



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:40 PM UTC

PDB ID : 4TUG / pdb_00004tug
Title : Crystal structure of MjMre11-DNA2 complex
Authors : Sung, S.; Cho, Y.
Deposited on : 2014-06-24
Resolution : 3.55 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:

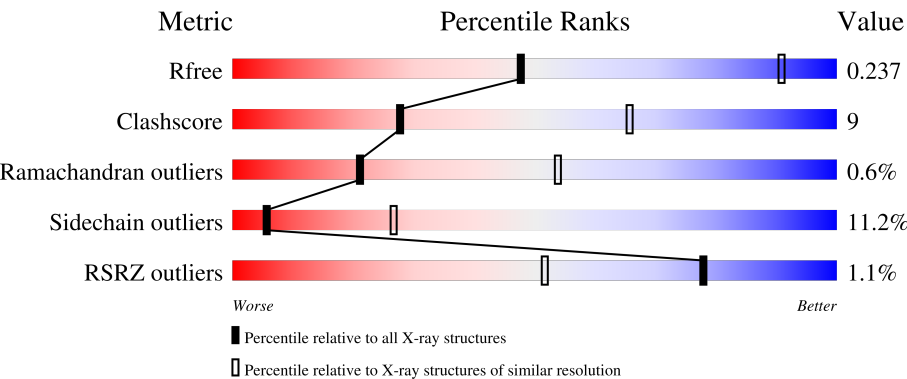
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.55 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	1410 (3.62-3.50)
Clashscore	190562	1480 (3.62-3.50)
Ramachandran outliers	187476	1440 (3.62-3.50)
Sidechain outliers	187428	1441 (3.62-3.50)
RSRZ outliers	180081	1409 (3.62-3.50)

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	337	% 63%26%5%6%
1	B	337	% 64%23%•9%
1	C	337	% 65%25%•7%
1	D	337	% 62%26%•9%
1	E	337	% 64%26%•7%

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Mol	Chain	Length	Quality of chain
1	F	337	% 61% 25% • 9%
2	G	14	7% 50% 50%
3	H	15	7% 80% 13% 7%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 16104 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 DNA double-strand break repair protein Mre11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	313	Total	C	N	O	S	0	0	0
			2601	1680	434	476	11			
1	A	317	Total	C	N	O	S	0	0	0
			2635	1698	444	482	11			
1	B	306	Total	C	N	O	S	0	0	0
			2546	1644	427	465	10			
1	D	307	Total	C	N	O	S	0	0	0
			2556	1650	430	466	10			
1	E	314	Total	C	N	O	S	0	0	0
			2605	1682	437	476	10			
1	F	305	Total	C	N	O	S	0	0	0
			2537	1640	428	459	10			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-3	ARG	-	expression tag	UNP Q58719
C	-2	GLY	-	expression tag	UNP Q58719
C	-1	SER	-	expression tag	UNP Q58719
C	0	HIS	-	expression tag	UNP Q58719
A	-3	ARG	-	expression tag	UNP Q58719
A	-2	GLY	-	expression tag	UNP Q58719
A	-1	SER	-	expression tag	UNP Q58719
A	0	HIS	-	expression tag	UNP Q58719
B	-3	ARG	-	expression tag	UNP Q58719
B	-2	GLY	-	expression tag	UNP Q58719
B	-1	SER	-	expression tag	UNP Q58719
B	0	HIS	-	expression tag	UNP Q58719
D	-3	ARG	-	expression tag	UNP Q58719
D	-2	GLY	-	expression tag	UNP Q58719
D	-1	SER	-	expression tag	UNP Q58719
D	0	HIS	-	expression tag	UNP Q58719
E	-3	ARG	-	expression tag	UNP Q58719

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-2	GLY	-	expression tag	UNP Q58719
E	-1	SER	-	expression tag	UNP Q58719
E	0	HIS	-	expression tag	UNP Q58719
F	-3	ARG	-	expression tag	UNP Q58719
F	-2	GLY	-	expression tag	UNP Q58719
F	-1	SER	-	expression tag	UNP Q58719
F	0	HIS	-	expression tag	UNP Q58719

- Molecule 2 is a DNA chain called DNA (5'-D(P*GP*CP*AP*CP*GP*TP*AP*GP*GP*AP*CP*AP*GP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	G	14	Total	C	N	O	P	0	0	0
			290	136	59	81	14			

- Molecule 3 is a DNA chain called DNA (5'-D(P*CP*TP*GP*TP*CP*CP*TP*AP*CP*GP*TP*GP*CP*CP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	15	Total	C	N	O	P	0	0	0
			302	144	51	92	15			

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	2	Total	Mg	0	0
			2	2		
4	A	2	Total	Mg	0	0
			2	2		
4	B	2	Total	Mg	0	0
			2	2		
4	D	2	Total	Mg	0	0
			2	2		
4	E	2	Total	Mg	0	0
			2	2		
4	F	2	Total	Mg	0	0
			2	2		

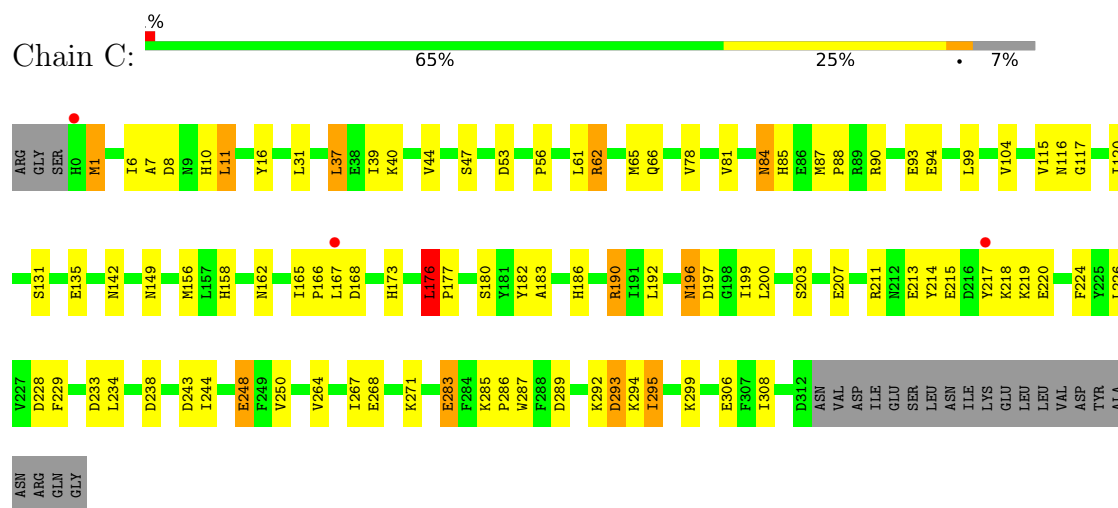
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	2	Total 2	O 2	0	0
5	A	7	Total 7	O 7	0	0
5	B	5	Total 5	O 5	0	0
5	D	2	Total 2	O 2	0	0
5	E	3	Total 3	O 3	0	0
5	F	1	Total 1	O 1	0	0

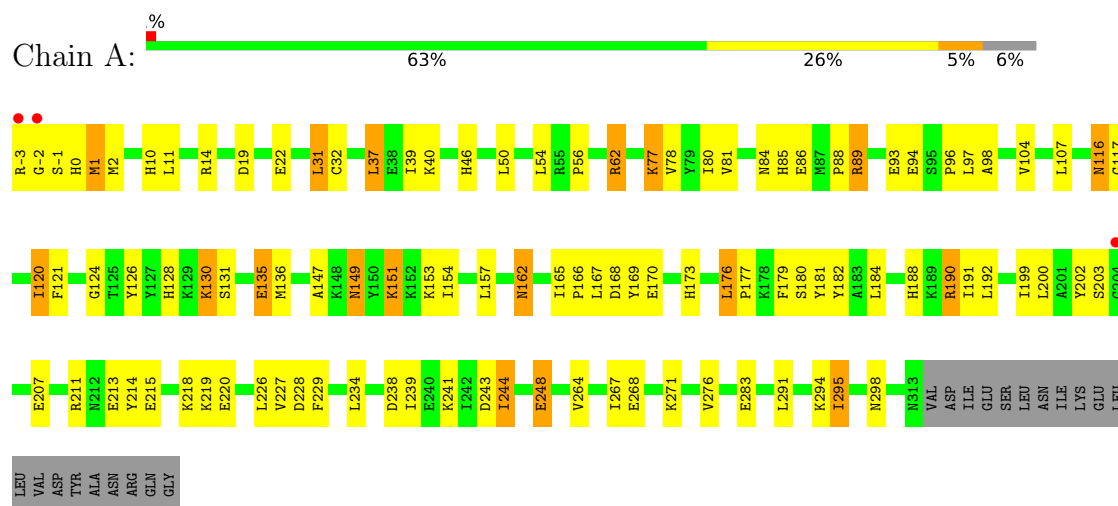
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: DNA double-strand break repair protein Mre11

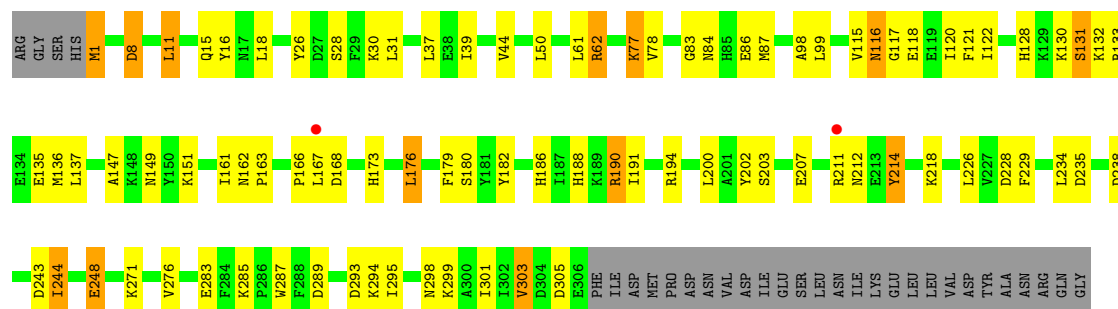


- Molecule 1: DNA double-strand break repair protein Mre11

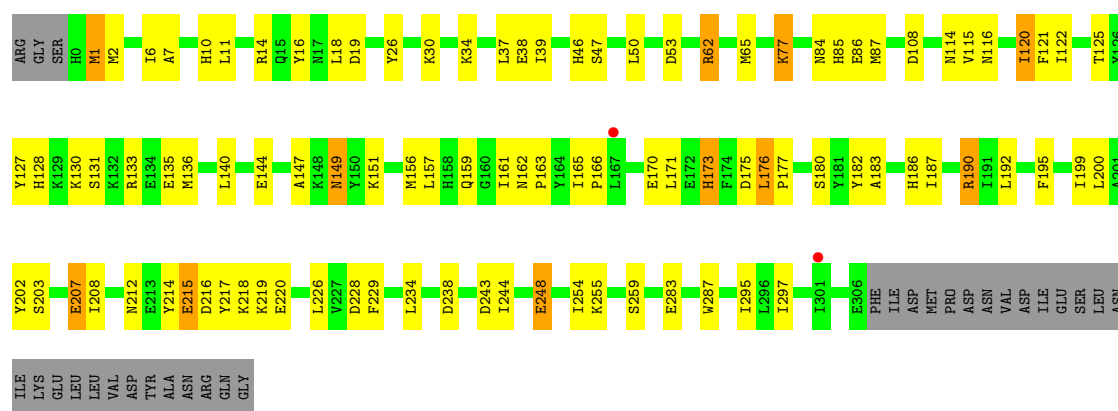


- Molecule 1: DNA double-strand break repair protein Mre11

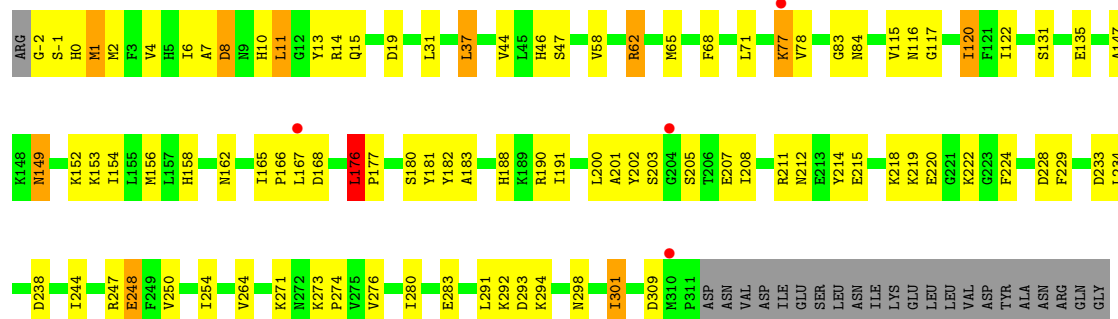




- Molecule 1: DNA double-strand break repair protein Mre11

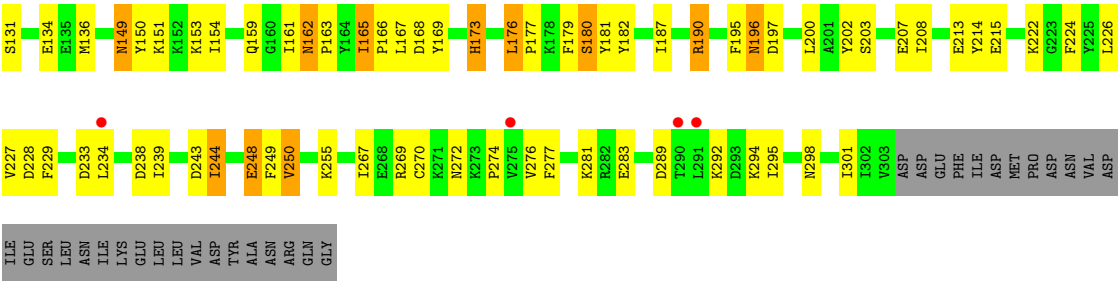


- Molecule 1: DNA double-strand break repair protein Mre11

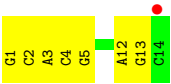


- Molecule 1: DNA double-strand break repair protein Mre11

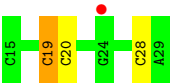
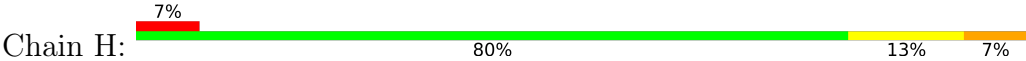




• Molecule 2: DNA (5'-D(P*GP*CP*AP*CP*GP*TP*AP*GP*GP*AP*CP*AP*GP*C)-3')



• Molecule 3: DNA (5'-D(P*CP*TP*GP*TP*CP*CP*TP*AP*CP*GP*TP*GP*CP*CP*A)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	91.53Å 185.62Å 106.21Å 90.00° 99.86° 90.00°	Depositor
Resolution (Å)	42.42 – 3.55 42.42 – 3.55	Depositor EDS
% Data completeness (in resolution range)	94.1 (42.42-3.55) 94.1 (42.42-3.55)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.31 (at 3.57Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8_1069)	Depositor
R, R_{free}	0.184 , 0.236 0.189 , 0.237	Depositor DCC
R_{free} test set	1996 reflections (4.75%)	wwPDB-VP
Wilson B-factor (Å ²)	79.8	Xtriage
Anisotropy	0.097	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 99.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	16104	wwPDB-VP
Average B, all atoms (Å ²)	144.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.00% 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

Bond lengths and bond angles in the following residue types are not validated in this section: MG

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	0.43	0/2692	0.89	9/3613 (0.2%)
1	B	0.41	0/2600	0.88	4/3489 (0.1%)
1	C	0.44	0/2657	0.90	6/3567 (0.2%)
1	D	0.39	0/2611	0.91	9/3504 (0.3%)
1	E	0.41	0/2662	0.90	5/3574 (0.1%)
1	F	0.36	0/2592	0.90	5/3478 (0.1%)
2	G	0.23	0/326	0.99	0/501
3	H	0.30	0/336	1.25	4/515 (0.8%)
All	All	0.40	0/16476	0.91	42/22241 (0.2%)

There are no bond length outliers.

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	176	LEU	CA-C-N	9.73	131.00	120.11
1	C	176	LEU	C-N-CA	9.73	131.00	120.11
1	E	176	LEU	CA-C-N	9.45	130.70	120.11
1	E	176	LEU	C-N-CA	9.45	130.70	120.11
1	A	176	LEU	CA-C-N	8.11	129.20	120.11
1	A	176	LEU	C-N-CA	8.11	129.20	120.11
1	F	87	MET	CA-C-N	8.10	128.15	119.89
1	F	87	MET	C-N-CA	8.10	128.15	119.89
1	D	176	LEU	CA-C-N	7.80	128.85	120.11
1	D	176	LEU	C-N-CA	7.80	128.85	120.11
1	D	165	ILE	CA-C-N	7.60	129.34	119.84
1	D	165	ILE	C-N-CA	7.60	129.34	119.84
1	D	87	MET	CA-C-N	6.87	126.90	119.89
1	D	87	MET	C-N-CA	6.87	126.90	119.89
1	E	162	ASN	N-CA-C	6.79	122.08	113.25
1	F	165	ILE	CA-C-N	6.74	128.27	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	165	ILE	C-N-CA	6.74	128.27	119.84
1	E	165	ILE	CA-C-N	6.52	128.00	119.84
1	E	165	ILE	C-N-CA	6.52	128.00	119.84
1	D	162	ASN	N-CA-C	6.41	121.58	113.25
1	A	165	ILE	CA-C-N	6.35	127.78	119.84
1	A	165	ILE	C-N-CA	6.35	127.78	119.84
1	A	0	HIS	N-CA-C	-6.35	104.23	112.23
3	H	19	DC	C5'-C4'-C3'	-6.24	105.55	114.90
1	A	56	PRO	CA-C-N	6.20	126.17	119.78
1	A	56	PRO	C-N-CA	6.20	126.17	119.78
1	A	162	ASN	N-CA-C	6.03	121.09	113.25
1	C	165	ILE	CA-C-N	5.83	127.13	119.84
1	C	165	ILE	C-N-CA	5.83	127.13	119.84
1	C	162	ASN	N-CA-C	5.58	120.50	113.25
1	C	186	HIS	N-CA-C	5.55	119.53	112.87
3	H	19	DC	N1-C1'-C2'	-5.40	105.41	113.50
1	B	303	VAL	N-CA-C	5.30	115.22	107.37
1	B	162	ASN	N-CA-C	5.26	120.09	113.25
1	D	120	ILE	N-CA-C	5.24	116.27	108.46
1	A	214	TYR	N-CA-C	5.14	116.57	111.07
1	B	176	LEU	CA-C-N	5.13	125.86	120.11
1	B	176	LEU	C-N-CA	5.13	125.86	120.11
3	H	28	DC	C4'-C3'-O3'	-5.13	102.31	110.00
1	D	255	LYS	N-CA-C	-5.11	108.07	114.56
1	F	162	ASN	N-CA-C	5.04	119.80	113.25
3	H	28	DC	C5'-C4'-O4'	5.03	116.94	109.40

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	2635	0	2650	60	0
1	B	2546	0	2572	40	0
1	C	2601	0	2618	43	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	2556	0	2579	44	0
1	E	2605	0	2620	47	0
1	F	2537	0	2570	62	0
2	G	290	0	156	7	0
3	H	302	0	170	1	0
4	A	2	0	0	0	0
4	B	2	0	0	0	0
4	C	2	0	0	0	0
4	D	2	0	0	0	0
4	E	2	0	0	0	0
4	F	2	0	0	0	0
5	A	7	0	0	1	0
5	B	5	0	0	0	0
5	C	2	0	0	0	0
5	D	2	0	0	0	0
5	E	3	0	0	0	0
5	F	1	0	0	0	0
All	All	16104	0	15935	287	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 9.

All (287) 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:B:1:MET:HG3	1:B:115:VAL:HG21	1.57	0.84
1:E:65:MET:HE1	1:F:65:MET:HE1	1.66	0.77
1:F:179:PHE:HB2	1:F:182:TYR:HE1	1.53	0.74
2:G:2:DC:H2'	2:G:3:DA:H8	1.54	0.73
1:F:127:TYR:HA	1:F:159:GLN:HE22	1.55	0.71
1:A:14:ARG:NH2	1:A:19:ASP:OD2	2.26	0.69
1:F:31:LEU:HD23	1:F:244:ILE:HD12	1.75	0.69
1:B:131:SER:HB2	1:B:132:LYS:HD2	1.75	0.69
2:G:2:DC:H2'	2:G:3:DA:C8	2.27	0.68
1:F:2:MET:HA	1:F:228:ASP:HA	1.76	0.67
1:F:121:PHE:HB2	1:F:150:TYR:HD1	1.60	0.67
1:C:44:VAL:HB	1:C:78:VAL:HG22	1.77	0.67
1:A:14:ARG:HG3	1:A:22:GLU:HB2	1.78	0.66
1:C:84:ASN:OD1	1:C:158:HIS:NE2	2.29	0.64
1:E:8:ASP:OD2	1:E:10:HIS:NE2	2.28	0.64
1:C:88:PRO:HB3	1:C:93:GLU:HB3	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:190:ARG:NH2	1:F:243:ASP:OD2	2.28	0.64
1:F:182:TYR:HB2	1:F:200:LEU:HA	1.80	0.63
1:F:44:VAL:HB	1:F:78:VAL:HG22	1.81	0.63
1:A:191:ILE:HB	1:A:202:TYR:HB2	1.81	0.63
1:F:179:PHE:HB2	1:F:182:TYR:CE1	2.34	0.62
1:D:228:ASP:OD1	1:D:229:PHE:N	2.33	0.61
1:D:219:LYS:HG2	1:D:220:GLU:HG2	1.82	0.61
1:C:65:MET:HE1	1:D:65:MET:HE1	1.83	0.60
1:F:289:ASP:HA	1:F:292:LYS:HB3	1.84	0.60
1:F:274:PRO:HG2	1:F:295:ILE:HG13	1.82	0.60
1:E:68:PHE:HD1	1:E:71:LEU:HD12	1.67	0.60
1:A:153:LYS:HG3	1:A:179:PHE:HA	1.83	0.59
1:E:182:TYR:HB2	1:E:200:LEU:HA	1.84	0.59
1:E:264:VAL:HG22	1:E:291:LEU:HD11	1.85	0.58
1:A:-3:ARG:O	1:A:-1:SER:OG	2.22	0.58
1:E:44:VAL:HB	1:E:78:VAL:HG22	1.84	0.58
1:F:187:ILE:HD12	1:F:202:TYR:CZ	2.39	0.58
1:F:39:ILE:HD13	1:F:226:LEU:HD22	1.86	0.58
1:A:-2:GLY:N	1:F:166:PRO:O	2.37	0.57
1:F:154:ILE:HG23	1:F:181:TYR:HB3	1.87	0.57
1:F:161:ILE:HG22	1:F:163:PRO:HD2	1.86	0.57
1:C:90:ARG:NH2	1:C:93:GLU:OE2	2.37	0.57
1:B:228:ASP:OD1	1:B:229:PHE:N	2.37	0.57
1:E:8:ASP:OD2	1:E:188:HIS:NE2	2.38	0.57
1:A:1:MET:HE3	1:A:120:ILE:HD13	1.86	0.57
1:C:306:GLU:HB2	1:B:62:ARG:HD2	1.86	0.57
1:D:215:GLU:OE1	1:D:216:ASP:N	2.38	0.57
1:C:6:ILE:HG13	1:C:224:PHE:HB3	1.86	0.57
1:B:194:ARG:NE	1:B:235:ASP:OD1	2.37	0.56
1:A:211:ARG:NH2	5:A:502:HOH:O	2.38	0.56
1:B:39:ILE:HD13	1:B:226:LEU:HD22	1.87	0.56
1:D:62:ARG:HH11	1:D:62:ARG:CG	2.19	0.55
1:C:182:TYR:HB2	1:C:200:LEU:HA	1.88	0.55
1:C:293:ASP:N	1:C:293:ASP:OD1	2.40	0.55
1:A:98:ALA:HB1	1:B:62:ARG:NH2	2.21	0.55
1:D:127:TYR:HA	1:D:159:GLN:HE22	1.71	0.55
1:F:248:GLU:OE2	1:F:272:ASN:N	2.36	0.55
1:E:147:ALA:HB1	1:E:153:LYS:HD2	1.89	0.55
1:C:228:ASP:OD1	1:C:229:PHE:N	2.40	0.54
1:A:128:HIS:HB2	1:A:136:MET:HE2	1.89	0.54
1:A:190:ARG:NH2	1:A:243:ASP:OD1	2.41	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:18:LEU:HD11	1:D:297:ILE:HD12	1.89	0.54
1:C:39:ILE:HD13	1:C:226:LEU:HD22	1.90	0.54
1:E:0:HIS:HB2	1:E:1:MET:HE2	1.90	0.54
1:E:228:ASP:OD1	1:E:229:PHE:N	2.40	0.54
1:F:182:TYR:HD2	1:F:200:LEU:HD12	1.73	0.54
1:F:153:LYS:HG3	1:F:179:PHE:HA	1.90	0.54
1:A:228:ASP:OD1	1:A:229:PHE:N	2.40	0.53
1:E:215:GLU:O	1:E:219:LYS:HB3	2.08	0.53
1:B:8:ASP:OD2	1:B:188:HIS:NE2	2.41	0.53
1:F:228:ASP:OD1	1:F:229:PHE:N	2.41	0.53
1:D:163:PRO:HG3	1:D:173:HIS:CD2	2.43	0.53
1:B:44:VAL:HB	1:B:78:VAL:HG22	1.91	0.53
1:B:26:TYR:O	1:B:30:LYS:HG3	2.09	0.53
1:B:61:LEU:HD22	1:B:99:LEU:HD12	1.91	0.53
1:B:116:ASN:OD1	1:B:116:ASN:N	2.40	0.52
1:D:121:PHE:CE2	1:D:147:ALA:HB2	2.45	0.52
1:E:68:PHE:CD1	1:E:71:LEU:HD12	2.45	0.52
1:E:248:GLU:OE2	1:E:271:LYS:N	2.41	0.52
1:D:218:LYS:NZ	1:D:248:GLU:OE1	2.39	0.52
1:E:14:ARG:NH2	1:E:19:ASP:OD2	2.43	0.52
1:F:8:ASP:OD2	1:F:10:HIS:NE2	2.32	0.52
1:F:59:LYS:O	1:F:63:ILE:HG13	2.10	0.52
1:D:16:TYR:HB2	1:D:18:LEU:HD13	1.92	0.52
1:F:29:PHE:O	1:F:33:ILE:HG12	2.10	0.51
1:C:16:TYR:CD1	1:C:299:LYS:HE3	2.45	0.51
1:A:227:VAL:HG22	1:A:239:ILE:HG12	1.92	0.51
1:A:62:ARG:NH1	1:B:98:ALA:HB1	2.26	0.51
1:A:192:LEU:HD11	1:A:199:ILE:HD11	1.92	0.51
1:C:292:LYS:HE3	1:A:135:GLU:HG3	1.92	0.51
1:B:128:HIS:HB2	1:B:136:MET:HE2	1.92	0.51
1:D:156:MET:HG2	1:D:183:ALA:HB3	1.93	0.50
1:F:127:TYR:HA	1:F:159:GLN:NE2	2.22	0.50
1:F:128:HIS:HB2	1:F:136:MET:HE2	1.92	0.50
1:A:31:LEU:HB3	1:A:244:ILE:HD12	1.93	0.50
1:D:62:ARG:HH11	1:D:62:ARG:HG3	1.75	0.50
2:G:4:DC:H2''	2:G:5:DG:C8	2.46	0.50
1:B:50:LEU:N	1:B:86:GLU:OE2	2.45	0.50
1:E:276:VAL:HG13	1:E:298:ASN:HA	1.94	0.49
1:A:181:TYR:CE2	1:A:227:VAL:HG21	2.47	0.49
1:F:167:LEU:HD12	1:F:168:ASP:HB3	1.95	0.49
1:E:205:SER:O	1:E:222:LYS:HD3	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:219:LYS:HG2	1:E:220:GLU:HG2	1.95	0.49
1:C:286:PRO:HB3	1:A:126:TYR:CZ	2.47	0.49
1:E:177:PRO:O	1:E:182:TYR:OH	2.25	0.49
1:C:61:LEU:HD22	1:C:99:LEU:HD12	1.95	0.48
1:F:173:HIS:HB2	1:F:195:PHE:CE2	2.48	0.48
1:C:176:LEU:HA	1:C:177:PRO:HD3	1.61	0.48
1:D:39:ILE:HD13	1:D:226:LEU:HD22	1.94	0.48
1:C:167:LEU:HA	1:C:168:ASP:HA	1.57	0.48
1:A:116:ASN:HA	1:A:117:GLY:HA2	1.57	0.48
1:A:276:VAL:HG13	1:A:298:ASN:HA	1.95	0.48
1:C:94:GLU:OE1	1:D:62:ARG:NH1	2.47	0.48
1:B:115:VAL:HG23	1:B:118:GLU:H	1.77	0.48
1:F:149:ASN:OD1	1:F:149:ASN:N	2.38	0.48
1:C:215:GLU:O	1:C:219:LYS:HB3	2.14	0.48
1:D:77:LYS:HD2	1:D:77:LYS:HA	1.57	0.48
1:C:192:LEU:HD11	1:C:199:ILE:HD11	1.95	0.47
1:C:219:LYS:HG2	1:C:220:GLU:HG2	1.95	0.47
1:A:37:LEU:O	1:A:40:LYS:HG2	2.14	0.47
1:A:248:GLU:OE2	1:A:271:LYS:N	2.46	0.47
1:D:149:ASN:OD1	1:D:149:ASN:N	2.36	0.47
1:C:248:GLU:OE2	1:C:271:LYS:N	2.47	0.47
1:E:1:MET:HG2	1:E:115:VAL:HG21	1.95	0.47
1:B:276:VAL:HG13	1:B:298:ASN:HA	1.96	0.47
1:F:249:PHE:HB3	1:F:277:PHE:CE2	2.49	0.47
1:A:267:ILE:HD13	1:A:295:ILE:HD11	1.96	0.47
1:F:153:LYS:H	1:F:180:SER:HG	1.62	0.47
1:F:167:LEU:HA	1:F:168:ASP:HA	1.49	0.47
1:B:116:ASN:HA	1:B:117:GLY:HA2	1.51	0.47
1:C:62:ARG:O	1:C:66:GLN:HB2	2.15	0.47
1:A:10:HIS:NE2	1:A:188:HIS:CE1	2.83	0.47
1:A:182:TYR:HB2	1:A:200:LEU:HA	1.97	0.47
1:D:108:ASP:OD1	1:D:125:THR:OG1	2.29	0.47
1:E:156:MET:HG2	1:E:183:ALA:HB3	1.95	0.47
1:E:13:TYR:CE2	1:E:15:GLN:HB2	2.50	0.46
1:F:2:MET:HG2	1:F:228:ASP:HB2	1.96	0.46
1:F:222:LYS:O	1:F:244:ILE:N	2.48	0.46
1:A:154:ILE:HG23	1:A:181:TYR:HB3	1.97	0.46
1:A:-3:ARG:O	1:F:167:LEU:HD13	2.15	0.46
1:E:6:ILE:HG13	1:E:224:PHE:HB3	1.97	0.46
1:D:14:ARG:NH2	1:D:19:ASP:OD2	2.48	0.46
1:D:140:LEU:HD21	1:D:157:LEU:HD11	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:6:ILE:O	1:F:46:HIS:HA	2.16	0.46
1:C:7:ALA:HB2	1:C:47:SER:HB2	1.98	0.46
1:A:10:HIS:CE1	1:A:85:HIS:CD2	3.04	0.46
1:A:88:PRO:HB3	1:A:93:GLU:HB3	1.98	0.46
1:D:187:ILE:HD12	1:D:202:TYR:CE2	2.51	0.46
2:G:1:DG:H2'	2:G:1:DG:N3	2.30	0.46
1:C:264:VAL:O	1:C:268:GLU:HG2	2.16	0.46
1:E:176:LEU:HA	1:E:177:PRO:HD3	1.61	0.46
1:A:264:VAL:HG22	1:A:291:LEU:HD11	1.98	0.46
1:D:207:GLU:OE1	1:D:208:ILE:N	2.49	0.46
1:D:254:ILE:HA	1:D:259:SER:OG	2.16	0.46
1:B:179:PHE:HB2	1:B:182:TYR:CE1	2.50	0.46
2:G:12:DA:H2''	2:G:13:DG:H5''	1.98	0.45
1:B:167:LEU:HA	1:B:168:ASP:HA	1.59	0.45
1:D:127:TYR:HA	1:D:159:GLN:NE2	2.31	0.45
1:D:144:GLU:OE1	1:D:177:PRO:HB3	2.16	0.45
1:D:192:LEU:HD11	1:D:199:ILE:HD11	1.97	0.45
1:F:270:CYS:SG	1:F:274:PRO:HB3	2.57	0.45
1:A:-3:ARG:HG2	1:F:169:TYR:H	1.80	0.45
1:E:116:ASN:HA	1:E:117:GLY:HA2	1.49	0.45
1:F:276:VAL:HG13	1:F:298:ASN:HA	1.98	0.45
1:C:78:VAL:HG12	1:C:104:VAL:HG13	1.98	0.45
1:A:130:LYS:HG3	1:A:170:GLU:HA	1.99	0.45
1:E:254:ILE:HD12	1:E:280:ILE:HG12	1.97	0.45
1:C:190:ARG:NH2	1:C:243:ASP:OD1	2.49	0.45
1:F:196:ASN:HB3	1:F:197:ASP:H	1.67	0.45
1:C:56:PRO:HG2	1:C:61:LEU:HD21	1.99	0.45
1:A:149:ASN:OD1	1:A:149:ASN:N	2.36	0.45
1:F:208:ILE:HG23	1:F:213:GLU:HB2	1.99	0.45
1:E:11:LEU:HD12	1:E:11:LEU:HA	1.77	0.45
1:F:162:ASN:HB2	1:F:169:TYR:CE2	2.52	0.45
1:A:39:ILE:HD13	1:A:226:LEU:HD22	1.99	0.44
1:B:121:PHE:CE2	1:B:147:ALA:HB2	2.53	0.44
1:F:97:LEU:HB3	1:F:106:ILE:HD11	1.98	0.44
1:A:46:HIS:HB3	1:A:80:ILE:HG13	1.99	0.44
1:A:78:VAL:HG12	1:A:104:VAL:HG13	1.99	0.44
1:B:77:LYS:HD2	1:B:77:LYS:HA	1.63	0.44
1:D:7:ALA:HB2	1:D:47:SER:HB2	1.99	0.44
1:F:173:HIS:HB2	1:F:195:PHE:CZ	2.52	0.44
1:C:177:PRO:O	1:C:182:TYR:OH	2.30	0.44
1:B:83:GLY:O	1:B:87:MET:HG2	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:6:ILE:O	1:E:46:HIS:HA	2.17	0.44
1:F:187:ILE:HD12	1:F:202:TYR:CE2	2.52	0.44
1:C:167:LEU:HD12	1:C:168:ASP:HA	1.99	0.44
1:B:28:SER:HB2	1:B:244:ILE:HG21	2.00	0.44
1:F:244:ILE:HD13	1:F:244:ILE:HA	1.84	0.44
1:D:6:ILE:O	1:D:46:HIS:HA	2.17	0.44
1:E:191:ILE:HB	1:E:202:TYR:HB2	1.99	0.44
1:A:167:LEU:HA	1:A:168:ASP:HA	1.57	0.44
1:D:62:ARG:HH11	1:D:62:ARG:CB	2.31	0.44
1:E:2:MET:HA	1:E:228:ASP:HA	2.00	0.44
1:F:176:LEU:HA	1:F:177:PRO:HD3	1.79	0.44
1:F:227:VAL:HG22	1:F:239:ILE:HG12	2.00	0.44
1:B:293:ASP:N	1:B:293:ASP:OD2	2.51	0.44
1:D:190:ARG:NH2	1:D:243:ASP:OD1	2.51	0.44
1:C:1:MET:SD	1:C:115:VAL:HG21	2.58	0.43
1:B:191:ILE:HB	1:B:202:TYR:HB2	1.99	0.43
1:D:2:MET:HA	1:D:228:ASP:HA	1.98	0.43
1:D:26:TYR:O	1:D:30:LYS:HG3	2.18	0.43
1:C:116:ASN:HA	1:C:117:GLY:HA2	1.54	0.43
1:A:177:PRO:O	1:A:182:TYR:OH	2.32	0.43
1:B:285:LYS:NZ	1:B:289:ASP:OD1	2.43	0.43
1:D:34:LYS:HE2	1:D:38:GLU:OE2	2.18	0.43
1:E:62:ARG:HE	1:E:62:ARG:HB3	1.56	0.43
1:F:120:ILE:HD11	1:F:154:ILE:HD11	2.00	0.43
3:H:19:DC:H2'	3:H:20:DC:C6	2.53	0.43
1:E:7:ALA:HB2	1:E:47:SER:HB2	2.00	0.43
1:E:208:ILE:HD11	1:E:214:TYR:HD1	1.83	0.43
1:C:37:LEU:O	1:C:40:LYS:HG2	2.19	0.43
1:A:89:ARG:NH2	2:G:4:DC:O3'	2.51	0.43
1:D:10:HIS:CE1	1:D:85:HIS:CD2	3.06	0.43
1:F:250:VAL:HG21	1:F:267:ILE:HG12	2.00	0.43
1:E:167:LEU:HA	1:E:168:ASP:HA	1.60	0.43
1:E:274:PRO:HD2	1:E:294:LYS:O	2.18	0.43
1:C:214:TYR:O	1:C:217:TYR:HB3	2.19	0.43
1:B:161:ILE:HG22	1:B:163:PRO:HD2	2.01	0.43
1:A:-3:ARG:NE	1:F:169:TYR:O	2.51	0.43
1:A:264:VAL:O	1:A:268:GLU:HG2	2.19	0.43
1:B:190:ARG:NH2	1:B:243:ASP:OD2	2.51	0.43
1:F:77:LYS:HA	1:F:77:LYS:HD2	1.66	0.43
1:F:208:ILE:HD13	1:F:214:TYR:HA	2.00	0.43
1:E:37:LEU:HD13	1:E:37:LEU:HA	1.82	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:241:LYS:HE2	1:F:134:GLU:OE2	2.19	0.42
1:F:255:LYS:HA	1:F:281:LYS:HD2	2.00	0.42
1:A:97:LEU:HD23	1:A:97:LEU:HA	1.84	0.42
1:A:151:LYS:H	1:A:151:LYS:HG2	1.44	0.42
1:B:248:GLU:OE2	1:B:271:LYS:N	2.51	0.42
1:B:132:LYS:HD2	1:B:132:LYS:N	2.35	0.42
1:E:-2:GLY:HA3	1:E:-1:SER:HA	1.57	0.42
2:G:12:DA:H2''	2:G:13:DG:C5'	2.49	0.42
1:E:1:MET:HB2	1:E:1:MET:HE3	1.87	0.42
1:E:58:VAL:HG11	1:F:96:PRO:HA	2.01	0.42
1:C:156:MET:HG2	1:C:183:ALA:HB3	2.01	0.42
1:D:182:TYR:HB2	1:D:200:LEU:HA	2.02	0.42
1:E:120:ILE:HG13	1:E:152:LYS:O	2.19	0.42
1:F:153:LYS:N	1:F:180:SER:OG	2.49	0.42
1:C:283:GLU:HG2	1:A:94:GLU:HG2	2.01	0.42
1:E:247:ARG:NH1	1:E:273:LYS:O	2.51	0.42
1:D:133:ARG:NH2	1:D:175:ASP:OD2	2.43	0.41
1:C:10:HIS:CE1	1:C:85:HIS:CD2	3.08	0.41
1:A:32:CYS:SG	1:A:244:ILE:HG13	2.61	0.41
1:D:128:HIS:HB2	1:D:136:MET:HE2	2.02	0.41
1:A:120:ILE:HD11	1:A:154:ILE:HD11	2.00	0.41
1:F:7:ALA:HA	1:F:46:HIS:CE1	2.56	0.41
1:C:267:ILE:HD13	1:C:295:ILE:HD11	2.01	0.41
1:B:182:TYR:HB2	1:B:200:LEU:HA	2.01	0.41
1:D:173:HIS:HB2	1:D:195:PHE:CE2	2.55	0.41
1:E:83:GLY:HA2	1:E:158:HIS:CD2	2.55	0.41
1:B:11:LEU:HA	1:B:11:LEU:HD12	1.83	0.41
1:B:77:LYS:HD2	1:B:78:VAL:H	1.86	0.41
1:B:303:VAL:HG12	1:B:305:ASP:H	1.86	0.41
1:F:165:ILE:HA	1:F:166:PRO:HD2	1.88	0.41
1:A:77:LYS:HD2	1:A:77:LYS:HA	1.69	0.41
1:D:10:HIS:CE1	1:D:85:HIS:HD2	2.37	0.41
1:C:196:ASN:HB3	1:C:197:ASP:H	1.53	0.41
1:A:162:ASN:HB2	1:A:169:TYR:CD2	2.56	0.41
1:F:190:ARG:HH12	1:F:224:PHE:C	2.28	0.41
1:C:87:MET:HE3	1:C:87:MET:HB3	1.93	0.41
1:B:133:ARG:O	1:B:137:LEU:HG	2.21	0.41
1:D:50:LEU:HB3	1:D:86:GLU:OE2	2.20	0.41
1:C:11:LEU:HD12	1:C:11:LEU:HA	1.76	0.41
1:C:116:ASN:OD1	1:C:116:ASN:N	2.44	0.41
1:A:-3:ARG:HB3	1:F:167:LEU:HD12	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:-2:GLY:HA3	1:A:-1:SER:HA	1.86	0.41
1:A:50:LEU:N	1:A:86:GLU:OE2	2.53	0.41
1:A:244:ILE:HA	1:A:244:ILE:HD13	1.77	0.41
1:B:163:PRO:HG3	1:B:173:HIS:HB3	2.01	0.41
1:D:1:MET:HE2	1:D:115:VAL:HG21	2.03	0.41
1:D:116:ASN:OD1	1:D:116:ASN:N	2.41	0.41
1:E:301:ILE:H	1:E:301:ILE:HG13	1.73	0.41
1:A:107:LEU:HD13	1:A:124:GLY:N	2.36	0.41
1:F:87:MET:HA	1:F:88:PRO:HD3	1.83	0.41
1:A:54:LEU:O	1:A:96:PRO:HD3	2.21	0.40
1:A:219:LYS:HG2	1:A:220:GLU:HG2	2.02	0.40
1:D:214:TYR:O	1:D:217:TYR:HB3	2.21	0.40
1:E:77:LYS:HD2	1:E:77:LYS:HA	1.57	0.40
1:E:149:ASN:OD1	1:E:149:ASN:N	2.39	0.40
1:E:154:ILE:HG23	1:E:181:TYR:HB3	2.03	0.40
1:C:285:LYS:HZ3	1:C:289:ASP:CG	2.29	0.40
1:B:16:TYR:CD1	1:B:299:LYS:HE3	2.57	0.40
1:B:211:ARG:O	1:B:214:TYR:HB3	2.22	0.40
1:E:292:LYS:HG3	1:E:293:ASP:N	2.35	0.40
1:B:15:GLN:HG2	1:B:16:TYR:CD2	2.56	0.40
1:D:170:GLU:O	1:D:171:LEU:HD23	2.21	0.40
1:A:121:PHE:CE2	1:A:147:ALA:HB2	2.56	0.40
1:E:191:ILE:O	1:E:201:ALA:HA	2.20	0.40
1:A:2:MET:HE2	1:A:226:LEU:HD11	2.04	0.40
1:A:157:LEU:HB2	1:A:184:LEU:HD22	2.03	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	315/337 (94%)	299 (95%)	14 (4%)	2 (1%)	21 54

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	304/337 (90%)	290 (95%)	12 (4%)	2 (1%)	18	51
1	C	311/337 (92%)	293 (94%)	16 (5%)	2 (1%)	21	54
1	D	305/337 (90%)	290 (95%)	13 (4%)	2 (1%)	18	51
1	E	312/337 (93%)	295 (95%)	15 (5%)	2 (1%)	21	54
1	F	303/337 (90%)	289 (95%)	13 (4%)	1 (0%)	36	65
All	All	1850/2022 (92%)	1756 (95%)	83 (4%)	11 (1%)	21	54

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	186	HIS
1	E	309	ASP
1	C	166	PRO
1	B	186	HIS
1	C	213	GLU
1	A	213	GLU
1	A	166	PRO
1	B	166	PRO
1	D	166	PRO
1	E	166	PRO
1	F	301	ILE

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	290/308 (94%)	259 (89%)	31 (11%)	6	28
1	B	280/308 (91%)	246 (88%)	34 (12%)	5	24
1	C	286/308 (93%)	251 (88%)	35 (12%)	5	23
1	D	281/308 (91%)	250 (89%)	31 (11%)	6	27
1	E	286/308 (93%)	256 (90%)	30 (10%)	6	29
1	F	279/308 (91%)	249 (89%)	30 (11%)	6	28

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1702/1848 (92%)	1511 (89%)	191 (11%)	6 27

All (191) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	1	MET
1	C	8	ASP
1	C	11	LEU
1	C	31	LEU
1	C	37	LEU
1	C	53	ASP
1	C	62	ARG
1	C	81	VAL
1	C	84	ASN
1	C	120	ILE
1	C	131	SER
1	C	135	GLU
1	C	142	ASN
1	C	149	ASN
1	C	173	HIS
1	C	176	LEU
1	C	180	SER
1	C	190	ARG
1	C	196	ASN
1	C	203	SER
1	C	207	GLU
1	C	211	ARG
1	C	218	LYS
1	C	233	ASP
1	C	234	LEU
1	C	238	ASP
1	C	244	ILE
1	C	248	GLU
1	C	250	VAL
1	C	283	GLU
1	C	287	TRP
1	C	293	ASP
1	C	294	LYS
1	C	295	ILE
1	C	308	ILE
1	A	1	MET
1	A	11	LEU

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Mol	Chain	Res	Type
1	A	31	LEU
1	A	37	LEU
1	A	62	ARG
1	A	77	LYS
1	A	81	VAL
1	A	84	ASN
1	A	89	ARG
1	A	116	ASN
1	A	120	ILE
1	A	130	LYS
1	A	131	SER
1	A	135	GLU
1	A	149	ASN
1	A	151	LYS
1	A	173	HIS
1	A	176	LEU
1	A	180	SER
1	A	190	ARG
1	A	203	SER
1	A	207	GLU
1	A	215	GLU
1	A	218	LYS
1	A	234	LEU
1	A	238	ASP
1	A	244	ILE
1	A	248	GLU
1	A	283	GLU
1	A	294	LYS
1	A	295	ILE
1	B	1	MET
1	B	8	ASP
1	B	11	LEU
1	B	18	LEU
1	B	31	LEU
1	B	37	LEU
1	B	62	ARG
1	B	77	LYS
1	B	84	ASN
1	B	116	ASN
1	B	120	ILE
1	B	122	ILE
1	B	130	LYS

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Mol	Chain	Res	Type
1	B	131	SER
1	B	135	GLU
1	B	149	ASN
1	B	151	LYS
1	B	176	LEU
1	B	180	SER
1	B	190	ARG
1	B	203	SER
1	B	207	GLU
1	B	212	ASN
1	B	214	TYR
1	B	218	LYS
1	B	234	LEU
1	B	238	ASP
1	B	244	ILE
1	B	248	GLU
1	B	283	GLU
1	B	287	TRP
1	B	294	LYS
1	B	295	ILE
1	B	301	ILE
1	D	1	MET
1	D	11	LEU
1	D	37	LEU
1	D	53	ASP
1	D	62	ARG
1	D	77	LYS
1	D	84	ASN
1	D	114	ASN
1	D	120	ILE
1	D	122	ILE
1	D	130	LYS
1	D	131	SER
1	D	135	GLU
1	D	149	ASN
1	D	151	LYS
1	D	161	ILE
1	D	173	HIS
1	D	176	LEU
1	D	180	SER
1	D	190	ARG
1	D	203	SER

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Mol	Chain	Res	Type
1	D	207	GLU
1	D	212	ASN
1	D	215	GLU
1	D	234	LEU
1	D	238	ASP
1	D	244	ILE
1	D	248	GLU
1	D	283	GLU
1	D	287	TRP
1	D	295	ILE
1	E	1	MET
1	E	4	VAL
1	E	8	ASP
1	E	11	LEU
1	E	31	LEU
1	E	37	LEU
1	E	62	ARG
1	E	77	LYS
1	E	84	ASN
1	E	120	ILE
1	E	122	ILE
1	E	131	SER
1	E	135	GLU
1	E	149	ASN
1	E	176	LEU
1	E	180	SER
1	E	190	ARG
1	E	203	SER
1	E	207	GLU
1	E	211	ARG
1	E	212	ASN
1	E	218	LYS
1	E	233	ASP
1	E	234	LEU
1	E	238	ASP
1	E	244	ILE
1	E	248	GLU
1	E	250	VAL
1	E	283	GLU
1	E	301	ILE
1	F	8	ASP
1	F	11	LEU

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Mol	Chain	Res	Type
1	F	31	LEU
1	F	37	LEU
1	F	53	ASP
1	F	81	VAL
1	F	84	ASN
1	F	120	ILE
1	F	122	ILE
1	F	130	LYS
1	F	131	SER
1	F	149	ASN
1	F	151	LYS
1	F	173	HIS
1	F	176	LEU
1	F	180	SER
1	F	190	ARG
1	F	196	ASN
1	F	203	SER
1	F	207	GLU
1	F	215	GLU
1	F	233	ASP
1	F	234	LEU
1	F	238	ASP
1	F	244	ILE
1	F	248	GLU
1	F	250	VAL
1	F	269	ARG
1	F	283	GLU
1	F	294	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	75	ASN
1	C	85	HIS
1	C	162	ASN
1	C	253	ASN
1	A	0	HIS
1	A	162	ASN
1	B	66	GLN
1	B	162	ASN
1	B	253	ASN
1	D	85	HIS

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Mol	Chain	Res	Type
1	D	114	ASN
1	D	212	ASN
1	E	52	ASN
1	E	114	ASN
1	E	159	GLN
1	F	114	ASN
1	F	159	GLN
1	F	298	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 12 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	317/337 (94%)	-0.23	3 (0%) 81 56	70, 114, 179, 243	0
1	B	306/337 (90%)	-0.22	2 (0%) 84 61	66, 130, 208, 285	0
1	C	313/337 (92%)	-0.20	3 (0%) 79 54	78, 114, 178, 221	0
1	D	307/337 (91%)	-0.10	2 (0%) 84 61	78, 129, 232, 306	0
1	E	314/337 (93%)	-0.18	4 (1%) 75 47	79, 129, 208, 267	0
1	F	305/337 (90%)	0.21	5 (1%) 70 42	103, 191, 279, 339	0
2	G	14/14 (100%)	0.75	1 (7%) 22 13	131, 184, 233, 233	0
3	H	15/15 (100%)	0.78	1 (6%) 24 14	134, 174, 222, 393	0
All	All	1891/2051 (92%)	-0.11	21 (1%) 78 52	66, 132, 240, 393	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	H	24	DG	5.6
1	D	167	LEU	3.9
1	B	211	ARG	3.4
1	E	310	MET	3.4
1	C	167	LEU	3.0
1	A	-3	ARG	2.9
1	B	167	LEU	2.8
1	C	217	TYR	2.5
1	F	41	PRO	2.4
1	C	0	HIS	2.4
1	E	204	GLY	2.4
1	F	234	LEU	2.4
1	E	77	LYS	2.4
1	F	290	THR	2.3
1	A	-2	GLY	2.3
1	A	204	GLY	2.3

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Mol	Chain	Res	Type	RSRZ
2	G	14	DC	2.1
1	F	275	VAL	2.1
1	F	291	LEU	2.1
1	D	301	ILE	2.0
1	E	167	LEU	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	MG	B	402	1/1	0.94	0.10	77,77,77,77	0
4	MG	B	401	1/1	0.98	0.08	57,57,57,57	0
4	MG	A	401	1/1	0.98	0.08	54,54,54,54	0
4	MG	D	402	1/1	0.98	0.07	61,61,61,61	0
4	MG	E	402	1/1	0.98	0.09	102,102,102,102	0
4	MG	F	402	1/1	0.98	0.11	121,121,121,121	0
4	MG	D	401	1/1	0.99	0.09	45,45,45,45	0
4	MG	A	402	1/1	0.99	0.06	55,55,55,55	0
4	MG	E	401	1/1	0.99	0.05	70,70,70,70	0
4	MG	C	402	1/1	0.99	0.05	77,77,77,77	0
4	MG	F	401	1/1	0.99	0.05	77,77,77,77	0
4	MG	C	401	1/1	0.99	0.08	50,50,50,50	0

6.5 Other polymers [i](#)

There are no such residues in this entry.