



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 24, 2026 – 03:52 AM UTC

PDB ID : 9CY7 / pdb\_00009cy7  
Title : Structure of S.pombe Dis3L2 in complex with oligoU RNA substrate  
Authors : Garg, A.; Joshua-Tor, L.  
Deposited on : 2024-08-01  
Resolution : 3.52 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](mailto: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>.

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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

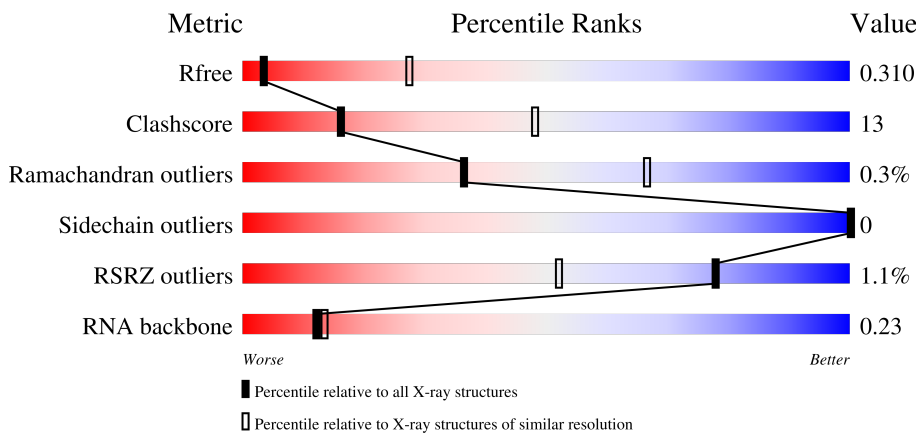
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.52 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	1025 (3.56-3.48)
Clashscore	190562	1079 (3.56-3.48)
Ramachandran outliers	187476	1052 (3.56-3.48)
Sidechain outliers	187428	1053 (3.56-3.48)
RSRZ outliers	180081	1024 (3.56-3.48)
RNA backbone	3983	1018 (4.00-3.04)

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  $\geq 3$ , 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	760	<div> <div> <div></div> <div>66%</div> <div>25%</div> <div>9%</div> </div> </div>
1	B	760	<div> <div> <div></div> <div>63%</div> <div>28%</div> <div>9%</div> </div> </div>
2	C	13	<div> <div> <div>15%</div> <div>31%</div> <div>46%</div> <div>8%</div> </div> </div>
2	D	13	<div> <div> <div>8%</div> <div>38%</div> <div>46%</div> <div>8%</div> </div> </div>

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 11555 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 DIS3-like exonuclease 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	693	Total	C	N	O	S	0	0	0
			5548	3515	960	1050	23			
1	B	693	Total	C	N	O	S	0	0	0
			5524	3492	956	1053	23			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	461	ASN	ASP	conflict	UNP O14040
B	461	ASN	ASP	conflict	UNP O14040

- Molecule 2 is a RNA chain called U13 RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	12	Total	C	N	O	P	0	0	0
			240	108	24	96	12			
2	D	12	Total	C	N	O	P	0	0	0
			240	108	24	96	12			

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		
3	B	1	Total	Mg	0	0
			1	1		

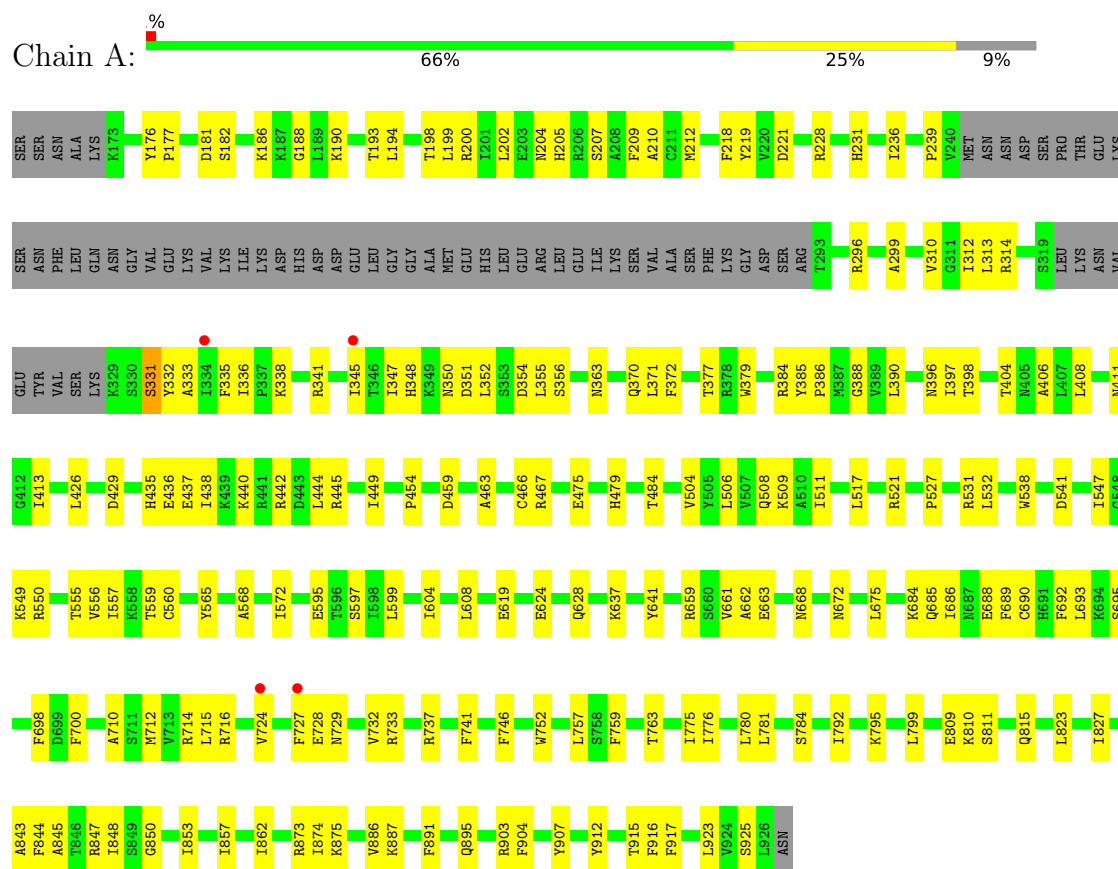
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	1	Total	O	0	0
			1	1		

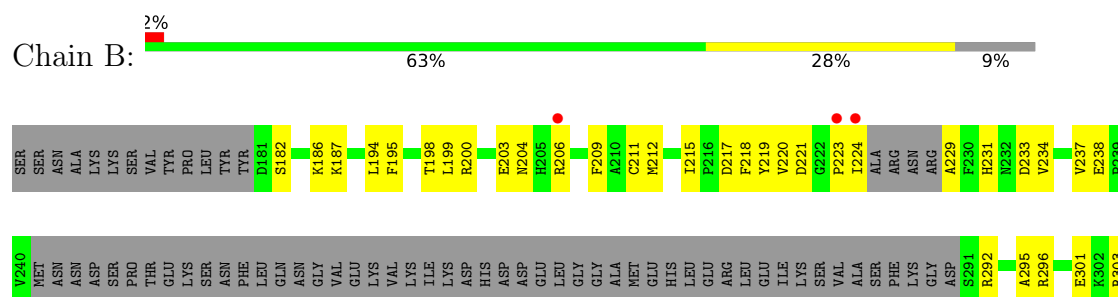
### 3 Residue-property plots

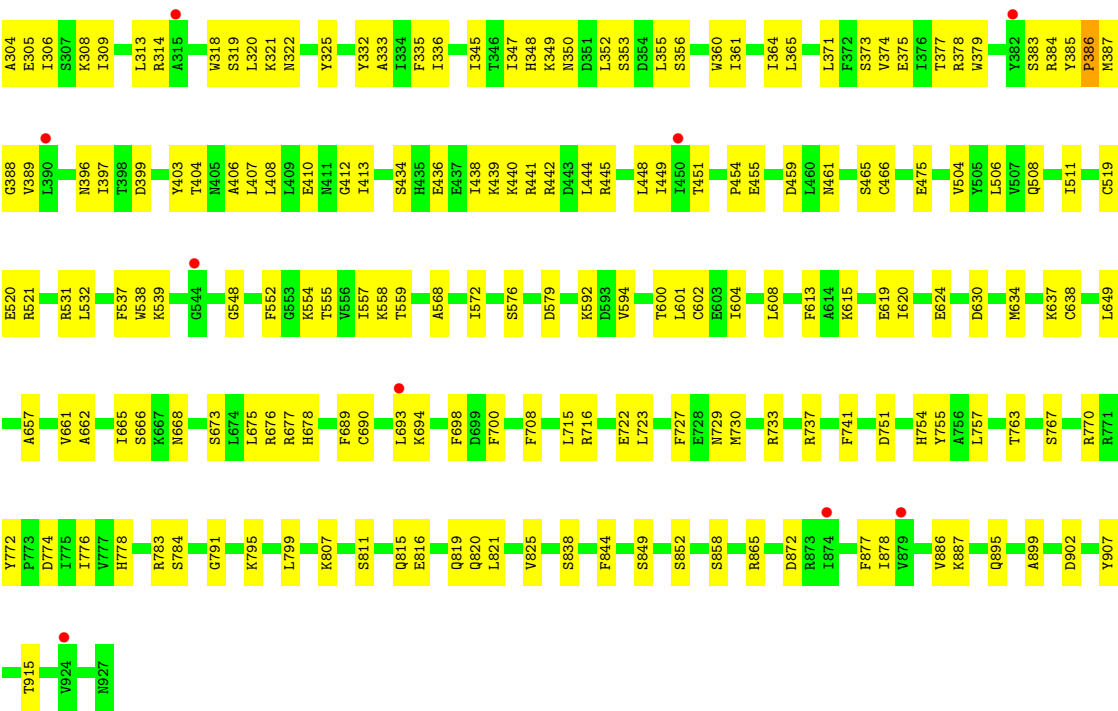
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: DIS3-like exonuclease 2



#### • Molecule 1: DIS3-like exonuclease 2





• Molecule 2: U13 RNA



• Molecule 2: U13 RNA



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	122.84Å 122.84Å 296.77Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.79 – 3.52 29.79 – 3.52	Depositor EDS
% Data completeness (in resolution range)	99.7 (29.79-3.52) 92.0 (29.79-3.52)	Depositor EDS
$R_{merge}$	0.15	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.59 (at 3.55Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487:0000)	Depositor
R, $R_{free}$	0.282 , 0.310 0.282 , 0.310	Depositor DCC
$R_{free}$ test set	2000 reflections (6.89%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	112.7	Xtriage
Anisotropy	0.501	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 74.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	11555	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	138.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.58% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

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	0.11	0/5665	0.35	0/7659
1	B	0.15	0/5636	0.36	0/7618
2	C	0.23	0/263	0.63	0/404
2	D	0.20	0/263	0.53	0/404
All	All	0.14	0/11827	0.37	0/16085

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5548	0	5504	131	0
1	B	5524	0	5470	160	0
2	C	240	0	121	13	0
2	D	240	0	121	13	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	D	1	0	0	0	0
All	All	11555	0	11216	298	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 13.

All (298) 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:212:MET:HB3	1:B:215:ILE:HG13	1.55	0.88
1:B:308:LYS:HD2	1:B:373:SER:HB2	1.60	0.83
1:B:442:ARG:HB3	1:B:555:THR:HG22	1.61	0.82
1:B:319:SER:H	1:B:322:ASN:HB2	1.48	0.78
1:A:442:ARG:HB3	1:A:555:THR:HG22	1.67	0.76
1:B:349:LYS:HG3	1:B:387:MET:HE1	1.69	0.75
1:B:539:LYS:HD2	1:B:548:GLY:HA3	1.70	0.74
1:A:188:GLY:HA3	1:A:194:LEU:HD23	1.68	0.73
1:B:754:HIS:NE2	2:D:10:U:OP1	2.23	0.72
1:B:677:ARG:NH1	1:B:751:ASP:O	2.26	0.68
1:A:413:ILE:HD13	1:A:504:VAL:HG21	1.76	0.68
1:B:620:ILE:HD11	1:B:708:PHE:HE1	1.57	0.68
1:B:733:ARG:HG2	2:D:7:U:H4'	1.75	0.67
1:A:628:GLN:HB2	1:A:637:LYS:HB3	1.77	0.66
1:B:441:ARG:HG2	1:B:554:LYS:HG3	1.78	0.66
1:A:444:LEU:HB2	1:A:557:ILE:HG22	1.77	0.66
1:B:849:SER:HG	1:B:852:SER:HG	1.42	0.66
1:A:194:LEU:HD12	1:A:236:ILE:HG22	1.77	0.66
1:B:220:VAL:HG13	1:B:223:PRO:HD2	1.78	0.66
1:B:375:GLU:O	1:B:388:GLY:HA3	1.96	0.65
1:A:310:VAL:HG11	1:A:406:ALA:HB1	1.79	0.65
1:B:200:ARG:HG3	1:B:231:HIS:HE1	1.62	0.64
1:B:237:VAL:HB	1:B:295:ALA:HB1	1.78	0.64
1:A:672:ASN:ND2	1:A:792:ILE:O	2.31	0.63
1:B:413:ILE:HD13	1:B:504:VAL:HG21	1.80	0.63
1:A:312:ILE:HG12	1:A:915:THR:HG21	1.81	0.63
1:B:361:ILE:O	1:B:365:LEU:HG	1.98	0.63
1:B:774:ASP:O	1:B:778:HIS:ND1	2.27	0.62
1:A:347:ILE:HG22	1:A:351:ASP:HB3	1.82	0.62
1:A:204:ASN:HB3	2:C:2:U:H1'	1.80	0.61
1:A:371:LEU:HG	1:A:397:ILE:HG12	1.82	0.61
1:B:466:CYS:HB2	1:B:601:LEU:HD21	1.82	0.61
1:B:335:PHE:HB3	1:B:345:ILE:HB	1.82	0.60
1:B:356:SER:HB3	1:B:360:TRP:HA	1.83	0.60
1:B:878:ILE:HG22	1:B:887:LYS:HB3	1.83	0.60
1:A:352:LEU:HD23	1:A:355:LEU:HB2	1.84	0.60
1:A:333:ALA:HB3	1:A:347:ILE:HB	1.84	0.59
1:A:355:LEU:HD22	1:A:363:ASN:HB3	1.83	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:737:ARG:HG2	2:D:8:U:H1'	1.84	0.59
1:A:210:ALA:HB3	1:A:218:PHE:HB2	1.85	0.59
1:A:352:LEU:HD13	1:A:390:LEU:HB3	1.83	0.58
1:A:541:ASP:HB3	1:A:547:ILE:HD11	1.85	0.58
1:A:438:ILE:HD11	1:A:445:ARG:HE	1.68	0.58
1:B:838:SER:HB2	1:B:907:TYR:HB3	1.84	0.58
1:B:321:LYS:HE2	1:B:321:LYS:HA	1.85	0.57
1:B:348:HIS:CE1	1:B:387:MET:HE3	2.39	0.57
1:B:613:PHE:CZ	1:B:619:GLU:HB2	2.39	0.57
1:B:238:GLU:O	1:B:295:ALA:HA	2.05	0.57
1:B:441:ARG:HD2	1:B:532:LEU:HD22	1.87	0.57
1:A:690:CYS:HA	1:A:693:LEU:HD12	1.87	0.56
1:B:449:ILE:HG12	1:B:465:SER:HB3	1.87	0.56
1:B:377:THR:O	1:B:386:PRO:HB2	2.06	0.56
1:B:195:PHE:HB3	1:B:212:MET:HE1	1.88	0.56
1:B:445:ARG:HG3	1:B:558:LYS:HB3	1.87	0.56
1:B:899:ALA:H	1:B:902:ASP:CG	2.13	0.56
1:A:729:ASN:O	1:A:733:ARG:HG3	2.07	0.55
1:B:336:ILE:HG21	1:B:915:THR:HG21	1.87	0.55
1:A:313:LEU:HD11	1:A:333:ALA:HB1	1.87	0.55
1:A:352:LEU:HB3	1:A:356:SER:HB2	1.88	0.55
1:A:181:ASP:OD1	1:A:182:SER:N	2.40	0.55
1:A:231:HIS:CD2	1:A:341:ARG:HD2	2.42	0.55
1:B:715:LEU:HD21	1:B:727:PHE:HD2	1.72	0.54
1:A:205:HIS:C	1:A:207:SER:H	2.15	0.54
1:B:352:LEU:HG	1:B:389:VAL:HA	1.89	0.54
1:A:467:ARG:HB3	1:A:475:GLU:HB3	1.89	0.54
1:A:710:ALA:O	1:A:714:ARG:HG3	2.07	0.54
1:A:795:LYS:O	1:A:799:LEU:HG	2.07	0.54
1:A:685:GLN:NE2	2:C:6:U:O2	2.40	0.54
1:B:449:ILE:HG21	1:B:557:ILE:HD12	1.90	0.54
2:C:8:U:H4'	2:C:9:U:OP2	2.06	0.54
1:A:396:ASN:ND2	1:A:398:THR:OG1	2.41	0.54
1:B:878:ILE:O	1:B:886:VAL:HA	2.08	0.54
1:B:360:TRP:O	1:B:364:ILE:HG13	2.08	0.53
1:A:746:PHE:HB2	1:A:752:TRP:HE1	1.72	0.53
1:B:887:LYS:HE3	1:B:895:GLN:HB3	1.89	0.53
1:A:449:ILE:O	1:A:560:CYS:N	2.40	0.53
1:B:314:ARG:HG2	1:B:318:TRP:CD2	2.44	0.53
1:A:219:TYR:CE2	1:A:221:ASP:HB2	2.43	0.53
1:A:875:LYS:HD2	1:A:891:PHE:HA	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:347:ILE:HB	1:B:352:LEU:CD2	2.39	0.53
1:B:195:PHE:HB2	1:B:237:VAL:HG23	1.90	0.53
1:B:521:ARG:O	1:B:531:ARG:NH2	2.42	0.53
1:A:350:ASN:O	1:A:356:SER:OG	2.26	0.52
2:C:10:U:H2'	2:C:11:U:H6	1.74	0.52
1:B:693:LEU:HD13	1:B:698:PHE:HD2	1.75	0.52
1:A:811:SER:O	1:A:815:GLN:HG3	2.10	0.52
1:A:354:ASP:O	1:A:355:LEU:HD23	2.10	0.52
2:C:6:U:H4'	2:C:7:U:OP2	2.10	0.52
1:A:338:LYS:HA	1:A:916:PHE:CE1	2.45	0.51
1:B:754:HIS:HE2	2:D:10:U:P	2.33	0.51
1:A:527:PRO:HA	1:A:559:THR:OG1	2.11	0.51
1:B:198:THR:HB	1:B:211:CYS:HB2	1.93	0.51
1:B:630:ASP:OD1	1:B:634:MET:N	2.44	0.51
1:A:659:ARG:HD2	1:A:759:PHE:CE1	2.46	0.51
1:B:408:LEU:HD13	1:B:511:ILE:HD11	1.92	0.51
1:A:689:PHE:CE2	1:A:693:LEU:HD11	2.46	0.50
1:B:347:ILE:HB	1:B:352:LEU:HD22	1.92	0.50
1:A:202:LEU:HD11	1:A:209:PHE:HE1	1.76	0.50
1:A:372:PHE:HB3	1:A:390:LEU:HD21	1.94	0.50
1:A:200:ARG:HG3	1:A:231:HIS:NE2	2.26	0.50
1:A:436:GLU:O	1:A:440:LYS:HG2	2.11	0.50
1:A:538:TRP:HD1	1:A:550:ARG:HG2	1.77	0.50
1:A:404:THR:O	1:A:408:LEU:HG	2.12	0.50
1:B:233:ASP:OD1	1:B:303:ARG:HA	2.12	0.50
1:A:847:ARG:NH1	2:C:5:U:OP1	2.39	0.50
1:B:229:ALA:N	1:B:384:ARG:HE	2.10	0.50
1:B:360:TRP:CE2	1:B:361:ILE:HG13	2.47	0.50
1:B:209:PHE:CZ	2:D:4:U:H2'	2.47	0.50
1:B:412:GLY:O	1:B:807:LYS:NZ	2.45	0.49
1:B:434:SER:O	1:B:438:ILE:HG13	2.12	0.49
1:A:714:ARG:HG2	1:B:615:LYS:HG2	1.93	0.49
1:B:187:LYS:HD3	1:B:187:LYS:C	2.37	0.49
1:A:438:ILE:HD11	1:A:445:ARG:NE	2.26	0.49
1:A:408:LEU:HD13	1:A:511:ILE:HD11	1.93	0.49
1:A:209:PHE:HZ	2:C:4:U:H1'	1.78	0.49
1:A:312:ILE:HG12	1:A:915:THR:CG2	2.42	0.49
1:A:312:ILE:HD11	1:A:336:ILE:HD12	1.94	0.49
1:B:204:ASN:ND2	1:B:206:ARG:HB2	2.27	0.49
1:B:877:PHE:HA	1:B:887:LYS:O	2.13	0.49
1:B:455:GLU:CD	1:B:455:GLU:H	2.19	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:448:LEU:HD11	1:B:594:VAL:HG22	1.95	0.49
1:A:521:ARG:O	1:A:531:ARG:NH2	2.45	0.48
1:B:403:TYR:O	1:B:407:LEU:HG	2.12	0.48
1:B:454:PRO:HD3	2:D:12:U:O2'	2.13	0.48
1:B:620:ILE:HD11	1:B:708:PHE:CE1	2.44	0.48
1:A:186:LYS:O	1:A:190:LYS:HG2	2.14	0.48
1:A:568:ALA:O	1:A:572:ILE:HG13	2.13	0.48
1:B:198:THR:HG22	1:B:199:LEU:H	1.79	0.48
1:B:206:ARG:HH22	2:D:2:U:P	2.36	0.48
1:B:436:GLU:O	1:B:440:LYS:HG3	2.12	0.48
1:B:200:ARG:HD3	2:D:6:U:C2	2.49	0.48
1:A:408:LEU:HD21	1:A:506:LEU:HD11	1.95	0.48
1:B:313:LEU:HD11	1:B:333:ALA:HB1	1.95	0.48
1:A:236:ILE:HB	1:A:299:ALA:HB3	1.96	0.48
1:A:335:PHE:HB2	1:A:347:ILE:HD11	1.96	0.48
1:A:729:ASN:O	1:A:732:VAL:HG12	2.14	0.48
1:B:408:LEU:HD21	1:B:506:LEU:HD11	1.96	0.48
1:B:406:ALA:O	1:B:410:GLU:HG3	2.15	0.47
1:B:795:LYS:O	1:B:799:LEU:HG	2.14	0.47
2:D:8:U:H4'	2:D:9:U:OP1	2.14	0.47
1:A:435:HIS:O	1:A:438:ILE:HG22	2.14	0.47
1:A:737:ARG:NE	2:C:8:U:O2	2.47	0.47
1:B:319:SER:N	1:B:322:ASN:HB2	2.23	0.47
1:A:231:HIS:HD2	1:A:341:ARG:HD2	1.77	0.47
1:A:332:TYR:CD2	1:A:348:HIS:HA	2.50	0.47
1:A:538:TRP:CD1	1:A:550:ARG:HG2	2.50	0.47
1:B:303:ARG:HH21	1:B:305:GLU:HA	1.80	0.47
1:B:729:ASN:O	1:B:733:ARG:HG3	2.15	0.47
1:B:441:ARG:NH1	1:B:554:LYS:HB2	2.29	0.47
1:B:371:LEU:HG	1:B:397:ILE:HG12	1.95	0.47
1:B:602:CYS:HA	1:B:649:LEU:HD12	1.95	0.47
1:B:811:SER:O	1:B:815:GLN:HG3	2.15	0.47
1:B:872:ASP:N	1:B:872:ASP:OD1	2.48	0.47
1:A:411:ASN:HB3	1:A:810:LYS:HB3	1.96	0.47
1:B:693:LEU:HD13	1:B:698:PHE:CD2	2.50	0.47
1:A:887:LYS:HE2	1:A:895:GLN:OE1	2.15	0.46
1:A:907:TYR:CZ	1:A:925:SER:HB2	2.50	0.46
1:B:690:CYS:O	1:B:694:LYS:HG2	2.14	0.46
1:B:436:GLU:HA	1:B:439:LYS:HD2	1.97	0.46
1:A:377:THR:HG23	1:A:388:GLY:HA2	1.97	0.46
1:B:451:THR:HG23	1:B:559:THR:HG21	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:10:U:H2'	2:C:11:U:C6	2.50	0.46
1:B:678:HIS:ND1	1:B:755:TYR:HB3	2.31	0.46
1:B:592:LYS:H	1:B:592:LYS:HD3	1.80	0.46
1:B:865:ARG:NH1	2:D:4:U:OP1	2.44	0.46
1:B:352:LEU:HD21	1:B:388:GLY:O	2.14	0.46
1:A:823:LEU:O	1:A:827:ILE:HG13	2.16	0.46
1:A:313:LEU:O	1:A:314:ARG:NH1	2.48	0.45
1:A:853:ILE:HD11	1:A:886:VAL:HG11	1.99	0.45
1:A:331:SER:OG	1:A:332:TYR:N	2.48	0.45
1:A:692:PHE:O	1:A:695:SER:OG	2.32	0.45
1:A:757:LEU:HB3	1:A:759:PHE:CD2	2.51	0.45
1:B:461:ASN:HD22	1:B:767:SER:HA	1.82	0.45
1:B:475:GLU:HG2	1:B:537:PHE:HD2	1.81	0.45
1:B:237:VAL:HA	1:B:296:ARG:O	2.15	0.45
1:B:552:PHE:HE2	1:B:661:VAL:HG22	1.81	0.45
1:B:459:ASP:OD1	1:B:459:ASP:N	2.39	0.45
1:A:843:ALA:HB2	1:A:857:ILE:HG13	1.99	0.45
1:B:309:ILE:HB	1:B:374:VAL:HG12	1.97	0.45
1:B:613:PHE:CE1	1:B:619:GLU:HB2	2.52	0.45
1:A:200:ARG:HH12	2:C:6:U:H5'	1.82	0.44
1:A:219:TYR:HE2	1:A:221:ASP:HB2	1.81	0.44
1:A:684:LYS:O	1:A:688:GLU:HG3	2.17	0.44
1:A:662:ALA:HB2	1:A:763:THR:HG22	1.98	0.44
1:B:234:VAL:O	1:B:301:GLU:N	2.49	0.44
1:B:519:CYS:SG	1:B:520:GLU:HG3	2.58	0.44
1:B:349:LYS:O	1:B:352:LEU:HB3	2.16	0.44
1:B:619:GLU:O	1:B:716:ARG:NH2	2.36	0.44
1:A:459:ASP:OD1	1:A:459:ASP:N	2.48	0.44
1:B:444:LEU:HB2	1:B:557:ILE:HG22	2.00	0.44
1:B:730:MET:HA	1:B:733:ARG:HD2	2.00	0.44
1:B:770:ARG:NH2	2:D:12:U:OP2	2.51	0.44
2:D:6:U:H4'	2:D:7:U:OP2	2.15	0.44
1:A:659:ARG:O	1:A:663:GLU:HG3	2.17	0.44
1:A:844:PHE:CZ	1:A:903:ARG:HB2	2.53	0.44
1:B:844:PHE:HA	1:B:902:ASP:O	2.18	0.44
1:A:182:SER:O	1:A:186:LYS:HG3	2.18	0.44
1:B:200:ARG:NH2	1:B:209:PHE:HD2	2.16	0.44
1:B:322:ASN:HB3	1:B:325:TYR:HB3	2.00	0.44
1:A:624:GLU:HG3	1:A:641:TYR:HB3	2.00	0.43
1:A:426:LEU:HD13	1:A:517:LEU:HD23	1.99	0.43
1:A:437:GLU:HG2	1:A:532:LEU:HD21	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:712:MET:O	1:A:716:ARG:HG3	2.18	0.43
1:B:662:ALA:HB2	1:B:763:THR:HG22	1.99	0.43
1:B:665:ILE:O	1:B:673:SER:OG	2.37	0.43
1:A:193:THR:O	1:A:239:PRO:HD3	2.19	0.43
1:B:314:ARG:HG2	1:B:318:TRP:CE2	2.53	0.43
1:B:378:ARG:HG2	1:B:379:TRP:N	2.34	0.43
1:B:308:LYS:HD3	1:B:309:ILE:H	1.83	0.43
1:B:693:LEU:HD12	1:B:700:PHE:HB2	2.00	0.43
1:A:198:THR:HG22	1:A:199:LEU:N	2.32	0.43
1:A:438:ILE:HG13	1:A:556:VAL:HG11	2.01	0.43
1:B:219:TYR:CG	1:B:220:VAL:N	2.87	0.43
1:B:816:GLU:O	1:B:820:GLN:HG2	2.19	0.43
1:A:348:HIS:CG	1:A:351:ASP:HB2	2.53	0.43
1:B:234:VAL:HG12	1:B:301:GLU:HB3	2.01	0.43
1:B:182:SER:O	1:B:186:LYS:HG3	2.19	0.43
1:B:384:ARG:HA	1:B:384:ARG:HH11	1.83	0.43
1:A:508:GLN:HE21	1:A:509:LYS:HE2	1.84	0.43
1:A:693:LEU:HD13	1:A:700:PHE:CD1	2.54	0.43
1:B:783:ARG:CZ	1:B:791:GLY:HA3	2.49	0.43
1:A:384:ARG:HD2	1:A:385:TYR:CZ	2.54	0.43
1:A:668:ASN:OD1	1:A:784:SER:HB2	2.18	0.42
1:A:912:TYR:HE1	1:A:917:PHE:CD2	2.37	0.42
1:B:404:THR:HG21	1:B:508:GLN:HE22	1.84	0.42
1:A:715:LEU:HD21	1:A:727:PHE:HD2	1.84	0.42
1:B:350:ASN:ND2	1:B:360:TRP:CD2	2.86	0.42
1:B:722:GLU:HB3	1:B:858:SER:CB	2.49	0.42
1:B:772:TYR:CZ	1:B:776:ILE:HD11	2.54	0.42
2:C:3:U:O2'	2:C:4:U:H5	2.01	0.42
1:A:737:ARG:NH2	1:A:809:GLU:HG3	2.34	0.42
1:A:176:TYR:HB2	1:A:228:ARG:HH11	1.84	0.42
1:B:349:LYS:HB2	1:B:352:LEU:HB2	2.01	0.42
1:B:371:LEU:HD21	1:B:397:ILE:HG23	2.02	0.42
1:A:661:VAL:HG13	1:A:781:LEU:HD22	2.01	0.42
1:A:675:LEU:HB2	1:A:741:PHE:CZ	2.54	0.42
1:B:637:LYS:HD3	1:B:638:CYS:H	1.84	0.42
1:A:348:HIS:CD2	1:A:351:ASP:HB2	2.54	0.42
1:A:757:LEU:HD11	2:C:10:U:H4'	2.01	0.42
1:B:676:ARG:NH1	2:D:10:U:OP2	2.53	0.42
1:B:220:VAL:HG13	1:B:223:PRO:CD	2.47	0.42
1:B:303:ARG:NE	1:B:304:ALA:O	2.52	0.42
1:B:396:ASN:HB2	1:B:399:ASP:HB2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:723:LEU:HD23	1:B:723:LEU:HA	1.88	0.42
1:B:754:HIS:CE1	1:B:757:LEU:HG	2.55	0.42
1:A:686:ILE:HD12	1:A:686:ILE:HA	1.90	0.42
1:B:238:GLU:CD	1:B:296:ARG:HH21	2.28	0.42
1:B:689:PHE:HE1	1:B:730:MET:HB2	1.85	0.42
2:C:9:U:H2'	2:C:10:U:H6	1.85	0.42
1:A:463:ALA:HB3	1:A:479:HIS:HB2	2.02	0.42
1:A:595:GLU:O	1:A:599:LEU:HG	2.19	0.42
1:A:693:LEU:HB3	1:A:698:PHE:HB3	2.01	0.42
1:A:845:ALA:HB2	1:A:904:PHE:HE1	1.85	0.42
1:A:212:MET:HG2	1:A:218:PHE:HE1	1.85	0.42
1:A:716:ARG:HA	1:A:724:VAL:HG21	2.02	0.42
1:A:873:ARG:H	1:A:873:ARG:HG3	1.64	0.42
1:B:203:GLU:HG2	1:B:204:ASN:N	2.35	0.42
1:B:212:MET:HE3	1:B:218:PHE:HD2	1.83	0.42
1:B:689:PHE:CE2	1:B:693:LEU:HD11	2.55	0.41
1:A:335:PHE:HB3	1:A:345:ILE:HB	2.03	0.41
1:A:874:ILE:HD11	1:A:923:LEU:HD23	2.02	0.41
1:B:224:ILE:HB	1:B:384:ARG:HH21	1.85	0.41
1:B:353:SER:C	1:B:355:LEU:H	2.27	0.41
1:B:194:LEU:HD13	1:B:238:GLU:HG3	2.02	0.41
1:A:379:TRP:CG	1:A:386:PRO:HB3	2.56	0.41
1:A:466:CYS:SG	1:A:597:SER:OG	2.73	0.41
1:B:217:ASP:O	1:B:292:ARG:HD3	2.21	0.41
1:B:568:ALA:O	1:B:572:ILE:HG13	2.20	0.41
1:B:604:ILE:O	1:B:608:LEU:HG	2.20	0.41
1:A:454:PRO:HD3	1:A:565:TYR:HE2	1.85	0.41
1:A:776:ILE:O	1:A:780:LEU:HG	2.20	0.41
1:B:200:ARG:HH21	1:B:209:PHE:HD2	1.67	0.41
1:B:538:TRP:HZ2	1:B:657:ALA:HA	1.86	0.41
1:A:351:ASP:O	1:A:352:LEU:HB2	2.19	0.41
1:A:712:MET:HE1	1:A:728:GLU:HG2	2.02	0.41
1:A:370:GLN:HB2	1:A:372:PHE:CZ	2.56	0.41
1:B:383:SER:HB3	1:B:385:TYR:CE1	2.56	0.41
1:B:404:THR:O	1:B:408:LEU:HG	2.19	0.41
1:B:666:SER:HB3	1:B:675:LEU:HD11	2.03	0.41
1:A:429:ASP:OD1	1:A:429:ASP:N	2.54	0.41
1:A:619:GLU:O	1:A:716:ARG:NH2	2.54	0.41
1:B:576:SER:HB2	1:B:579:ASP:OD2	2.20	0.41
1:B:668:ASN:ND2	1:B:784:SER:HB2	2.36	0.41
1:A:177:PRO:HD3	1:A:296:ARG:HH21	1.86	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:823:LEU:HD21	1:A:862:ILE:HD13	2.03	0.40
1:A:848:ILE:HD13	1:A:886:VAL:HG12	2.02	0.40
1:B:308:LYS:NZ	1:B:375:GLU:OE2	2.53	0.40
1:B:332:TYR:CG	1:B:348:HIS:HB3	2.55	0.40
1:B:624:GLU:HG3	1:B:819:GLN:HE22	1.85	0.40
1:A:549:LYS:HG2	1:A:550:ARG:O	2.22	0.40
1:A:604:ILE:O	1:A:608:LEU:HG	2.21	0.40
1:B:313:LEU:O	1:B:314:ARG:NH1	2.49	0.40
1:B:600:THR:O	1:B:604:ILE:HG13	2.22	0.40
1:B:821:LEU:O	1:B:825:VAL:HG23	2.22	0.40
1:A:484:THR:HG22	1:A:775:ILE:HG23	2.03	0.40
1:B:219:TYR:CE2	1:B:221:ASP:HB2	2.56	0.40
1:B:319:SER:OG	1:B:320:LEU:N	2.55	0.40
1:B:675:LEU:HB2	1:B:741:PHE:CE1	2.56	0.40

There are no symmetry-related clashes.

## 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	687/760 (90%)	655 (95%)	30 (4%)	2 (0%)	36	66
1	B	687/760 (90%)	649 (94%)	36 (5%)	2 (0%)	36	66
All	All	1374/1520 (90%)	1304 (95%)	66 (5%)	4 (0%)	36	66

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	850	GLY
1	B	386	PRO
1	A	331	SER
1	B	306	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	619/679 (91%)	619 (100%)	0	100	100
1	B	617/679 (91%)	617 (100%)	0	100	100
All	All	1236/1358 (91%)	1236 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	628	GLN
1	A	648	HIS
1	A	761	HIS
1	B	363	ASN
1	B	753	HIS

### 5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	C	11/13 (84%)	5 (45%)	2 (18%)
2	D	11/13 (84%)	8 (72%)	2 (18%)
All	All	22/26 (84%)	13 (59%)	4 (18%)

All (13) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	C	3	U
2	C	4	U
2	C	7	U
2	C	8	U
2	C	9	U
2	D	3	U
2	D	4	U

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Mol	Chain	Res	Type
2	D	5	U
2	D	6	U
2	D	7	U
2	D	9	U
2	D	10	U
2	D	11	U

All (4) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	C	6	U
2	C	8	U
2	D	6	U
2	D	8	U

## 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 2 ligands modelled in this entry, 2 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 [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	693/760 (91%)	0.03	4 (0%) 85 64	80, 135, 193, 230	0
1	B	693/760 (91%)	0.09	12 (1%) 69 41	79, 133, 194, 221	0
2	C	12/13 (92%)	0.47	0 100 100	86, 161, 217, 227	0
2	D	12/13 (92%)	0.53	0 100 100	104, 143, 202, 205	0
All	All	1410/1546 (91%)	0.07	16 (1%) 78 52	79, 134, 196, 230	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	334	ILE	2.8
1	B	315	ALA	2.8
1	A	727	PHE	2.8
1	B	693	LEU	2.7
1	B	382	TYR	2.7
1	B	223	PRO	2.6
1	B	390	LEU	2.5
1	B	206	ARG	2.5
1	A	724	VAL	2.4
1	B	874	ILE	2.3
1	B	924	VAL	2.2
1	B	450	ILE	2.2
1	B	224	ILE	2.1
1	B	544	GLY	2.1
1	A	345	ILE	2.0
1	B	879	VAL	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, 95<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
3	MG	A	1001	1/1	0.92	0.08	116,116,116,116	0
3	MG	B	1001	1/1	0.97	0.04	166,166,166,166	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.