



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 02:17 PM UTC

PDB ID : 5MS4 / pdb_00005ms4
Title : Kallikrein-related peptidase 8 leupeptin inhibitor complex
Authors : Debela, M.; Magdolen, V.; Skala, W.; Bode, W.; Brandstetter, H.; Goettig, P.
Deposited on : 2016-12-30
Resolution : 2.10 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
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

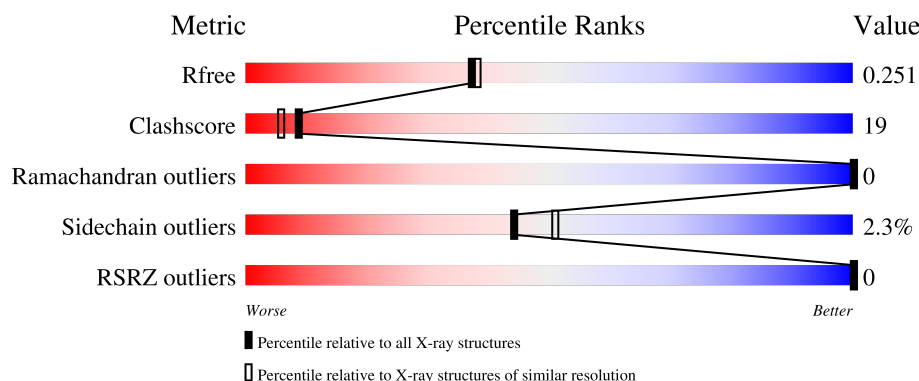
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	228	
1	B	228	
1	C	228	
1	D	228	
2	E	4	

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Mol	Chain	Length	Quality of chain
2	F	4	 50% 50%
2	G	4	 50% 25% 25%
2	H	4	 75% 25%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8120 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.

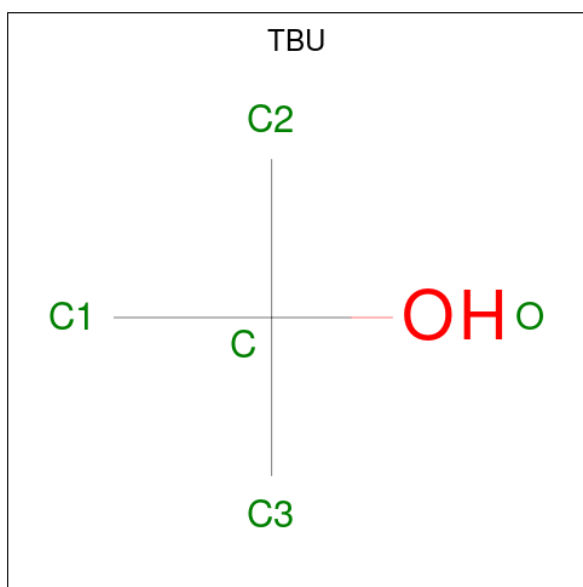
- Molecule 1 is a protein called Kallikrein-8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	228	Total	C	N	O	S	0	0	0
			1711	1061	303	332	15			
1	B	228	Total	C	N	O	S	0	0	0
			1711	1061	303	332	15			
1	C	228	Total	C	N	O	S	0	0	0
			1711	1061	303	332	15			
1	D	228	Total	C	N	O	S	0	0	0
			1711	1061	303	332	15			

- Molecule 2 is a protein called LEUPEPTIN.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	4	Total	C	N	O	0	0	0
			30	20	6	4			
2	F	4	Total	C	N	O	0	0	0
			30	20	6	4			
2	G	4	Total	C	N	O	0	0	0
			30	20	6	4			
2	H	4	Total	C	N	O	0	0	0
			30	20	6	4			

- Molecule 3 is TERTIARY-BUTYL ALCOHOL (CCD ID: TBU) (formula: C₄H₁₀O).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 5 4 1	0	0
3	A	1	Total C O 5 4 1	0	0
3	A	1	Total C O 5 4 1	0	0

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Ca 1 1	0	0
4	B	1	Total Ca 1 1	0	0
4	C	1	Total Ca 1 1	0	0
4	D	1	Total Ca 1 1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	277	Total O 277 277	0	0
5	B	276	Total O 276 276	0	0

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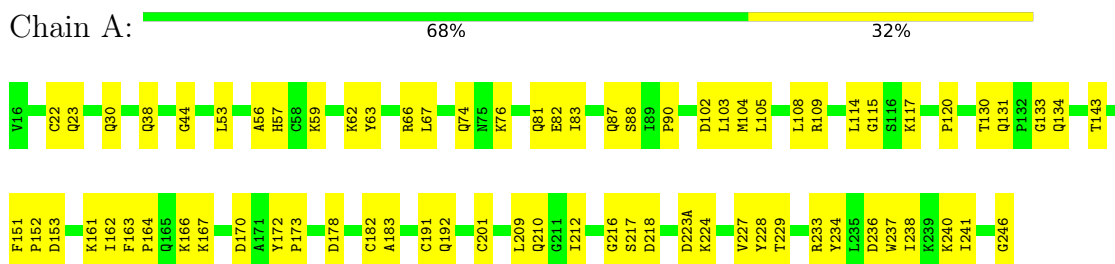
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	277	Total 277	O 277	0	0
5	D	293	Total 293	O 293	0	0
5	E	2	Total 2	O 2	0	0
5	F	4	Total 4	O 4	0	0
5	G	4	Total 4	O 4	0	0
5	H	4	Total 4	O 4	0	0

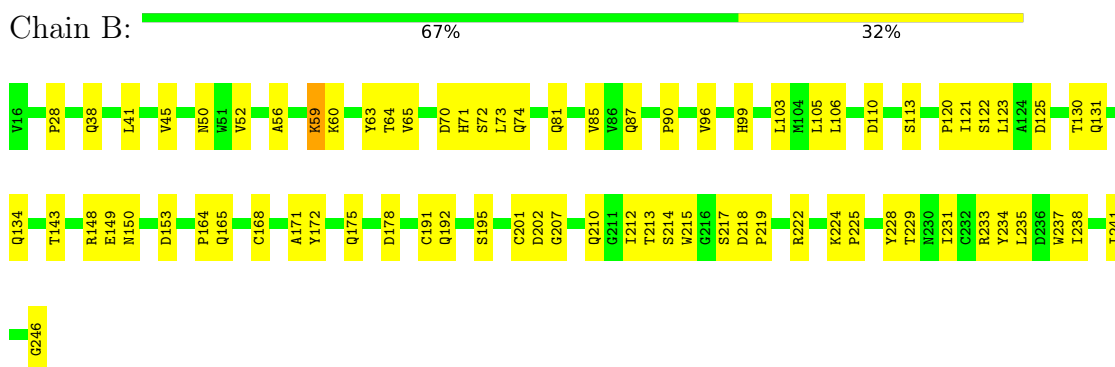
3 Residue-property plots

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.

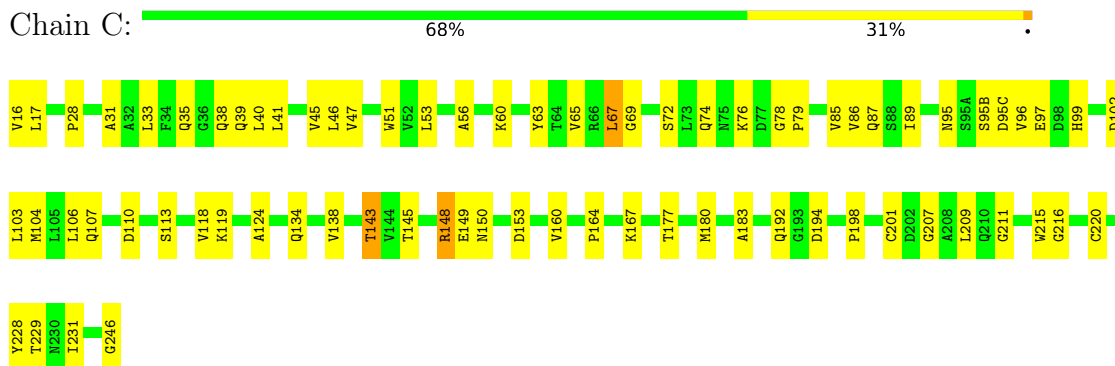
• Molecule 1: Kallikrein-8



• Molecule 1: Kallikrein-8



• Molecule 1: Kallikrein-8



• Molecule 1: Kallikrein-8

Y228	L108	V16
T229	S113	L17
H230		Q23
I231	S116	
		P28
I238	K119	Q29
	P120	Q30
K245	I121	A31
G246		A32
	T130	L33
		F34
	Q134	
	S139	Q38
	G140	Q39
	W141	L40
	G142	L41
	T143	G42
	V144	G43
	T145	N50
		W51
	R148	V52
	E149	L53
	N150	T54
	F151	A55
		A56
	E159	H57
		C58
	T162	K62
	F163	V63
	P164	T64
	K167	V65
		R66
	A171	
	Q175	D70
		L73
	A183	Q74
		N75
	T190	K76
	C191	D77
	Q192	G78
	G193	P79
	L194	E80
	S195	O81
		E82
	C201	
	Q210	Q87
	G211	P95C
	T212	
	T213	H99
	S214	
	W215	L103
		M104
	D223A	L105
	K224	L106
	P225	G107

Chain E: 25% 75%

ACE1	
L2	
L3	
R4	

Chain F: 50% 50%

ACE1	
L2	
L3	
R4	

Chain G:

ACE1	
L2	
L3	
R4	

Chain H: 75% 25%

ACE1

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	83.10Å 46.05Å 103.35Å 90.00° 91.40° 90.00°	Depositor
Resolution (Å)	65.53 – 2.10 65.53 – 2.10	Depositor EDS
% Data completeness (in resolution range)	97.7 (65.53-2.10) 91.0 (65.53-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.18	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.63 (at 2.10Å)	Xtriage
Refinement program	PHENIX 1.11.1-2575	Depositor
R, R_{free}	0.212 , 0.259 0.215 , 0.251	Depositor DCC
R_{free} test set	2287 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	14.3	Xtriage
Anisotropy	0.682	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 47.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.38$, $\langle L^2 \rangle = 0.21$	Xtriage
Estimated twinning fraction	0.237 for h,-k,-l	Xtriage
Reported twinning fraction	0.720 for h,-k,-l	Depositor
Outliers	0 of 45153 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8120	wwPDB-VP
Average B, all atoms (Å ²)	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 46.19 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.1856e-04.*

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: CA, ACE, AR7, TBU

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	0.16	0/1752	0.39	0/2378
1	B	0.14	0/1752	0.38	0/2378
1	C	0.16	0/1752	0.37	0/2378
1	D	0.18	0/1752	0.39	0/2378
2	E	0.15	0/16	0.53	0/21
2	F	0.56	0/16	1.55	0/21
2	G	1.48	0/16	1.08	0/21
2	H	0.19	0/16	0.98	0/21
All	All	0.18	0/7072	0.40	0/9596

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	D	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	145	THR	Peptide

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	1711	0	1644	70	0
1	B	1711	0	1644	62	0
1	C	1711	0	1644	65	0
1	D	1711	0	1644	68	0
2	E	30	0	38	2	0
2	F	30	0	38	8	0
2	G	30	0	38	7	0
2	H	30	0	38	2	0
3	A	15	0	30	8	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	277	0	0	30	0
5	B	276	0	0	21	0
5	C	277	0	0	15	1
5	D	293	0	0	16	1
5	E	2	0	0	0	0
5	F	4	0	0	0	0
5	G	4	0	0	0	0
5	H	4	0	0	0	0
All	All	8120	0	6758	257	2

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 19.

All (257) 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:C:215:TRP:HB3	2:G:3:LEU:HD23	1.20	1.08
1:D:95(C):ASP:HB2	5:D:427:HOH:O	1.55	1.03
1:D:38:GLN:HB3	5:D:572:HOH:O	1.76	0.84
1:C:215:TRP:CB	2:G:3:LEU:HD23	2.07	0.84
1:D:145:THR:HG21	5:D:519:HOH:O	1.76	0.84
1:A:153:ASP:H	3:A:303:TBU:HO	1.29	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:192:GLN:HG2	5:A:432:HOH:O	1.81	0.80
1:A:173:PRO:HD3	5:A:428:HOH:O	1.84	0.75
1:B:218:ASP:HB2	2:F:2:LEU:HD22	1.69	0.74
1:A:246:GLY:HA2	5:A:512:HOH:O	1.87	0.74
1:D:66:ARG:NH2	1:D:70:ASP:OD2	2.17	0.74
1:A:237:TRP:HB2	3:A:302:TBU:H23	1.70	0.74
5:B:501:HOH:O	2:F:3:LEU:HD11	1.89	0.72
1:A:82:GLU:OE2	1:B:150:ASN:ND2	2.22	0.72
1:A:153:ASP:N	3:A:303:TBU:O	2.19	0.69
1:D:149:GLU:OE1	5:D:401:HOH:O	2.10	0.69
1:A:66:ARG:NH2	5:A:407:HOH:O	2.23	0.68
1:C:51:TRP:HE1	1:C:246:GLY:HA3	1.58	0.68
1:C:124:ALA:N	5:C:407:HOH:O	2.27	0.68
1:A:105:LEU:HD11	1:A:238:ILE:HG23	1.75	0.67
1:A:62:LYS:HD2	5:B:432:HOH:O	1.92	0.67
1:A:236:ASP:OD1	3:A:302:TBU:O	2.12	0.67
1:B:175:GLN:NE2	5:B:409:HOH:O	2.23	0.66
1:C:103:LEU:HG	1:C:229:THR:HG21	1.77	0.66
5:A:471:HOH:O	1:B:38:GLN:HB3	1.95	0.65
1:D:105:LEU:HD11	1:D:238:ILE:HA	1.78	0.65
1:B:103:LEU:HG	1:B:229:THR:HG21	1.78	0.65
1:D:175:GLN:HG2	5:D:510:HOH:O	1.96	0.65
1:B:134:GLN:NE2	1:B:202:ASP:OD2	2.31	0.64
1:A:114:LEU:HB3	5:A:447:HOH:O	1.97	0.63
1:C:215:TRP:HB3	2:G:3:LEU:CD2	2.12	0.63
1:C:134:GLN:NE2	5:C:415:HOH:O	2.31	0.63
1:B:87:GLN:HA	1:C:74:GLN:HA	1.81	0.63
1:D:212:ILE:HG12	1:D:231:ILE:HD11	1.81	0.63
1:A:234:TYR:HD1	3:A:302:TBU:H13	1.65	0.62
1:D:23:GLN:NE2	5:D:418:HOH:O	2.33	0.62
1:A:67:LEU:O	5:A:401:HOH:O	2.16	0.61
1:C:113:SER:HB2	5:C:539:HOH:O	1.99	0.61
1:A:217:SER:N	5:A:420:HOH:O	2.33	0.61
1:A:192:GLN:NE2	5:A:422:HOH:O	2.33	0.61
1:A:117:LYS:HB2	3:A:301:TBU:H23	1.82	0.61
1:D:52:VAL:HG21	1:D:65:VAL:HG11	1.83	0.61
1:A:133:GLY:O	1:A:161:LYS:NZ	2.34	0.60
1:B:122:SER:OG	5:B:401:HOH:O	2.17	0.60
1:B:73:LEU:O	5:B:402:HOH:O	2.17	0.60
1:B:219:PRO:HD2	1:B:222:ARG:HD2	1.84	0.59
1:D:78:GLY:N	1:D:80:GLU:OE1	2.30	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:85:VAL:HG22	1:C:87:GLN:H	1.67	0.59
1:C:47:VAL:HG11	1:C:53:LEU:HD12	1.85	0.58
1:A:164:PRO:HG2	1:A:167:LYS:HE2	1.85	0.58
1:D:213:THR:HG22	1:D:228:TYR:CE2	2.38	0.58
1:C:192:GLN:HG2	5:C:543:HOH:O	2.02	0.58
1:B:192:GLN:HE22	2:F:2:LEU:HG	1.68	0.58
1:A:134:GLN:N	5:A:404:HOH:O	2.37	0.58
1:B:81:GLN:NE2	1:B:113:SER:O	2.32	0.58
1:A:87:GLN:NE2	5:A:412:HOH:O	2.30	0.57
1:A:109:ARG:NH1	5:A:421:HOH:O	2.33	0.57
1:A:53:LEU:HD13	1:A:105:LEU:HG	1.87	0.57
1:B:63:TYR:OH	1:B:106:LEU:HD22	2.05	0.57
1:B:191:CYS:HB3	5:B:426:HOH:O	2.04	0.57
1:D:201:CYS:HB2	1:D:210:GLN:HG3	1.87	0.57
1:D:142:GLY:N	1:D:194:ASP:OD1	2.33	0.56
1:D:171:ALA:HB1	1:D:224:LYS:HG3	1.87	0.56
1:A:217:SER:HB2	1:A:224:LYS:HB2	1.88	0.56
1:D:95(C):ASP:N	5:D:427:HOH:O	2.39	0.55
1:B:237:TRP:NE1	5:B:420:HOH:O	2.34	0.55
1:B:246:GLY:HA3	1:C:38:GLN:N	2.21	0.55
1:A:83:ILE:HG21	1:A:108:LEU:HB3	1.87	0.55
1:B:233:ARG:NH1	5:B:423:HOH:O	2.38	0.55
1:D:77:ASP:HB3	5:D:478:HOH:O	2.05	0.55
1:A:192:GLN:NE2	5:A:431:HOH:O	2.40	0.55
1:A:66:ARG:NH1	5:A:413:HOH:O	2.30	0.54
1:B:143:THR:N	5:B:426:HOH:O	2.41	0.54
1:C:95(B):SER:OG	5:C:401:HOH:O	2.17	0.54
1:D:43:GLY:N	1:D:195:SER:O	2.41	0.54
1:D:63:TYR:HE2	1:D:106:LEU:HD13	1.71	0.54
1:A:191:CYS:SG	5:A:572:HOH:O	2.59	0.54
1:C:138:VAL:HB	5:C:497:HOH:O	2.06	0.53
1:D:213:THR:HG22	1:D:228:TYR:HE2	1.72	0.53
1:A:114:LEU:HD22	1:A:120:PRO:HD3	1.91	0.53
1:B:172:TYR:OH	1:B:225:PRO:O	2.24	0.53
1:D:143:THR:OG1	1:D:145:THR:O	2.24	0.53
1:C:145:THR:HG22	1:C:150:ASN:H	1.73	0.53
1:A:166:LYS:NZ	1:A:170:ASP:OD1	2.26	0.52
1:A:209:LEU:HD21	1:A:212:ILE:HD11	1.91	0.52
1:D:164:PRO:HG2	1:D:167:LYS:HG2	1.92	0.52
1:D:34:PHE:HZ	1:D:73:LEU:HD12	1.74	0.52
2:E:3:LEU:H	2:E:3:LEU:HD12	1.75	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:143:THR:HG21	1:C:149:GLU:HA	1.91	0.52
1:C:97:GLU:N	5:C:425:HOH:O	2.39	0.52
1:D:29:TRP:HB3	1:D:121:ILE:HB	1.90	0.52
1:B:207:GLY:O	5:B:401:HOH:O	2.19	0.52
1:B:214:SER:O	5:B:403:HOH:O	2.18	0.52
1:C:209:LEU:N	5:C:407:HOH:O	2.43	0.52
1:A:81:GLN:N	5:A:401:HOH:O	2.39	0.52
1:C:67:LEU:HD22	1:C:118:VAL:HG13	1.92	0.51
1:D:50:ASN:HA	1:D:108:LEU:HD12	1.92	0.51
1:D:130:THR:HG22	5:D:472:HOH:O	2.10	0.51
1:B:120:PRO:O	5:B:404:HOH:O	2.19	0.51
1:D:28:PRO:HB2	1:D:119:LYS:HB3	1.92	0.51
1:A:74:GLN:HE22	1:D:62:LYS:HA	1.75	0.51
1:A:44:GLY:O	5:A:402:HOH:O	2.19	0.51
1:B:72:SER:OG	1:B:153:ASP:OD2	2.19	0.51
1:C:99:HIS:CD2	2:G:3:LEU:HD11	2.45	0.51
1:B:63:TYR:CE2	1:B:106:LEU:HD13	2.46	0.50
5:A:414:HOH:O	1:D:87:GLN:NE2	2.43	0.50
1:B:235:LEU:HB2	5:B:519:HOH:O	2.11	0.50
1:A:178:ASP:OD2	1:A:233:ARG:NH2	2.43	0.50
1:A:115:GLY:HA3	3:A:301:TBU:H13	1.93	0.50
1:B:212:ILE:HG12	1:B:231:ILE:HD11	1.93	0.50
1:A:183:ALA:HB3	1:A:228:TYR:CE1	2.47	0.50
1:B:125:ASP:OD1	5:B:405:HOH:O	2.20	0.50
1:C:164:PRO:HB2	1:C:167:LYS:HG3	1.93	0.50
1:A:218:ASP:HB2	2:E:1:ACE:H1	1.93	0.50
1:A:236:ASP:O	1:A:240:LYS:HG2	2.11	0.49
1:C:211:GLY:HA2	1:C:229:THR:O	2.12	0.49
1:A:59:LYS:HE3	1:A:90:PRO:HG3	1.95	0.49
1:C:110:ASP:OD2	5:C:402:HOH:O	2.19	0.49
1:D:194:ASP:O	1:D:213:THR:OG1	2.28	0.49
1:B:130:THR:OG1	1:B:131:GLN:N	2.34	0.49
1:D:116:SER:OG	5:D:402:HOH:O	2.19	0.49
1:D:17:LEU:HD11	1:D:191:CYS:HB2	1.94	0.48
1:A:162:ILE:N	5:A:404:HOH:O	2.22	0.48
1:B:52:VAL:HG21	1:B:65:VAL:HG11	1.95	0.48
1:B:123:LEU:HD11	1:B:238:ILE:HG21	1.94	0.48
1:A:56:ALA:HA	1:A:104:MET:HB2	1.95	0.48
1:A:130:THR:OG1	1:A:131:GLN:N	2.47	0.48
1:D:159:GLU:OE1	5:D:403:HOH:O	2.19	0.48
1:A:201:CYS:HB2	1:A:210:GLN:HG3	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:28:PRO:HB2	1:C:119:LYS:HG2	1.95	0.48
1:C:99:HIS:CG	2:G:3:LEU:HD21	2.48	0.48
1:C:16:VAL:N	5:C:418:HOH:O	2.46	0.48
1:D:148:ARG:O	1:D:149:GLU:HB2	2.13	0.48
1:A:152:PRO:HB3	5:A:545:HOH:O	2.13	0.48
1:B:217:SER:HB2	1:B:224:LYS:HB2	1.95	0.48
1:A:38:GLN:CB	1:D:107:GLN:HE22	2.27	0.48
1:C:63:TYR:CE2	1:C:85:VAL:HG11	2.48	0.48
1:D:34:PHE:N	5:D:439:HOH:O	2.47	0.47
1:D:74:GLN:HE21	1:D:75:ASN:CG	2.22	0.47
1:B:74:GLN:NE2	5:B:432:HOH:O	2.47	0.47
1:D:66:ARG:HD2	1:D:82:GLU:HG2	1.96	0.47
1:D:81:GLN:NE2	1:D:113:SER:O	2.40	0.47
1:A:227:VAL:HG23	5:A:405:HOH:O	2.14	0.47
1:D:223(A):ASP:OD1	1:D:224:LYS:N	2.48	0.47
1:C:104:MET:HE3	1:C:106:LEU:HD21	1.97	0.47
5:B:430:HOH:O	2:F:1:ACE:CH3	2.62	0.47
1:B:110:ASP:OD1	1:C:39:GLN:NE2	2.47	0.47
1:C:85:VAL:HG21	1:C:106:LEU:HB3	1.97	0.47
1:D:63:TYR:HB2	5:D:439:HOH:O	2.14	0.47
1:B:191:CYS:HA	2:F:4:AR7:HD2	1.96	0.47
1:A:38:GLN:HE21	1:D:245:LYS:HE3	1.79	0.47
1:C:17:LEU:HD11	1:C:220:CYS:HB2	1.97	0.47
1:D:163:PHE:HD2	1:D:225:PRO:HG3	1.79	0.47
1:C:86:VAL:O	1:C:87:GLN:HB2	2.14	0.46
1:D:87:GLN:HB2	1:D:107:GLN:NE2	2.30	0.46
1:D:211:GLY:HA2	1:D:231:ILE:HG12	1.97	0.46
1:D:215:TRP:HA	2:H:4:AR7:HG2	1.97	0.46
1:C:16:VAL:N	1:C:194:ASP:OD2	2.49	0.46
1:D:32:ALA:HB2	1:D:141:TRP:HZ3	1.81	0.46
1:C:56:ALA:N	1:C:102:ASP:OD1	2.48	0.46
1:C:60:LYS:NZ	5:C:408:HOH:O	2.25	0.46
1:A:163:PHE:HB2	1:A:182:CYS:HB2	1.97	0.46
1:C:138:VAL:HG23	1:C:198:PRO:O	2.16	0.46
1:B:59:LYS:NZ	1:B:90:PRO:HD3	2.30	0.46
1:D:139:SER:O	5:D:404:HOH:O	2.21	0.46
1:C:31:ALA:HB1	1:C:65:VAL:HG13	1.98	0.46
1:D:149:GLU:OE2	5:D:405:HOH:O	2.21	0.45
1:C:45:VAL:HG11	1:C:209:LEU:HD22	1.98	0.45
1:C:85:VAL:O	5:C:403:HOH:O	2.21	0.45
1:D:134:GLN:NE2	1:D:201:CYS:HB3	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:240:LYS:HE2	5:A:498:HOH:O	2.15	0.45
1:C:53:LEU:HD22	1:C:103:LEU:HD22	1.98	0.45
1:C:63:TYR:CZ	1:C:85:VAL:HG11	2.51	0.45
1:D:143:THR:HA	1:D:151:PHE:HA	1.98	0.45
1:D:162:ILE:HD13	1:D:183:ALA:HB2	1.97	0.45
1:A:23:GLN:NE2	5:A:429:HOH:O	2.39	0.45
1:B:28:PRO:O	5:B:406:HOH:O	2.21	0.45
1:B:41:LEU:HD13	1:B:60:LYS:HG3	1.98	0.45
1:C:35:GLN:N	1:C:39:GLN:O	2.40	0.45
1:C:183:ALA:HB3	1:C:228:TYR:CE1	2.52	0.45
1:D:103:LEU:HD21	1:D:238:ILE:HD11	1.98	0.45
1:C:46:LEU:HD23	1:C:119:LYS:C	2.42	0.45
1:D:171:ALA:O	1:D:224:LYS:HE3	2.17	0.45
1:B:63:TYR:HE2	1:B:106:LEU:HD13	1.82	0.44
1:B:99:HIS:HD2	1:B:215:TRP:CE2	2.35	0.44
1:C:51:TRP:CZ3	1:C:107:GLN:HB2	2.52	0.44
1:B:234:TYR:O	1:B:238:ILE:HG12	2.17	0.44
1:C:85:VAL:HB	5:C:618:HOH:O	2.17	0.44
1:B:201:CYS:SG	1:B:210:GLN:HG3	2.56	0.44
1:C:85:VAL:HG13	5:C:442:HOH:O	2.17	0.44
1:D:183:ALA:HB3	1:D:228:TYR:CE1	2.52	0.44
1:B:164:PRO:HB2	5:B:494:HOH:O	2.18	0.44
1:C:177:THR:HG23	1:C:180:MET:HE3	1.99	0.44
1:B:149:GLU:C	1:B:150:ASN:HD22	2.25	0.44
1:A:105:LEU:HD12	1:A:241:ILE:HD11	2.00	0.44
1:B:192:GLN:NE2	2:F:3:LEU:O	2.50	0.44
1:C:33:LEU:O	1:C:40:LEU:HD12	2.18	0.44
1:D:39:GLN:OE1	1:D:40:LEU:N	2.50	0.44
1:D:74:GLN:HE21	1:D:75:ASN:ND2	2.16	0.44
1:C:216:GLY:HA3	2:G:4:AR7:CZ	2.48	0.44
1:B:50:ASN:HD22	1:C:38:GLN:NE2	2.16	0.44
1:C:60:LYS:HD3	1:C:60:LYS:HA	1.88	0.44
1:A:117:LYS:HB2	3:A:301:TBU:C2	2.48	0.44
1:B:59:LYS:HE3	1:B:59:LYS:HB2	1.41	0.44
1:C:87:GLN:HG2	1:C:89:ILE:HG12	2.00	0.44
1:B:56:ALA:HB2	1:B:103:LEU:O	2.18	0.43
1:B:105:LEU:HD12	1:B:241:ILE:HD12	1.99	0.43
1:A:30:GLN:HA	5:A:402:HOH:O	2.17	0.43
1:A:59:LYS:HE2	1:A:88:SER:HB3	2.00	0.43
1:A:62:LYS:NZ	5:A:417:HOH:O	2.32	0.43
1:C:215:TRP:CB	2:G:3:LEU:HA	2.49	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:240:LYS:HD3	1:A:240:LYS:HA	1.75	0.43
1:D:41:LEU:HD21	1:D:58:CYS:O	2.19	0.43
1:C:76:LYS:O	1:C:78:GLY:N	2.51	0.43
1:D:56:ALA:HA	1:D:104:MET:HB2	1.99	0.43
1:A:216:GLY:HA2	5:A:405:HOH:O	2.18	0.43
1:A:38:GLN:HB3	1:D:107:GLN:NE2	2.34	0.43
1:A:151:PHE:HZ	5:A:432:HOH:O	2.01	0.43
1:A:38:GLN:HB3	1:D:107:GLN:HE22	1.83	0.42
1:A:76:LYS:HE3	1:A:76:LYS:HB3	1.90	0.42
1:A:224:LYS:NZ	5:A:411:HOH:O	2.29	0.42
1:B:165:GLN:HA	1:B:168:CYS:HB3	2.01	0.42
1:B:171:ALA:HB1	1:B:224:LYS:HD3	2.01	0.42
1:B:195:SER:HA	1:B:213:THR:HB	2.01	0.42
1:C:56:ALA:HA	1:C:104:MET:HB2	2.01	0.42
1:B:218:ASP:OD2	2:F:2:LEU:HD13	2.19	0.42
1:D:54:THR:OG1	1:D:55:ALA:N	2.53	0.42
1:B:201:CYS:HB2	1:B:210:GLN:HG3	2.01	0.42
1:B:178:ASP:O	5:B:407:HOH:O	2.21	0.42
1:A:57:HIS:ND1	1:A:102:ASP:OD2	2.50	0.42
1:A:223(A):ASP:HB3	5:A:411:HOH:O	2.19	0.42
1:B:63:TYR:CE2	1:B:85:VAL:HG21	2.55	0.42
1:B:64:THR:C	5:B:433:HOH:O	2.62	0.42
1:B:215:TRP:HA	2:F:4:AR7:HG1	2.02	0.42
1:D:50:ASN:O	1:D:107:GLN:HA	2.19	0.41
1:B:50:ASN:OD1	1:B:50:ASN:N	2.52	0.41
1:C:69:GLY:HA3	1:C:79:PRO:O	2.20	0.41
1:C:145:THR:HG21	1:C:148:ARG:HB3	2.02	0.41
1:B:70:ASP:OD1	1:B:71:HIS:N	2.52	0.41
1:C:51:TRP:NE1	1:C:246:GLY:HA3	2.29	0.41
1:A:38:GLN:CG	1:D:107:GLN:HE22	2.34	0.41
1:A:172:TYR:OH	1:A:224:LYS:HB3	2.21	0.41
1:B:213:THR:HA	1:B:228:TYR:CD2	2.55	0.41
1:B:121:ILE:O	5:B:408:HOH:O	2.21	0.41
1:D:103:LEU:HD13	1:D:229:THR:HG21	2.03	0.41
1:C:160:VAL:HG21	1:C:183:ALA:HB1	2.03	0.41
1:A:103:LEU:HB2	1:A:229:THR:HG21	2.03	0.41
1:C:95:ASN:OD1	1:C:95(C):ASP:OD1	2.38	0.41
1:D:167:LYS:HG2	1:D:167:LYS:H	1.59	0.41
1:D:192:GLN:HB2	5:D:524:HOH:O	2.20	0.40
1:A:246:GLY:O	5:A:403:HOH:O	2.21	0.40
1:B:99:HIS:HD2	1:B:215:TRP:CD2	2.40	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:41:LEU:HD22	1:C:60:LYS:HE3	2.03	0.40
1:C:207:GLY:O	5:C:404:HOH:O	2.22	0.40
1:D:190:THR:HG23	2:H:4:AR7:HNE	1.86	0.40
1:A:103:LEU:HD13	1:A:229:THR:HG21	2.04	0.40
1:C:72:SER:OG	1:C:153:ASP:OD2	2.31	0.40

All (2) 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)
5:C:431:HOH:O	5:C:478:HOH:O[2_654]	2.10	0.10
5:D:575:HOH:O	5:D:611:HOH:O[1_545]	2.18	0.02

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	226/228 (99%)	219 (97%)	7 (3%)	0	100	100
1	B	226/228 (99%)	212 (94%)	14 (6%)	0	100	100
1	C	226/228 (99%)	213 (94%)	13 (6%)	0	100	100
1	D	226/228 (99%)	216 (96%)	10 (4%)	0	100	100
2	E	2/4 (50%)	1 (50%)	1 (50%)	0	100	100
2	F	2/4 (50%)	2 (100%)	0	0	100	100
2	G	2/4 (50%)	1 (50%)	1 (50%)	0	100	100
2	H	2/4 (50%)	2 (100%)	0	0	100	100
All	All	912/928 (98%)	866 (95%)	46 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	190/190 (100%)	187 (98%)	3 (2%)	55	64
1	B	190/190 (100%)	186 (98%)	4 (2%)	47	54
1	C	190/190 (100%)	184 (97%)	6 (3%)	34	38
1	D	190/190 (100%)	186 (98%)	4 (2%)	47	54
2	E	2/2 (100%)	2 (100%)	0	100	100
2	F	2/2 (100%)	1 (50%)	1 (50%)	0	0
2	G	2/2 (100%)	2 (100%)	0	100	100
2	H	2/2 (100%)	2 (100%)	0	100	100
All	All	768/768 (100%)	750 (98%)	18 (2%)	44	51

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	CYS
1	A	63	TYR
1	A	143	THR
1	B	45	VAL
1	B	59	LYS
1	B	96	VAL
1	B	148	ARG
1	C	67	LEU
1	C	96	VAL
1	C	143	THR
1	C	148	ARG
1	C	201	CYS
1	C	231	ILE
1	D	30	GLN
1	D	99	HIS
1	D	106	LEU
1	D	143	THR
2	F	2	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	38	GLN
1	A	74	GLN
1	A	150	ASN
1	A	156	ASN
1	A	192	GLN
1	B	23	GLN
1	B	74	GLN
1	B	95	ASN
1	B	99	HIS
1	B	150	ASN
1	B	192	GLN
1	B	210	GLN
1	C	35	GLN
1	C	38	GLN
1	C	99	HIS
1	D	30	GLN
1	D	74	GLN
1	D	87	GLN
1	D	107	GLN
1	D	230	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

4 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	AR7	G	4	1	10,10,11	0.22	0	9,11,13	1.29	1 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	AR7	H	4	2,1	10,10,11	0.26	0	9,11,13	1.59	1 (11%)
2	AR7	E	4	2,1	10,10,11	0.29	0	9,11,13	1.37	2 (22%)
2	AR7	F	4	2,1	10,10,11	0.24	0	9,11,13	1.10	1 (11%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	AR7	G	4	1	-	4/9/9/11	-
2	AR7	H	4	2,1	-	4/9/9/11	-
2	AR7	E	4	2,1	-	2/9/9/11	-
2	AR7	F	4	2,1	-	2/9/9/11	-

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	4	AR7	CG-CD-NE	-4.00	100.97	112.20
2	G	4	AR7	CG-CD-NE	-3.39	102.69	112.20
2	E	4	AR7	CG-CD-NE	-3.27	103.02	112.20
2	F	4	AR7	CG-CD-NE	-2.41	105.43	112.20
2	E	4	AR7	CB-CA-C	-2.02	109.49	112.25

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	E	4	AR7	O-C-CA-N
2	E	4	AR7	O-C-CA-CB
2	G	4	AR7	O-C-CA-N
2	H	4	AR7	NE-CD-CG-CB
2	G	4	AR7	NE-CD-CG-CB
2	H	4	AR7	CA-CB-CG-CD
2	G	4	AR7	O-C-CA-CB
2	H	4	AR7	O-C-CA-N
2	H	4	AR7	CG-CD-NE-CZ
2	F	4	AR7	NE-CD-CG-CB

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Mol	Chain	Res	Type	Atoms
2	F	4	AR7	O-C-CA-N
2	G	4	AR7	N-CA-CB-CG

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	G	4	AR7	1	0
2	H	4	AR7	2	0
2	F	4	AR7	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 4 are monoatomic - leaving 3 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	TBU	A	301	-	4,4,4	0.24	0	6,6,6	0.77	0
3	TBU	A	302	-	4,4,4	0.31	0	6,6,6	0.54	0
3	TBU	A	303	-	4,4,4	0.45	0	6,6,6	0.28	0

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.

3 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	301	TBU	3	0
3	A	302	TBU	3	0
3	A	303	TBU	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
2	G	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	G	3:LEU	C	4:AR7	N	1.19

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	228/228 (100%)	-1.25	0 100 100	5, 11, 24, 41	0
1	B	228/228 (100%)	-1.19	0 100 100	3, 13, 28, 41	0
1	C	228/228 (100%)	-1.21	0 100 100	5, 14, 27, 37	0
1	D	228/228 (100%)	-1.15	0 100 100	6, 14, 36, 60	0
2	E	2/4 (50%)	-1.23	0 100 100	17, 17, 17, 18	0
2	F	2/4 (50%)	-0.55	0 100 100	30, 30, 30, 35	0
2	G	2/4 (50%)	-1.13	0 100 100	12, 12, 12, 21	0
2	H	2/4 (50%)	-1.20	0 100 100	14, 14, 14, 24	0
All	All	920/928 (99%)	-1.20	0 100 100	3, 13, 29, 60	0

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains [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
2	AR7	G	4	11/12	0.98	0.04	9,11,21,43	0
2	AR7	H	4	11/12	0.98	0.04	8,13,24,25	0
2	AR7	E	4	11/12	0.99	0.03	6,10,15,26	0
2	AR7	F	4	11/12	0.99	0.03	7,11,19,24	0

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
3	TBU	A	302	5/5	0.98	0.04	10,10,18,20	0
4	CA	D	301	1/1	0.98	0.07	49,49,49,49	0
3	TBU	A	303	5/5	0.99	0.06	8,26,29,36	0
4	CA	B	301	1/1	0.99	0.02	29,29,29,29	0
4	CA	C	301	1/1	0.99	0.03	24,24,24,24	0
3	TBU	A	301	5/5	0.99	0.04	8,8,20,30	0
4	CA	A	304	1/1	1.00	0.01	15,15,15,15	0

6.5 Other polymers [i](#)

There are no such residues in this entry.