



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:37 PM GMT

PDB ID : 3RIP  
Title : Crystal Structure of human gamma-tubulin complex protein 4 (GCP4)  
Authors : Gregory-Paaron, L.; Guillet, V.; Mourey, L.  
Deposited on : 2011-04-14  
Resolution : 2.30 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  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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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

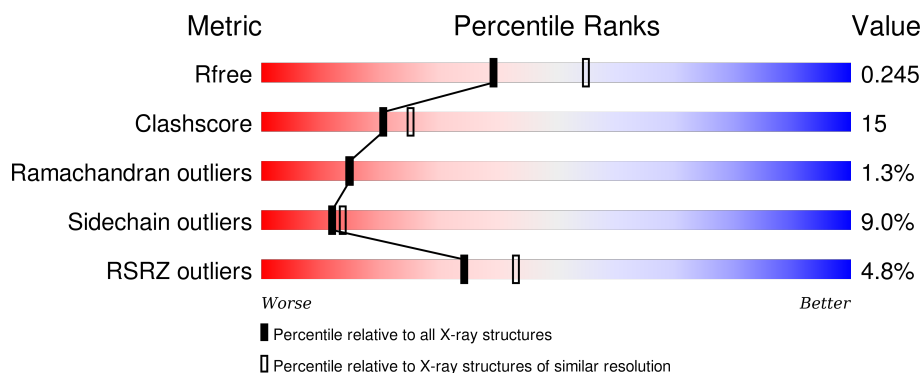
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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  $\geq 3$ , 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	677	<div> <div>4%</div> <div>61%</div> <div>18%</div> <div>5%</div> <div>16%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	GOL	A	683	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	GOL	A	685	-	-	-	X
3	GOL	A	686	-	-	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4843 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 Gamma-tubulin complex component 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	568	4572	2963	774	816	19	0	4	0

There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	667	CYS	-	EXPRESSION TAG	UNP Q9UGJ1
A	668	GLY	-	EXPRESSION TAG	UNP Q9UGJ1
A	669	ARG	-	EXPRESSION TAG	UNP Q9UGJ1
A	670	LEU	-	EXPRESSION TAG	UNP Q9UGJ1
A	671	GLU	-	EXPRESSION TAG	UNP Q9UGJ1
A	672	HIS	-	EXPRESSION TAG	UNP Q9UGJ1
A	673	HIS	-	EXPRESSION TAG	UNP Q9UGJ1
A	674	HIS	-	EXPRESSION TAG	UNP Q9UGJ1
A	675	HIS	-	EXPRESSION TAG	UNP Q9UGJ1
A	676	HIS	-	EXPRESSION TAG	UNP Q9UGJ1
A	677	HIS	-	EXPRESSION TAG	UNP Q9UGJ1

- Molecule 2 is (4R)-2-METHYLPENTANE-2,4-DIOL (three-letter code: MRD) (formula: C<sub>6</sub>H<sub>14</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			8	6	2		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		

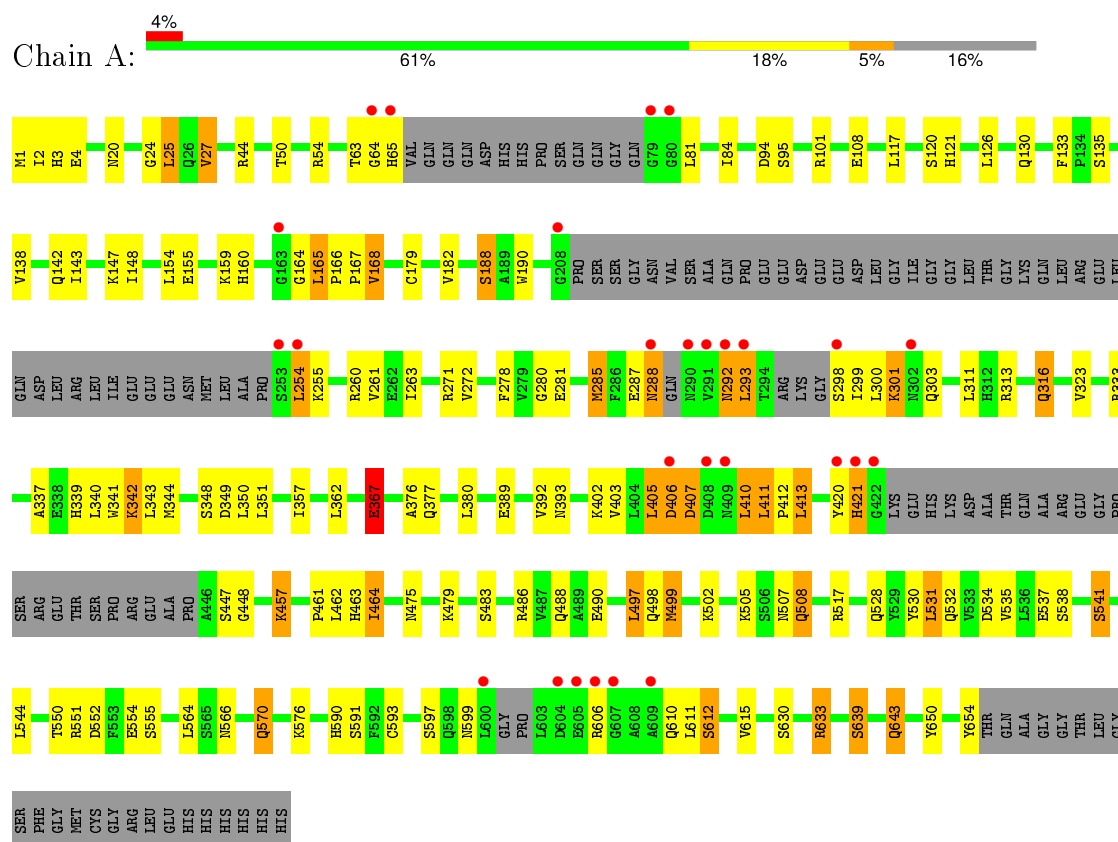
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	215	Total	O	0	2
			215	215		

### 3 Residue-property plots [i](#)

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: Gamma-tubulin complex component 4



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 63 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	214.95Å 214.95Å 128.66Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.30 34.51 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.5 (20.00-2.30) 99.4 (34.51-2.20)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.85 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.5.0072	Depositor
R, $R_{free}$	0.227 , 0.260 0.216 , 0.245	Depositor DCC
$R_{free}$ test set	3848 reflections (5.25%)	DCC
Wilson B-factor (Å <sup>2</sup> )	49.4	Xtriage
Anisotropy	0.371	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 55.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 88024 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	4843	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.54% 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, 0.375 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: MRD, GOL

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	1.01	2/4676 (0.0%)	1.01	16/6334 (0.3%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	367[A]	GLU	CD-OE2	8.72	1.35	1.25
1	A	367[B]	GLU	CD-OE2	8.72	1.35	1.25

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	633	ARG	NE-CZ-NH2	-16.28	112.16	120.30
1	A	633	ARG	NE-CZ-NH1	12.51	126.56	120.30
1	A	367[A]	GLU	CG-CD-OE1	10.78	139.85	118.30
1	A	367[B]	GLU	CG-CD-OE1	10.78	139.85	118.30
1	A	367[A]	GLU	OE1-CD-OE2	-8.23	113.42	123.30
1	A	367[B]	GLU	OE1-CD-OE2	-8.23	113.42	123.30
1	A	413	LEU	CA-CB-CG	7.82	133.29	115.30
1	A	27	VAL	CB-CA-C	-7.23	97.66	111.40
1	A	54	ARG	NE-CZ-NH2	-6.65	116.98	120.30
1	A	260	ARG	NE-CZ-NH1	6.21	123.41	120.30
1	A	367[A]	GLU	CG-CD-OE2	-5.93	106.45	118.30
1	A	367[B]	GLU	CG-CD-OE2	-5.93	106.45	118.30
1	A	260	ARG	NE-CZ-NH2	-5.39	117.61	120.30
1	A	54	ARG	NE-CZ-NH1	5.23	122.92	120.30
1	A	643	GLN	CB-CA-C	-5.21	99.99	110.40
1	A	133	PHE	C-N-CD	5.19	139.30	128.40

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	4572	0	4505	132	0
2	A	8	0	14	4	0
3	A	48	0	64	4	0
4	A	215	0	0	11	0
All	All	4843	0	4583	134	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 15.

All (134) 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:A:612:SER:HB3	4:A:888:HOH:O	1.44	1.14
1:A:288:ASN:HD22	1:A:288:ASN:H	1.11	0.99
1:A:410:LEU:HD23	1:A:411:LEU:N	1.80	0.96
1:A:550:THR:HG22	1:A:552:ASP:H	1.35	0.91
1:A:639:SER:O	1:A:643:GLN:HG2	1.73	0.89
1:A:1[B]:MET:HE2	1:A:3:HIS:H	1.39	0.85
1:A:497:LEU:HD23	1:A:497:LEU:C	1.98	0.83
1:A:164:GLY:O	1:A:165:LEU:HG	1.80	0.82
1:A:316:GLN:OE1	1:A:316:GLN:HA	1.81	0.81
1:A:288:ASN:ND2	1:A:288:ASN:H	1.79	0.79
1:A:117:LEU:HD11	1:A:121:HIS:HD2	1.48	0.79
1:A:606:ARG:O	1:A:610:GLN:HG2	1.85	0.77
1:A:534:ASP:OD2	1:A:633:ARG:CD	2.34	0.75
1:A:502[A]:LYS:HG2	4:A:901:HOH:O	1.86	0.74
1:A:188:SER:HB2	1:A:311:LEU:HD12	1.70	0.74
1:A:165:LEU:HB2	1:A:168:VAL:HG13	1.72	0.71
1:A:316:GLN:OE1	1:A:316:GLN:CA	2.39	0.70
1:A:505:LYS:NZ	1:A:599:ASN:OD1	2.22	0.70
1:A:285:MET:SD	1:A:285:MET:C	2.70	0.70
1:A:508:GLN:HA	1:A:508:GLN:NE2	2.06	0.70
1:A:534:ASP:OD2	1:A:633:ARG:HD2	1.93	0.69
2:A:678:MRD:H5C2	2:A:678:MRD:O2	1.92	0.69
1:A:117:LEU:HD11	1:A:121:HIS:CD2	2.29	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:135:SER:O	1:A:138:VAL:HG22	1.95	0.67
1:A:497:LEU:HD23	1:A:497:LEU:O	1.96	0.65
1:A:534:ASP:OD2	1:A:633:ARG:HD3	1.97	0.65
1:A:142:GLN:OE1	1:A:160:HIS:NE2	2.28	0.64
1:A:554:GLU:OE1	3:A:685:GOL:O3	2.15	0.64
1:A:84:ILE:HD12	1:A:182:VAL:HG13	1.80	0.64
1:A:20:ASN:O	1:A:24:GLY:N	2.27	0.64
1:A:650:TYR:CE1	1:A:654:TYR:HB2	2.34	0.63
1:A:81:LEU:HD21	1:A:147:LYS:HA	1.80	0.62
1:A:63:THR:O	1:A:65:HIS:N	2.26	0.61
1:A:285:MET:SD	1:A:285:MET:O	2.60	0.59
1:A:566:ASN:O	1:A:570:GLN:HG2	2.03	0.59
1:A:1[B]:MET:CE	1:A:3:HIS:H	2.11	0.59
1:A:190:TRP:CE2	1:A:280:GLY:HA3	2.37	0.59
1:A:287:GLU:HG3	1:A:287:GLU:O	2.04	0.58
1:A:410:LEU:HD23	1:A:411:LEU:H	1.66	0.57
1:A:497:LEU:CD2	1:A:497:LEU:C	2.70	0.57
1:A:550:THR:HG21	1:A:555:SER:CB	2.34	0.57
1:A:550:THR:HG21	1:A:555:SER:HB2	1.86	0.57
1:A:410:LEU:HD23	1:A:410:LEU:C	2.25	0.56
1:A:508:GLN:NE2	1:A:508:GLN:CA	2.68	0.56
1:A:288:ASN:ND2	1:A:288:ASN:N	2.51	0.55
1:A:531:LEU:O	1:A:535:VAL:HB	2.06	0.54
1:A:411:LEU:N	1:A:412:PRO:HD2	2.23	0.54
1:A:344:MET:HE2	1:A:461:PRO:HG2	1.89	0.53
1:A:287:GLU:CG	1:A:287:GLU:O	2.56	0.53
1:A:281:GLU:OE2	4:A:834:HOH:O	2.18	0.53
2:A:678:MRD:O2	2:A:678:MRD:C5	2.56	0.53
1:A:508:GLN:CA	1:A:508:GLN:HE21	2.21	0.53
1:A:143:ILE:HA	1:A:148:ILE:HG13	1.91	0.53
1:A:488:GLN:OE1	1:A:528:GLN:NE2	2.36	0.53
1:A:271:ARG:HG2	1:A:272:VAL:N	2.23	0.52
1:A:1[B]:MET:HE2	1:A:2:ILE:N	2.25	0.52
1:A:155:GLU:OE2	1:A:159:LYS:HE2	2.09	0.52
1:A:166:PRO:HB2	1:A:167:PRO:HD3	1.90	0.52
1:A:505:LYS:HE2	4:A:692:HOH:O	2.09	0.51
1:A:188:SER:HB2	1:A:311:LEU:CD1	2.41	0.50
1:A:108:GLU:HG3	4:A:792:HOH:O	2.10	0.50
1:A:341:TRP:NE1	1:A:550:THR:O	2.45	0.50
1:A:550:THR:HG22	1:A:551:ARG:N	2.27	0.49
1:A:420:TYR:O	1:A:421:HIS:HB2	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:593:CYS:SG	2:A:678:MRD:H5C1	2.53	0.49
1:A:84:ILE:HD12	1:A:182:VAL:CG1	2.41	0.49
1:A:537:GLU:O	1:A:541:SER:HB2	2.12	0.49
1:A:271:ARG:CG	1:A:272:VAL:N	2.75	0.49
1:A:486:ARG:O	1:A:490:GLU:HG3	2.12	0.49
1:A:1[B]:MET:HE2	1:A:3:HIS:N	2.18	0.49
1:A:502[B]:LYS:HE3	4:A:811:HOH:O	2.13	0.48
1:A:405:LEU:HD12	1:A:406:ASP:N	2.28	0.48
1:A:313:ARG:O	1:A:316:GLN:HB2	2.13	0.48
1:A:532:GLN:O	1:A:537:GLU:HG2	2.13	0.48
1:A:502[A]:LYS:HE3	4:A:727:HOH:O	2.13	0.48
1:A:254:LEU:O	1:A:255:LYS:HB2	2.14	0.48
1:A:126:LEU:O	1:A:130:GLN:HG3	2.14	0.48
1:A:340:LEU:HA	1:A:340:LEU:HD23	1.69	0.48
1:A:463:HIS:N	1:A:463:HIS:ND1	2.62	0.47
1:A:643:GLN:HG2	1:A:643:GLN:H	1.58	0.47
1:A:292:ASN:O	1:A:293:LEU:CB	2.62	0.47
1:A:389:GLU:O	1:A:393:ASN:ND2	2.41	0.47
1:A:351:LEU:HA	1:A:351:LEU:HD23	1.60	0.46
1:A:333:ARG:HD2	3:A:685:GOL:H11	1.97	0.46
1:A:479:LYS:NZ	4:A:830:HOH:O	2.49	0.46
1:A:1[B]:MET:HG2	1:A:4:GLU:HG2	1.98	0.45
1:A:285:MET:CG	1:A:285:MET:O	2.65	0.45
1:A:117:LEU:CD1	1:A:121:HIS:CD2	2.99	0.45
1:A:299:ILE:O	1:A:339:HIS:ND1	2.49	0.45
1:A:263:ILE:O	1:A:263:ILE:HG23	2.17	0.45
1:A:457:LYS:N	1:A:457:LYS:CD	2.80	0.45
1:A:341:TRP:HE3	1:A:342:LYS:N	2.14	0.45
1:A:498:GLN:CD	1:A:593:CYS:HB2	2.37	0.45
1:A:337:ALA:HB1	1:A:552:ASP:OD1	2.17	0.45
1:A:281:GLU:OE1	3:A:686:GOL:O2	2.28	0.45
1:A:410:LEU:CD2	1:A:411:LEU:N	2.66	0.45
1:A:351:LEU:HD22	1:A:544:LEU:HD22	1.98	0.45
1:A:566:ASN:O	1:A:570:GLN:CG	2.64	0.45
1:A:344:MET:HE2	1:A:462:LEU:HD23	1.97	0.45
1:A:154:LEU:HD23	1:A:179:CYS:HB3	1.98	0.45
1:A:406:ASP:O	1:A:407:ASP:C	2.55	0.45
1:A:570:GLN:H	1:A:570:GLN:HG2	1.59	0.44
1:A:278:PHE:HB2	3:A:686:GOL:H2	1.98	0.44
1:A:342:LYS:O	1:A:343:LEU:C	2.55	0.44
1:A:94:ASP:OD1	1:A:101:ARG:NH2	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:341:TRP:CZ2	1:A:551:ARG:NH1	2.85	0.44
1:A:507:ASN:OD1	1:A:507:ASN:C	2.56	0.44
1:A:344:MET:CE	1:A:461:PRO:HG2	2.48	0.44
1:A:155:GLU:O	1:A:159:LYS:HG2	2.18	0.44
1:A:499:MET:CE	1:A:517:ARG:HD3	2.48	0.44
1:A:25:LEU:HB2	1:A:50:THR:OG1	2.18	0.44
1:A:349:ASP:HA	4:A:720:HOH:O	2.18	0.43
1:A:461:PRO:O	1:A:464:ILE:HG23	2.19	0.43
1:A:499:MET:HE2	1:A:517:ARG:HD3	2.01	0.42
1:A:590:HIS:ND1	2:A:678:MRD:H5C3	2.35	0.42
1:A:344:MET:HE1	1:A:461:PRO:O	2.19	0.42
1:A:448:GLY:HA3	4:A:900:HOH:O	2.20	0.41
1:A:164:GLY:O	1:A:165:LEU:CG	2.60	0.41
1:A:350:LEU:N	4:A:720:HOH:O	2.54	0.41
1:A:611:LEU:O	1:A:615:VAL:HG23	2.20	0.41
1:A:402:LYS:HG2	1:A:402:LYS:O	2.20	0.41
1:A:380:LEU:HD21	1:A:392:VAL:HG21	2.03	0.41
1:A:164:GLY:O	1:A:165:LEU:CB	2.69	0.41
1:A:300:LEU:O	1:A:301:LYS:CB	2.69	0.40
1:A:530:TYR:CD1	1:A:633:ARG:HD3	2.56	0.40
1:A:376:ALA:O	1:A:377:GLN:C	2.58	0.40
1:A:475:ASN:O	1:A:479:LYS:HG3	2.22	0.40
1:A:457:LYS:HD2	1:A:457:LYS:N	2.36	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	557/677 (82%)	525 (94%)	25 (4%)	7 (1%)	15 15

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	165	LEU
1	A	292	ASN
1	A	293	LEU
1	A	301	LYS
1	A	407	ASP
1	A	64	GLY
1	A	421	HIS

### 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	494/604 (82%)	449 (91%)	45 (9%)	12	13

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	LEU
1	A	27	VAL
1	A	44	ARG
1	A	95	SER
1	A	120	SER
1	A	168	VAL
1	A	188	SER
1	A	254	LEU
1	A	261	VAL
1	A	285	MET
1	A	288	ASN
1	A	298	SER
1	A	303	GLN
1	A	316	GLN
1	A	323	VAL
1	A	342	LYS
1	A	348	SER
1	A	357	ILE
1	A	362	LEU
1	A	367[A]	GLU

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Mol	Chain	Res	Type
1	A	367[B]	GLU
1	A	403	VAL
1	A	405	LEU
1	A	406	ASP
1	A	410	LEU
1	A	411	LEU
1	A	413	LEU
1	A	447	SER
1	A	457	LYS
1	A	464	ILE
1	A	483	SER
1	A	497	LEU
1	A	499	MET
1	A	508	GLN
1	A	531	LEU
1	A	538	SER
1	A	541	SER
1	A	564	LEU
1	A	570	GLN
1	A	576	LYS
1	A	591	SER
1	A	597	SER
1	A	612	SER
1	A	630	SER
1	A	639	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	61	GLN
1	A	121	HIS
1	A	207	GLN
1	A	288	ASN
1	A	303	GLN
1	A	508	GLN
1	A	539	GLN
1	A	548	ASN
1	A	566	ASN
1	A	635	HIS
1	A	636	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

9 ligands 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	MRD	A	678	-	6,7,7	1.04	0	7,10,10	0.21	0
3	GOL	A	679	-	5,5,5	0.66	0	5,5,5	1.25	0
3	GOL	A	680	-	5,5,5	0.80	0	5,5,5	0.95	0
3	GOL	A	681	-	5,5,5	0.67	0	5,5,5	1.44	0
3	GOL	A	682	-	5,5,5	0.53	0	5,5,5	0.61	0
3	GOL	A	683	-	5,5,5	0.76	0	5,5,5	0.70	0
3	GOL	A	684	-	5,5,5	1.21	0	5,5,5	1.26	1 (20%)
3	GOL	A	685	-	5,5,5	0.56	0	5,5,5	0.69	0
3	GOL	A	686	-	5,5,5	0.41	0	5,5,5	0.62	0

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	MRD	A	678	-	-	0/5/5/5	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	A	679	-	-	0/4/4/4	0/0/0/0
3	GOL	A	680	-	-	0/4/4/4	0/0/0/0
3	GOL	A	681	-	-	0/4/4/4	0/0/0/0
3	GOL	A	682	-	-	0/4/4/4	0/0/0/0
3	GOL	A	683	-	-	0/4/4/4	0/0/0/0
3	GOL	A	684	-	-	0/4/4/4	0/0/0/0
3	GOL	A	685	-	-	0/4/4/4	0/0/0/0
3	GOL	A	686	-	-	0/4/4/4	0/0/0/0

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	684	GOL	C3-C2-C1	-2.56	101.07	111.12

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
2	A	678	MRD	4	0
3	A	685	GOL	2	0
3	A	686	GOL	2	0

## 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, 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	568/677 (83%)	-0.11	27 (4%) 34 43	29, 51, 82, 119	2 (0%)

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	253	SER	8.6
1	A	79	GLY	7.9
1	A	422	GLY	4.8
1	A	302	ASN	4.6
1	A	291	VAL	4.3
1	A	64	GLY	4.2
1	A	65	HIS	4.1
1	A	254	LEU	4.1
1	A	208	GLY	4.1
1	A	421	HIS	3.8
1	A	80	GLY	3.8
1	A	292	ASN	3.8
1	A	606	ARG	3.8
1	A	293	LEU	3.7
1	A	409	ASN	3.6
1	A	406	ASP	3.4
1	A	420	TYR	3.2
1	A	290	ASN	2.8
1	A	408	ASP	2.6
1	A	298	SER	2.6
1	A	600	LEU	2.5
1	A	163	GLY	2.3
1	A	605	GLU	2.3
1	A	604	ASP	2.2
1	A	288	ASN	2.2
1	A	607	GLY	2.2
1	A	609	ALA	2.1

## 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 carbohydrates 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. 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, 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	GOL	A	686	6/6	0.73	0.48	26.99	63,69,71,74	0
3	GOL	A	685	6/6	0.76	0.35	17.58	63,72,76,79	0
3	GOL	A	683	6/6	0.80	0.29	5.41	39,41,45,46	6
3	GOL	A	679	6/6	0.84	0.17	0.72	43,45,49,49	0
3	GOL	A	682	6/6	0.95	0.10	-1.19	39,43,46,46	0
3	GOL	A	681	6/6	0.92	0.09	-1.35	46,48,48,49	0
2	MRD	A	678	8/8	0.74	0.29	-	56,68,77,83	0
3	GOL	A	684	6/6	0.97	0.07	-	33,35,43,46	6
3	GOL	A	680	6/6	0.86	0.19	-	34,36,39,40	6

## 6.5 Other polymers [i](#)

There are no such residues in this entry.