:HB	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:8:PRO:O	3:L:102:THR:OG1	2.35	0.41
3:D:19:VAL:O	3:D:74:THR:HA	2.21	0.40
3:L:166:GLN:HG3	3:L:171:SER:HA	2.02	0.40
3:G:186:TYR:HA	3:G:192:TYR:OH	2.21	0.40
2:C:17:ARG:HA	2:C:17:ARG:HD3	1.86	0.40
2:H:63:ALA:O	2:H:64:GLN:C	2.63	0.40
3:I:159:SER:HA	3:I:178:THR:O	2.21	0.40
2:F:61:ALA:HB1	2:F:155:LYS:CG	2.51	0.40
1:B:35:SER:HA	1:B:50:SER:HA	2.04	0.40
2:H:135:GLU:O	2:H:168:THR:HB	2.22	0.40
2:F:57:CYS:SG	2:F:150:ASN:ND2	2.93	0.40
2:H:91:PHE:CD2	2:H:97:ILE:HB	2.56	0.40

All (1) 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)
2:F:96:ASN:ND2	1:J:31:ASP:OD1[2_645]	1.88	0.32

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	216/218 (99%)	208 (96%)	8 (4%)	0	100	100
1	B	216/218 (99%)	203 (94%)	12 (6%)	1 (0%)	24	57
1	E	216/218 (99%)	206 (95%)	10 (5%)	0	100	100
1	J	215/218 (99%)	196 (91%)	17 (8%)	2 (1%)	14	47
2	C	165/187 (88%)	160 (97%)	5 (3%)	0	100	100
2	F	165/187 (88%)	159 (96%)	5 (3%)	1 (1%)	21	54
2	H	165/187 (88%)	159 (96%)	6 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	K	165/187 (88%)	154 (93%)	8 (5%)	3 (2%)	6	34
3	D	195/213 (92%)	188 (96%)	7 (4%)	0	100	100
3	G	194/213 (91%)	188 (97%)	5 (3%)	1 (0%)	24	57
3	I	195/213 (92%)	190 (97%)	3 (2%)	2 (1%)	12	44
3	L	195/213 (92%)	180 (92%)	13 (7%)	2 (1%)	12	44
All	All	2302/2472 (93%)	2191 (95%)	99 (4%)	12 (0%)	24	57

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	K	140	PRO
2	F	44	ARG
3	G	138	ASN
1	B	51	ILE
3	I	68	GLY
3	I	211	ARG
1	J	41	PRO
1	J	156	PRO
3	L	142	ARG
2	K	143	PRO
3	L	144	ALA
2	K	142	SER

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	181/181 (100%)	176 (97%)	5 (3%)	38	61
1	B	178/181 (98%)	168 (94%)	10 (6%)	19	46
1	E	181/181 (100%)	173 (96%)	8 (4%)	25	51
1	J	179/181 (99%)	171 (96%)	8 (4%)	24	50
2	C	142/155 (92%)	130 (92%)	12 (8%)	10	34

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	F	141/155 (91%)	134 (95%)	7 (5%)	22	48
2	H	141/155 (91%)	133 (94%)	8 (6%)	18	45
2	K	140/155 (90%)	132 (94%)	8 (6%)	18	45
3	D	174/186 (94%)	167 (96%)	7 (4%)	28	54
3	G	173/186 (93%)	167 (96%)	6 (4%)	32	57
3	I	173/186 (93%)	170 (98%)	3 (2%)	53	70
3	L	173/186 (93%)	163 (94%)	10 (6%)	18	44
All	All	1976/2088 (95%)	1884 (95%)	92 (5%)	23	49

All (92) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	ARG
1	A	45	LEU
1	A	98	ARG
1	A	134	SER
1	A	198	THR
2	F	42	TYR
2	F	43	ILE
2	F	97	ILE
2	F	125	LYS
2	F	126	LEU
2	F	135	GLU
2	F	146	MET
3	G	39	LYS
3	G	42	LYS
3	G	45	LYS
3	G	48	ILE
3	G	49	SER
3	G	106	ILE
1	B	17	SER
1	B	18	LEU
1	B	35	SER
1	B	48	VAL
1	B	50	SER
1	B	51	ILE
1	B	52	SER
1	B	57	THR
1	B	63	SER
1	B	208	LYS

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Mol	Chain	Res	Type
2	C	31	GLN
2	C	36	ILE
2	C	121	VAL
2	C	125	LYS
2	C	127	LYS
2	C	131	VAL
2	C	132	ILE
2	C	133	ILE
2	C	137	TYR
2	C	142	SER
2	C	144	THR
2	C	146	MET
3	D	47	LEU
3	D	48	ILE
3	D	67	SER
3	D	72	THR
3	D	156	SER
3	D	158	ASN
3	D	159	SER
1	E	33	TYR
1	E	35	SER
1	E	100	GLU
1	E	134	SER
1	E	135	SER
1	E	136	LYS
1	E	142	THR
1	E	145	LEU
2	H	43	ILE
2	H	65	LEU
2	H	80	THR
2	H	84	LYS
2	H	95	LEU
2	H	153	SER
2	H	154	SER
2	H	155	LYS
3	I	188	LYS
3	I	190	LYS
3	I	207	LYS
1	J	38	ARG
1	J	135	SER
1	J	136	LYS
1	J	138	THR

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Mol	Chain	Res	Type
1	J	139	SER
1	J	142	THR
1	J	145	LEU
1	J	157	VAL
2	K	43	ILE
2	K	79	LEU
2	K	80	THR
2	K	96	ASN
2	K	142	SER
2	K	144	THR
2	K	154	SER
2	K	155	LYS
3	L	53	SER
3	L	69	THR
3	L	114	SER
3	L	117	ILE
3	L	133	VAL
3	L	143	GLU
3	L	146	VAL
3	L	206	THR
3	L	207	LYS
3	L	208	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	82	GLN
2	F	55	ASN
2	F	129	GLN
3	G	89	GLN
1	B	77	ASN
2	C	55	ASN
2	C	83	GLN
2	C	110	ASN
2	C	124	ASN
3	D	199	GLN
1	E	104	ASN
1	E	171	HIS
2	H	45	GLN
2	H	46	ASN
3	I	37	GLN
3	I	38	GLN

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Mol	Chain	Res	Type
3	I	137	ASN
3	I	155	GLN
2	K	40	ASN
3	L	100	GLN
3	L	124	GLN
3	L	189	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	218/218 (100%)	0.08	2 (0%) 81 58	28, 55, 80, 112	0
1	B	218/218 (100%)	0.20	5 (2%) 61 37	38, 63, 89, 115	0
1	E	218/218 (100%)	0.22	4 (1%) 67 42	38, 66, 121, 139	0
1	J	217/218 (99%)	0.33	2 (0%) 81 58	30, 66, 126, 149	0
2	C	167/187 (89%)	0.24	0 100 100	51, 75, 109, 139	0
2	F	167/187 (89%)	0.09	2 (1%) 76 52	34, 57, 96, 130	0
2	H	167/187 (89%)	0.50	2 (1%) 76 52	51, 92, 131, 154	0
2	K	167/187 (89%)	0.27	3 (1%) 67 42	41, 73, 115, 135	0
3	D	199/213 (93%)	0.19	3 (1%) 72 47	42, 66, 91, 99	0
3	G	198/213 (92%)	0.15	5 (2%) 58 35	30, 56, 80, 89	0
3	I	199/213 (93%)	0.47	5 (2%) 58 35	41, 79, 118, 129	0
3	L	199/213 (93%)	0.61	9 (4%) 38 20	44, 88, 139, 157	0
All	All	2334/2472 (94%)	0.28	42 (1%) 67 42	28, 67, 121, 157	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	L	117	ILE	3.1
3	D	29	ILE	2.8
3	G	67	SER	2.8
1	B	33	TYR	2.7
3	I	90	GLN	2.7
2	H	142	SER	2.7
3	I	29	ILE	2.7
1	B	195	SER	2.6
3	I	209	PHE	2.5
3	L	134	CYS	2.5
3	L	29	ILE	2.4

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Mol	Chain	Res	Type	RSRZ
3	L	153	ALA	2.4
2	H	170	ALA	2.4
3	G	127	SER	2.4
3	G	76	SER	2.4
3	D	67	SER	2.3
3	L	120	PRO	2.3
1	A	104	ASN	2.3
1	J	130	PRO	2.3
1	B	101	ALA	2.3
3	I	148	TRP	2.3
3	G	90	GLN	2.3
1	E	141	GLY	2.2
3	L	113	PRO	2.2
1	B	138	THR	2.2
2	F	119	SER	2.2
3	L	194	CYS	2.2
3	L	131	SER	2.2
2	K	142	SER	2.2
1	A	100	GLU	2.2
3	I	205	VAL	2.1
3	D	2	ILE	2.1
1	J	147	CYS	2.1
2	K	143	PRO	2.1
1	E	185	LEU	2.1
3	L	196	VAL	2.0
3	G	29	ILE	2.0
1	B	143	ALA	2.0
2	F	10	THR	2.0
1	E	140	GLY	2.0
1	E	146	GLY	2.0
2	K	141	GLY	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.