



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:36 PM GMT

PDB ID : 1L0Q
Title : Tandem YVTN beta-propeller and PKD domains from an archaeal surface layer protein
Authors : Jing, H.; Takagi, J.; Liu, J.-H.; Lindgren, S.; Zhang, R.-G.; Joachimiak, A.; Wang, J.-H.; Springer, T.A.
Deposited on : 2002-02-12
Resolution : 2.40 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

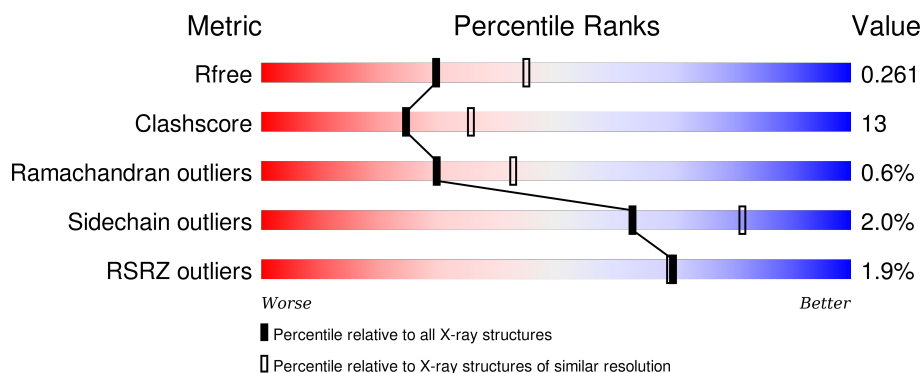
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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}	91344	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	391	<div> <div>%</div> <div> <div></div> <div>75%</div> <div>23%</div> <div>.</div> </div> </div>
1	B	391	<div> <div>%</div> <div> <div></div> <div>75%</div> <div>24%</div> <div>.</div> </div> </div>
1	C	391	<div> <div>5%</div> <div> <div></div> <div>76%</div> <div>22%</div> <div>.</div> </div> </div>
1	D	391	<div> <div>%</div> <div> <div></div> <div>76%</div> <div>23%</div> <div>.</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 12764 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 Surface layer protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	391	Total	C	N	O	S	Se	0	0	0
			2850	1782	466	596	1	5			
1	B	391	Total	C	N	O	S	Se	0	0	0
			2850	1782	466	596	1	5			
1	C	391	Total	C	N	O	S	Se	0	0	0
			2850	1782	466	596	1	5			
1	D	391	Total	C	N	O	S	Se	0	0	0
			2850	1782	466	596	1	5			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	35	MSE	MET	MODIFIED RESIDUE	UNP Q50245
A	93	MSE	MET	MODIFIED RESIDUE	UNP Q50245
A	180	MSE	MET	MODIFIED RESIDUE	UNP Q50245
A	228	MSE	MET	MODIFIED RESIDUE	UNP Q50245
A	265	YCM	CYS	MODIFIED RESIDUE	UNP Q50245
A	282	MSE	MET	MODIFIED RESIDUE	UNP Q50245
B	35	MSE	MET	MODIFIED RESIDUE	UNP Q50245
B	93	MSE	MET	MODIFIED RESIDUE	UNP Q50245
B	180	MSE	MET	MODIFIED RESIDUE	UNP Q50245
B	228	MSE	MET	MODIFIED RESIDUE	UNP Q50245
B	265	YCM	CYS	MODIFIED RESIDUE	UNP Q50245
B	282	MSE	MET	MODIFIED RESIDUE	UNP Q50245
C	35	MSE	MET	MODIFIED RESIDUE	UNP Q50245
C	93	MSE	MET	MODIFIED RESIDUE	UNP Q50245
C	180	MSE	MET	MODIFIED RESIDUE	UNP Q50245
C	228	MSE	MET	MODIFIED RESIDUE	UNP Q50245
C	265	YCM	CYS	MODIFIED RESIDUE	UNP Q50245
C	282	MSE	MET	MODIFIED RESIDUE	UNP Q50245
D	35	MSE	MET	MODIFIED RESIDUE	UNP Q50245
D	93	MSE	MET	MODIFIED RESIDUE	UNP Q50245
D	180	MSE	MET	MODIFIED RESIDUE	UNP Q50245

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Chain	Residue	Modelled	Actual	Comment	Reference
D	228	MSE	MET	MODIFIED RESIDUE	UNP Q50245
D	265	YCM	CYS	MODIFIED RESIDUE	UNP Q50245
D	282	MSE	MET	MODIFIED RESIDUE	UNP Q50245

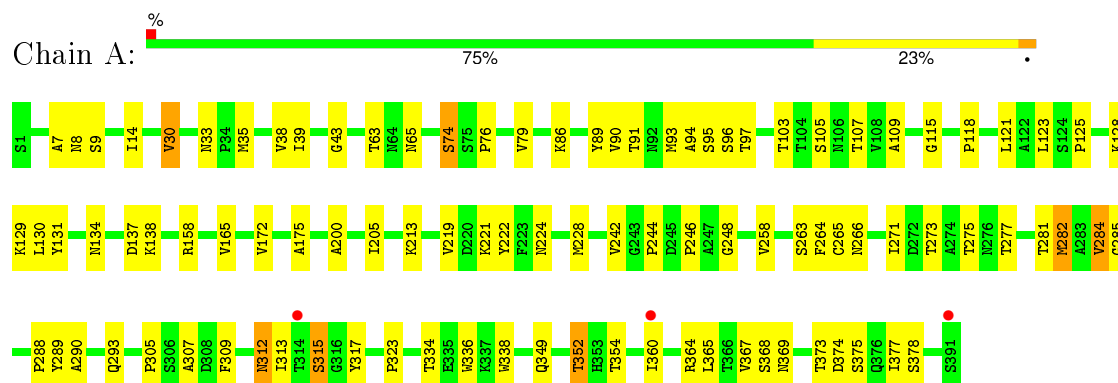
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	334	Total 334	O 334	0	0
2	B	348	Total 348	O 348	0	0
2	C	310	Total 310	O 310	0	0
2	D	372	Total 372	O 372	0	0

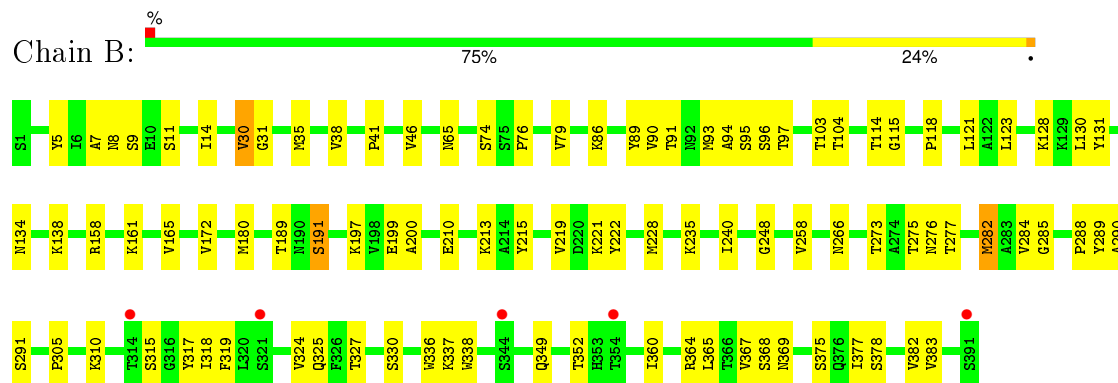
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

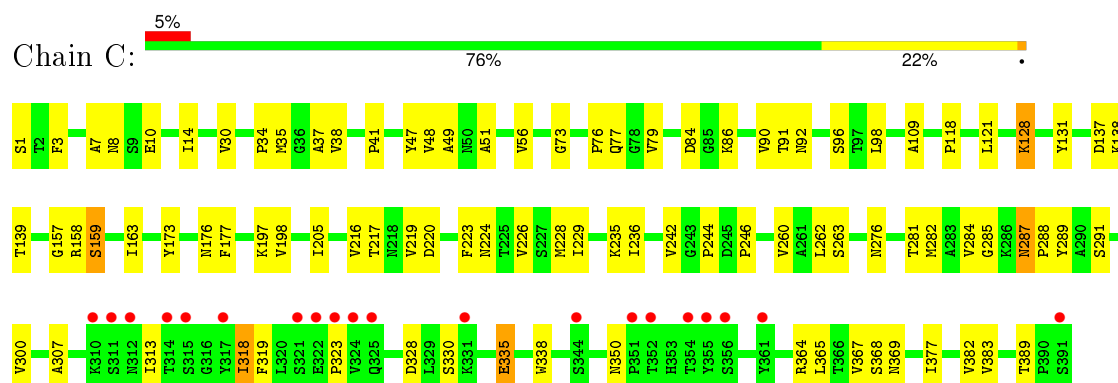
• Molecule 1: Surface layer protein



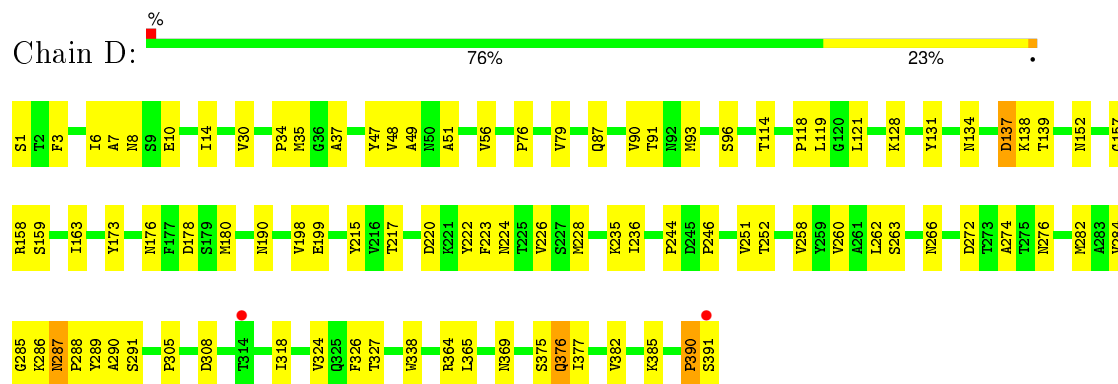
• Molecule 1: Surface layer protein



• Molecule 1: Surface layer protein



● Molecule 1: Surface layer protein



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	188.18Å 65.89Å 140.03Å 90.00° 124.25° 90.00°	Depositor
Resolution (Å)	20.00 – 2.40 41.42 – 2.40	Depositor EDS
% Data completeness (in resolution range)	93.9 (20.00-2.40) 93.9 (41.42-2.40)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.49 (at 2.39Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.214 , 0.263 0.214 , 0.261	Depositor DCC
R_{free} test set	2642 reflections (5.05%)	DCC
Wilson B-factor (Å ²)	19.5	Xtriage
Anisotropy	0.159	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 53.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	21 of 52426 reflections (0.040%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12764	wwPDB-VP
Average B, all atoms (Å ²)	22.0	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 63.86 % 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 9.0635e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: YCM

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	0.40	0/2885	0.70	0/3946
1	B	0.40	0/2885	0.69	0/3946
1	C	0.41	0/2885	0.68	0/3946
1	D	0.41	0/2885	0.68	0/3946
All	All	0.41	0/11540	0.69	0/15784

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	2850	0	2848	75	0
1	B	2850	0	2848	77	0
1	C	2850	0	2848	67	0
1	D	2850	0	2848	80	0
2	A	334	0	0	12	0
2	B	348	0	0	18	0
2	C	310	0	0	11	0
2	D	372	0	0	16	0
All	All	12764	0	11392	295	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 (295) 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:D:138:LYS:HE2	1:D:158:ARG:HG2	1.16	1.10
1:D:138:LYS:CE	1:D:158:ARG:HG2	1.96	0.95
1:B:318:ILE:HD13	1:B:324:VAL:HG22	1.47	0.93
1:B:213:LYS:HD3	1:B:228:MSE:HE3	1.51	0.91
1:B:123:LEU:HD13	1:B:130:LEU:HD13	1.53	0.89
1:D:318:ILE:HD13	1:D:324:VAL:HG22	1.53	0.89
1:D:114:THR:OG1	1:D:118:PRO:HG3	1.74	0.86
1:D:10:GLU:HG3	1:D:287:ASN:HB2	1.57	0.86
1:B:93:MSE:SE	2:B:566:HOH:O	2.45	0.84
1:B:235:LYS:HE2	2:B:687:HOH:O	1.77	0.84
1:C:8:ASN:CG	1:C:285:GLY:HA3	1.99	0.82
1:A:213:LYS:HD3	1:A:228:MSE:HE3	1.61	0.82
1:D:8:ASN:CG	1:D:285:GLY:HA3	1.99	0.82
1:C:10:GLU:HG3	1:C:287:ASN:HB2	1.62	0.82
1:B:213:LYS:HD3	1:B:228:MSE:CE	2.10	0.81
1:B:114:THR:OG1	1:B:118:PRO:HG3	1.82	0.79
1:C:220:ASP:HB3	1:C:223:PHE:O	1.85	0.77
1:B:9:SER:HB3	1:B:35:MSE:HE2	1.66	0.76
1:A:165:VAL:HG22	1:A:172:VAL:HG22	1.67	0.75
1:A:315:SER:HB2	1:A:317:TYR:CE2	2.22	0.75
1:D:220:ASP:HB3	1:D:223:PHE:O	1.86	0.74
1:A:263:SER:HA	1:A:288:PRO:HD2	1.68	0.74
1:C:282:MSE:HG3	2:C:606:HOH:O	1.88	0.72
1:C:1:SER:HB2	1:C:3:PHE:HE1	1.53	0.72
1:D:152:ASN:HB2	2:D:720:HOH:O	1.90	0.71
1:A:63:THR:HG21	1:B:180:MSE:HE1	1.73	0.70
1:C:364:ARG:CZ	1:C:377:ILE:HD11	2.22	0.70
1:C:128:LYS:HE3	2:C:689:HOH:O	1.91	0.69
1:D:199:GLU:HB3	2:D:637:HOH:O	1.92	0.68
1:B:104:THR:HG23	2:B:671:HOH:O	1.93	0.68
1:C:263:SER:HB2	1:C:287:ASN:ND2	2.10	0.67
1:D:235:LYS:HE2	2:D:513:HOH:O	1.94	0.67
1:D:263:SER:HA	1:D:288:PRO:HD2	1.77	0.66
1:D:1:SER:HB2	1:D:3:PHE:HE1	1.61	0.66
1:B:282:MSE:HA	1:B:282:MSE:HE3	1.78	0.66
1:D:364:ARG:CZ	1:D:377:ILE:HD11	2.26	0.66
1:A:275:THR:O	1:A:277:THR:HG23	1.95	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:228:MSE:HE1	1:C:276:ASN:HD21	1.60	0.65
1:A:79:VAL:HG22	1:A:90:VAL:HG22	1.79	0.65
1:A:7:ALA:HA	1:A:14:ILE:HD13	1.78	0.65
1:C:263:SER:HA	1:C:288:PRO:HD2	1.79	0.64
1:B:364:ARG:CZ	1:B:377:ILE:HD11	2.28	0.64
1:A:123:LEU:HD13	1:A:130:LEU:HD13	1.79	0.64
1:B:7:ALA:HA	1:B:14:ILE:HD13	1.79	0.64
1:B:364:ARG:HB2	1:B:377:ILE:CD1	2.28	0.63
1:B:138:LYS:HE2	1:B:158:ARG:HG2	1.81	0.63
1:B:165:VAL:HG22	1:B:172:VAL:HG22	1.79	0.63
1:A:213:LYS:HD3	1:A:228:MSE:CE	2.28	0.63
1:B:291:SER:N	2:B:581:HOH:O	2.31	0.63
1:B:79:VAL:HG22	1:B:90:VAL:HG22	1.81	0.62
1:B:315:SER:HB3	1:B:317:TYR:CE2	2.34	0.62
1:C:289:TYR:HB2	2:C:673:HOH:O	1.99	0.62
1:D:228:MSE:CE	1:D:276:ASN:HD21	2.12	0.62
1:C:229:ILE:HA	1:C:235:LYS:O	1.98	0.62
1:D:263:SER:HB2	1:D:287:ASN:ND2	2.14	0.62
1:B:221:LYS:O	1:B:222:TYR:HB2	2.00	0.61
1:A:352:THR:HG22	2:A:704:HOH:O	1.99	0.61
1:B:213:LYS:HE2	2:B:470:HOH:O	2.00	0.61
1:B:74:SER:HB3	1:B:94:ALA:HB3	1.83	0.61
1:A:293:GLN:HA	2:A:485:HOH:O	2.01	0.60
1:D:96:SER:HA	1:D:118:PRO:CD	2.31	0.60
1:D:93:MSE:HE3	1:D:119:LEU:HD13	1.82	0.60
1:B:5:TYR:CE2	1:B:46:VAL:HG21	2.36	0.60
1:C:226:VAL:HG21	1:C:260:VAL:HG21	1.83	0.60
1:A:221:LYS:O	1:A:222:TYR:HB2	2.02	0.60
1:B:96:SER:HA	1:B:118:PRO:CD	2.32	0.59
1:C:335:GLU:HG2	1:C:368:SER:HB2	1.82	0.59
1:C:1:SER:HB2	1:C:3:PHE:CE1	2.38	0.59
1:D:226:VAL:HG21	1:D:260:VAL:HG21	1.85	0.58
1:C:96:SER:HA	1:C:118:PRO:CD	2.33	0.58
1:B:123:LEU:HD13	1:B:130:LEU:CD1	2.32	0.58
1:D:79:VAL:HG12	1:D:90:VAL:HG22	1.86	0.58
1:A:289:TYR:HA	2:A:482:HOH:O	2.04	0.58
1:A:38:VAL:HG21	1:A:79:VAL:O	2.04	0.58
1:B:9:SER:CB	1:B:35:MSE:CE	2.81	0.57
1:B:336:TRP:CE2	1:B:349:GLN:HG3	2.39	0.57
1:D:1:SER:HB2	1:D:3:PHE:CE1	2.38	0.57
1:A:96:SER:HA	1:A:118:PRO:CD	2.35	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:65:ASN:ND2	2:B:401:HOH:O	2.32	0.56
1:C:300:VAL:HG11	1:D:180:MSE:HE1	1.88	0.56
1:A:93:MSE:SE	2:A:594:HOH:O	2.73	0.56
1:B:38:VAL:HG21	1:B:79:VAL:O	2.05	0.56
1:C:197:LYS:NZ	2:C:686:HOH:O	2.38	0.56
1:D:7:ALA:HA	1:D:14:ILE:HD13	1.86	0.56
1:A:364:ARG:CZ	1:A:377:ILE:HD11	2.36	0.56
1:B:310:LYS:HE2	1:B:327:THR:HG21	1.87	0.56
1:B:8:ASN:CG	1:B:285:GLY:HA3	2.27	0.55
1:B:337:LYS:HB2	2:B:564:HOH:O	2.06	0.55
1:A:8:ASN:CG	1:A:285:GLY:HA3	2.28	0.55
1:A:364:ARG:HB2	1:A:377:ILE:CD1	2.36	0.55
1:A:9:SER:HB3	1:A:35:MSE:HE2	1.88	0.54
1:C:96:SER:HA	1:C:118:PRO:HD2	1.88	0.54
1:D:385:LYS:HD3	2:D:626:HOH:O	2.07	0.54
1:C:7:ALA:HA	1:C:14:ILE:HD13	1.88	0.54
1:A:121:LEU:HA	1:A:131:TYR:O	2.07	0.54
1:B:128:LYS:HE2	2:B:629:HOH:O	2.06	0.54
1:D:190:ASN:ND2	2:D:533:HOH:O	2.30	0.53
1:B:9:SER:HB3	1:B:35:MSE:CE	2.33	0.53
1:B:364:ARG:HB2	1:B:377:ILE:HD13	1.90	0.53
1:A:282:MSE:HE3	1:A:282:MSE:HA	1.89	0.53
1:C:198:VAL:HG12	1:C:236:ILE:HD11	1.91	0.53
1:D:228:MSE:HE2	2:D:552:HOH:O	2.07	0.53
1:A:128:LYS:HE2	2:A:703:HOH:O	2.08	0.53
1:D:289:TYR:HB2	2:D:496:HOH:O	2.09	0.53
1:C:323:PRO:HA	2:C:608:HOH:O	2.09	0.53
1:B:275:THR:O	1:B:277:THR:HG23	2.08	0.53
1:C:1:SER:HB3	2:C:669:HOH:O	2.08	0.52
1:C:262:LEU:HD11	2:C:441:HOH:O	2.09	0.52
1:A:96:SER:HA	1:A:118:PRO:HD2	1.90	0.52
1:D:152:ASN:HD22	1:D:190:ASN:CB	2.22	0.52
1:D:305:PRO:HD3	1:D:369:ASN:ND2	2.23	0.52
1:B:189:THR:OG1	1:B:191:SER:HB2	2.09	0.52
1:D:152:ASN:ND2	1:D:190:ASN:HB3	2.25	0.52
1:D:263:SER:HB2	1:D:287:ASN:HD22	1.75	0.52
1:C:338:TRP:CE2	1:C:365:LEU:HD13	2.45	0.52
1:C:37:ALA:HA	1:C:47:TYR:O	2.10	0.52
1:B:95:SER:O	1:B:97:THR:HG23	2.09	0.52
1:D:34:PRO:HA	1:D:49:ALA:O	2.10	0.52
1:B:121:LEU:HA	1:B:131:TYR:O	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:377:ILE:CG2	1:B:378:SER:N	2.73	0.51
1:A:138:LYS:HE2	1:A:158:ARG:HG2	1.92	0.51
1:D:37:ALA:HA	1:D:47:TYR:O	2.10	0.51
1:C:138:LYS:HG2	1:C:158:ARG:HA	1.92	0.51
1:B:319:PHE:HA	1:B:383:VAL:O	2.10	0.51
1:C:319:PHE:HA	1:C:383:VAL:O	2.10	0.51
1:B:86:LYS:HB3	2:B:671:HOH:O	2.10	0.51
1:B:197:LYS:HG2	2:B:567:HOH:O	2.10	0.51
1:C:35:MSE:HG2	1:C:51:ALA:HB2	1.93	0.51
1:A:79:VAL:HA	1:A:89:TYR:O	2.10	0.51
1:C:263:SER:HB2	1:C:287:ASN:HD22	1.76	0.50
1:D:338:TRP:CE2	1:D:365:LEU:HD13	2.46	0.50
1:B:76:PRO:HA	1:B:91:THR:O	2.11	0.50
1:C:369:ASN:HB2	1:D:222:TYR:OH	2.12	0.50
1:A:95:SER:O	1:A:97:THR:HG23	2.12	0.50
1:D:114:THR:HG1	1:D:118:PRO:HG3	1.74	0.50
1:D:364:ARG:NH1	1:D:377:ILE:HD11	2.27	0.50
1:C:159:SER:HB3	1:C:177:PHE:HB3	1.93	0.50
1:D:159:SER:HB2	1:D:178:ASP:OD2	2.11	0.50
1:C:318:ILE:O	1:C:382:VAL:HA	2.12	0.50
1:D:228:MSE:HE3	1:D:276:ASN:HD21	1.76	0.50
1:D:76:PRO:HA	1:D:91:THR:O	2.11	0.50
1:C:38:VAL:HG12	1:C:79:VAL:HG23	1.93	0.50
1:A:86:LYS:HE2	2:A:694:HOH:O	2.11	0.50
1:D:10:GLU:HG3	1:D:287:ASN:CB	2.37	0.50
1:B:282:MSE:HA	1:B:282:MSE:CE	2.41	0.50
1:A:105:SER:O	1:A:107:THR:HG23	2.12	0.50
1:A:338:TRP:CE2	1:A:365:LEU:HD13	2.47	0.50
1:B:364:ARG:NH1	2:B:605:HOH:O	2.45	0.49
1:A:138:LYS:HG2	1:A:158:ARG:HA	1.94	0.49
1:B:93:MSE:HE1	2:B:684:HOH:O	2.12	0.49
1:D:223:PHE:CE1	2:D:716:HOH:O	2.55	0.49
1:C:38:VAL:HG12	1:C:79:VAL:CG2	2.42	0.49
1:A:266:ASN:HA	1:A:288:PRO:HD3	1.93	0.49
1:B:338:TRP:CE2	1:B:365:LEU:HD13	2.48	0.49
1:B:318:ILE:CD1	1:B:324:VAL:HG22	2.33	0.49
1:D:369:ASN:C	1:D:369:ASN:OD1	2.51	0.49
1:D:163:ILE:HA	1:D:173:TYR:O	2.13	0.49
1:D:138:LYS:HE3	2:D:569:HOH:O	2.12	0.49
1:D:93:MSE:SE	2:D:481:HOH:O	2.81	0.48
1:D:308:ASP:HA	1:D:376:GLN:HE21	1.77	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:217:THR:HB	1:D:246:PRO:HB2	1.94	0.48
1:A:336:TRP:CE2	1:A:349:GLN:HG3	2.49	0.48
1:A:264:PHE:HD2	1:A:265:YCM:HD3	1.78	0.48
1:B:377:ILE:HG22	1:B:378:SER:N	2.28	0.48
1:A:86:LYS:HA	1:A:103:THR:OG1	2.13	0.48
1:C:307:ALA:HB2	1:C:367:VAL:HG22	1.96	0.48
1:A:244:PRO:HB3	2:A:705:HOH:O	2.13	0.48
1:A:76:PRO:HA	1:A:91:THR:O	2.13	0.48
1:A:360:ILE:O	1:A:360:ILE:HG23	2.13	0.48
1:B:138:LYS:HG2	1:B:158:ARG:HA	1.96	0.48
1:D:215:TYR:CE2	1:D:228:MSE:HG3	2.48	0.48
1:C:41:PRO:HD2	2:C:399:HOH:O	2.13	0.48
1:C:79:VAL:HG12	1:C:90:VAL:HG22	1.95	0.47
1:D:224:ASN:ND2	1:D:244:PRO:HA	2.28	0.47
1:A:125:PRO:HG3	2:A:724:HOH:O	2.13	0.47
1:C:217:THR:HB	1:C:246:PRO:HB2	1.95	0.47
1:B:248:GLY:HA3	1:B:290:ALA:O	2.13	0.47
1:B:240:ILE:HD11	1:B:276:ASN:OD1	2.14	0.47
1:A:334:THR:HG21	2:A:473:HOH:O	2.13	0.47
1:A:74:SER:HB3	1:A:94:ALA:HB3	1.97	0.47
1:C:335:GLU:CG	1:C:368:SER:HB2	2.44	0.47
1:A:35:MSE:HE1	1:A:289:TYR:CE1	2.49	0.47
1:C:163:ILE:HA	1:C:173:TYR:O	2.15	0.47
1:D:87:GLN:HG3	2:D:680:HOH:O	2.14	0.47
1:A:93:MSE:HG3	2:A:594:HOH:O	2.15	0.47
1:D:318:ILE:CD1	1:D:324:VAL:HG22	2.36	0.46
1:D:128:LYS:HG3	2:D:748:HOH:O	2.14	0.46
1:D:391:SER:CB	2:D:741:HOH:O	2.63	0.46
1:A:377:ILE:CG2	1:A:378:SER:N	2.77	0.46
1:C:76:PRO:HA	1:C:91:THR:O	2.15	0.46
1:B:79:VAL:HA	1:B:89:TYR:O	2.15	0.46
1:C:137:ASP:O	1:C:139:THR:HG23	2.15	0.46
1:C:281:THR:O	1:C:282:MSE:HE2	2.16	0.46
1:D:152:ASN:HD22	1:D:190:ASN:HB3	1.80	0.46
1:D:152:ASN:ND2	1:D:190:ASN:CB	2.78	0.46
1:A:109:ALA:HA	1:C:109:ALA:O	2.16	0.46
1:C:291:SER:HB2	2:C:603:HOH:O	2.15	0.46
1:B:5:TYR:CD2	1:B:46:VAL:HG21	2.50	0.46
1:C:224:ASN:ND2	1:C:244:PRO:HA	2.31	0.46
1:B:86:LYS:HA	1:B:103:THR:OG1	2.16	0.45
1:C:228:MSE:HE3	1:C:276:ASN:OD1	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:310:LYS:HG3	1:B:327:THR:HG22	1.99	0.45
1:A:242:VAL:HG23	1:A:246:PRO:HG3	1.97	0.45
1:C:121:LEU:HA	1:C:131:TYR:O	2.17	0.45
1:B:291:SER:HB2	2:B:686:HOH:O	2.17	0.45
1:A:115:GLY:HA3	1:A:134:ASN:CG	2.37	0.45
1:A:9:SER:CB	1:A:35:MSE:CE	2.94	0.45
1:A:364:ARG:HB2	1:A:377:ILE:HD13	1.98	0.45
1:C:330:SER:OG	1:C:367:VAL:HG11	2.17	0.45
1:D:121:LEU:N	1:D:121:LEU:HD23	2.32	0.45
1:A:282:MSE:CE	1:A:282:MSE:HA	2.47	0.45
1:B:200:ALA:HB1	1:B:219:VAL:O	2.16	0.44
1:A:377:ILE:HG22	1:A:378:SER:N	2.32	0.44
1:D:251:VAL:HG22	1:D:258:VAL:HG22	1.97	0.44
1:D:390:PRO:HA	2:D:698:HOH:O	2.16	0.44
1:A:9:SER:HB3	1:A:35:MSE:CE	2.47	0.44
1:D:282:MSE:HG2	2:D:739:HOH:O	2.17	0.44
1:B:215:TYR:CE2	1:B:228:MSE:HG2	2.52	0.44
1:C:364:ARG:HB2	1:C:377:ILE:CD1	2.48	0.44
1:B:221:LYS:HG3	1:B:222:TYR:CD1	2.53	0.44
1:D:121:LEU:HA	1:D:131:TYR:O	2.18	0.44
1:C:328:ASP:N	1:C:350:ASN:OD1	2.43	0.44
1:C:157:GLY:H	1:C:176:ASN:ND2	2.16	0.44
1:A:248:GLY:HA3	1:A:290:ALA:O	2.18	0.43
1:A:65:ASN:ND2	2:A:504:HOH:O	2.51	0.43
1:A:323:PRO:HA	1:A:354:THR:HG22	2.00	0.43
1:D:10:GLU:HB2	1:D:286:LYS:CG	2.48	0.43
1:A:224:ASN:ND2	1:A:244:PRO:HA	2.32	0.43
1:A:175:ALA:HB2	1:A:205:ILE:HD11	2.00	0.43
1:B:318:ILE:O	1:B:382:VAL:HA	2.19	0.43
1:D:134:ASN:HB3	1:D:137:ASP:HB2	2.00	0.43
1:A:373:THR:CG2	1:A:374:ASP:N	2.81	0.43
1:B:360:ILE:HG23	1:B:360:ILE:O	2.19	0.43
1:C:84:ASP:CG	1:C:86:LYS:HG3	2.39	0.43
1:D:138:LYS:HG2	1:D:158:ARG:HA	2.01	0.43
1:C:137:ASP:O	1:C:138:LYS:HB2	2.19	0.43
1:C:389:THR:HG21	2:C:666:HOH:O	2.18	0.43
1:B:123:LEU:HD12	1:B:130:LEU:HA	2.00	0.43
1:B:161:LYS:NZ	2:B:690:HOH:O	2.37	0.43
1:A:258:VAL:HG23	1:A:273:THR:CG2	2.49	0.42
1:C:48:VAL:O	1:C:56:VAL:HA	2.19	0.42
1:C:34:PRO:HA	1:C:49:ALA:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:8:ASN:CB	1:D:285:GLY:HA3	2.49	0.42
1:D:272:ASP:OD1	1:D:274:ALA:HB3	2.20	0.42
1:B:305:PRO:HD3	1:B:369:ASN:ND2	2.34	0.42
1:D:326:PHE:HB3	1:D:338:TRP:CH2	2.55	0.42
1:B:210:GLU:HG3	2:B:587:HOH:O	2.18	0.42
1:A:9:SER:O	1:A:33:ASN:HA	2.20	0.42
1:B:327:THR:O	1:B:327:THR:HG23	2.19	0.42
1:D:246:PRO:HA	1:D:262:LEU:HD23	2.00	0.42
1:A:258:VAL:HB	1:A:271:ILE:HB	2.00	0.42
1:B:115:GLY:HA3	1:B:134:ASN:CG	2.40	0.42
1:C:8:ASN:CB	1:C:285:GLY:HA3	2.50	0.42
1:A:309:PHE:CZ	1:A:378:SER:HB2	2.55	0.42
1:D:266:ASN:HA	1:D:288:PRO:HD3	2.02	0.42
1:D:251:VAL:HG12	1:D:252:THR:O	2.19	0.42
1:A:312:ASN:HD22	1:A:313:ILE:N	2.18	0.42
1:D:318:ILE:O	1:D:382:VAL:HA	2.20	0.42
1:A:281:THR:O	1:A:282:MSE:HE3	2.20	0.42
1:B:258:VAL:HG23	1:B:273:THR:CG2	2.50	0.42
1:A:129:LYS:HG3	2:A:617:HOH:O	2.20	0.42
1:D:364:ARG:HB2	1:D:377:ILE:CD1	2.49	0.41
1:C:138:LYS:HD3	1:C:158:ARG:HG2	2.02	0.41
1:C:205:ILE:HG12	1:C:216:VAL:HG22	2.02	0.41
1:D:137:ASP:O	1:D:139:THR:HG23	2.19	0.41
1:D:157:GLY:H	1:D:176:ASN:ND2	2.17	0.41
1:A:221:LYS:O	1:A:222:TYR:CB	2.67	0.41
1:A:38:VAL:HG13	1:A:293:GLN:HB2	2.01	0.41
1:A:307:ALA:HB2	1:A:367:VAL:HG22	2.03	0.41
1:C:364:ARG:NH2	1:C:377:ILE:HD11	2.35	0.41
1:B:330:SER:OG	1:B:367:VAL:HG11	2.20	0.41
1:D:96:SER:HA	1:D:118:PRO:HD2	2.02	0.41
1:C:313:ILE:HG12	1:C:318:ILE:CD1	2.51	0.41
1:C:73:GLY:HA3	1:C:92:ASN:CG	2.41	0.41
1:B:35:MSE:HE1	1:B:289:TYR:CE1	2.56	0.41
1:A:258:VAL:HG23	1:A:273:THR:HG22	2.02	0.41
1:A:312:ASN:HD22	1:A:312:ASN:C	2.24	0.41
1:A:305:PRO:HD3	1:A:369:ASN:ND2	2.36	0.41
1:D:327:THR:O	1:D:327:THR:HG23	2.21	0.41
1:D:48:VAL:O	1:D:56:VAL:HA	2.21	0.41
1:A:200:ALA:HB1	1:A:219:VAL:O	2.20	0.41
1:D:198:VAL:HG12	1:D:236:ILE:HD11	2.02	0.41
1:D:6:ILE:HG23	1:D:290:ALA:HB2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:276:ASN:HB2	2:D:571:HOH:O	2.21	0.41
1:A:39:ILE:HD11	1:A:43:GLY:HA2	2.03	0.41
1:B:325:GLN:HB3	2:B:640:HOH:O	2.20	0.41
1:A:284:VAL:HB	1:A:285:GLY:H	1.53	0.40
1:C:77:GLN:NE2	2:C:678:HOH:O	2.53	0.40
1:B:41:PRO:HD2	2:B:398:HOH:O	2.20	0.40
1:B:30:VAL:HB	1:B:31:GLY:H	1.61	0.40
1:B:266:ASN:HA	1:B:288:PRO:HD3	2.04	0.40
1:B:199:GLU:HG3	2:B:502:HOH:O	2.20	0.40
1:D:35:MSE:HG2	1:D:51:ALA:HB2	2.04	0.40
1:C:242:VAL:CG2	1:C:246:PRO:HG3	2.51	0.40
1:C:219:VAL:HG22	1:C:246:PRO:HD2	2.03	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	388/391 (99%)	367 (95%)	19 (5%)	2 (0%)	34	48
1	B	388/391 (99%)	367 (95%)	19 (5%)	2 (0%)	34	48
1	C	388/391 (99%)	371 (96%)	15 (4%)	2 (0%)	34	48
1	D	388/391 (99%)	369 (95%)	16 (4%)	3 (1%)	24	35
All	All	1552/1564 (99%)	1474 (95%)	69 (4%)	9 (1%)	30	43

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	30	VAL
1	A	284	VAL
1	B	30	VAL

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Mol	Chain	Res	Type
1	B	284	VAL
1	C	284	VAL
1	C	30	VAL
1	D	284	VAL
1	D	30	VAL
1	D	390	PRO

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	331/326 (102%)	322 (97%)	9 (3%)	52	73
1	B	331/326 (102%)	325 (98%)	6 (2%)	66	84
1	C	331/326 (102%)	325 (98%)	6 (2%)	66	84
1	D	331/326 (102%)	326 (98%)	5 (2%)	72	87
All	All	1324/1304 (102%)	1298 (98%)	26 (2%)	63	81

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	VAL
1	A	74	SER
1	A	137	ASP
1	A	282	MSE
1	A	312	ASN
1	A	315	SER
1	A	352	THR
1	A	368	SER
1	A	375	SER
1	B	11	SER
1	B	191	SER
1	B	282	MSE
1	B	352	THR
1	B	368	SER
1	B	375	SER

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Mol	Chain	Res	Type
1	C	98	LEU
1	C	128	LYS
1	C	159	SER
1	C	287	ASN
1	C	318	ILE
1	C	335	GLU
1	D	137	ASP
1	D	287	ASN
1	D	291	SER
1	D	375	SER
1	D	376	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	65	ASN
1	A	77	GLN
1	A	176	ASN
1	A	190	ASN
1	A	312	ASN
1	B	22	ASN
1	B	65	ASN
1	B	176	ASN
1	B	353	HIS
1	B	371	ASN
1	C	65	ASN
1	C	176	ASN
1	C	190	ASN
1	C	287	ASN
1	C	371	ASN
1	D	65	ASN
1	D	152	ASN
1	D	176	ASN
1	D	287	ASN
1	D	371	ASN
1	D	376	GLN

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
1	YCM	A	265	1	8,9,10	0.43	0	5,10,12	1.12	1 (20%)
1	YCM	B	265	1	8,9,10	0.44	0	5,10,12	1.13	1 (20%)
1	YCM	C	265	1	8,9,10	0.40	0	5,10,12	1.05	1 (20%)
1	YCM	D	265	1	8,9,10	0.43	0	5,10,12	1.01	1 (20%)

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
1	YCM	A	265	1	-	0/6/8/10	0/0/0/0
1	YCM	B	265	1	-	0/6/8/10	0/0/0/0
1	YCM	C	265	1	-	0/6/8/10	0/0/0/0
1	YCM	D	265	1	-	0/6/8/10	0/0/0/0

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	265	YCM	O-C-CA	-2.40	119.25	125.49
1	A	265	YCM	O-C-CA	-2.26	119.61	125.49
1	D	265	YCM	O-C-CA	-2.03	120.20	125.49
1	C	265	YCM	O-C-CA	-2.02	120.23	125.49

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	265	YCM	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	385/391 (98%)	-0.42	3 (0%) 87 87	4, 19, 37, 57	0
1	B	385/391 (98%)	-0.40	5 (1%) 79 79	4, 18, 46, 63	0
1	C	385/391 (98%)	-0.21	20 (5%) 31 31	4, 19, 59, 70	0
1	D	385/391 (98%)	-0.46	2 (0%) 91 91	3, 16, 36, 55	0
All	All	1540/1564 (98%)	-0.37	30 (1%) 70 69	3, 18, 47, 70	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	321	SER	7.1
1	C	314	THR	5.5
1	B	314	THR	4.1
1	C	391	SER	3.9
1	B	321	SER	3.8
1	A	391	SER	3.4
1	C	323	PRO	3.4
1	D	391	SER	3.4
1	A	314	THR	3.3
1	C	311	SER	3.3
1	C	317	TYR	3.1
1	C	331	LYS	3.0
1	D	314	THR	3.0
1	C	352	THR	3.0
1	C	344	SER	2.8
1	C	356	SER	2.8
1	C	361	TYR	2.7
1	C	355	TYR	2.6
1	C	354	THR	2.5
1	C	315	SER	2.5
1	C	325	GLN	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	322	GLU	2.4
1	C	312	ASN	2.3
1	C	310	LYS	2.3
1	B	344	SER	2.2
1	C	351	PRO	2.2
1	B	354	THR	2.1
1	B	391	SER	2.1
1	C	324	VAL	2.1
1	A	360	ILE	2.0

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
1	YCM	C	265	10/11	0.97	0.09	-	11,17,31,32	0
1	YCM	D	265	10/11	0.96	0.10	-	11,14,24,25	0
1	YCM	A	265	10/11	0.96	0.12	-	15,20,28,31	0
1	YCM	B	265	10/11	0.95	0.11	-	13,17,25,28	0

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.