



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2014 – 12:30 AM GMT

PDB ID : 2J62  
Title : STRUCTURE OF A BACTERIAL O-GLCNACASE IN COMPLEX WITH  
GLCNACSTATIN  
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Deposited on : 2006-09-22  
Resolution : 2.26 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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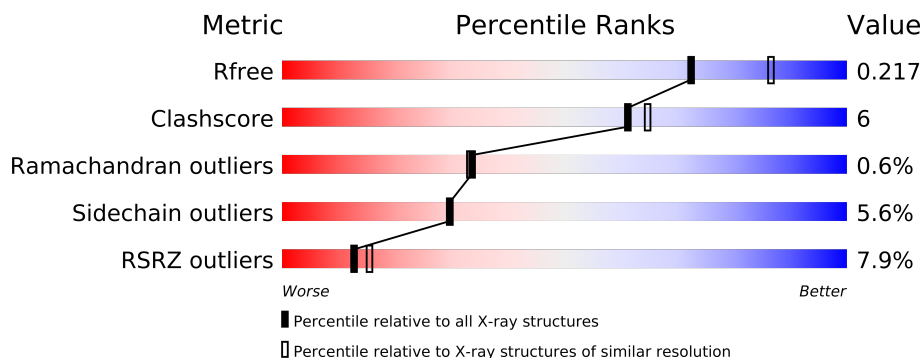
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.26 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))
$R_{free}$	66092	1108 (2.28-2.24)
Clashscore	79885	1326 (2.28-2.24)
Ramachandran outliers	78287	1291 (2.28-2.24)
Sidechain outliers	78261	1291 (2.28-2.24)
RSRZ outliers	66119	1110 (2.28-2.24)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	594	
1	B	594	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	CL	B	1626	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9774 atoms, of which 0 are hydrogen and 0 are deuterium.

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 HYALURONIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	585	Total	C	N	O	S	0	0	0
			4628	2913	756	942	17			
1	B	585	Total	C	N	O	S	0	0	0
			4629	2913	757	942	17			

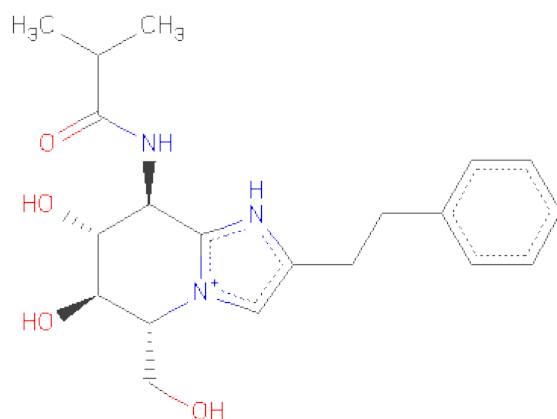
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	388	ASP	ASN	CONFLICT	UNP Q0TR53
B	388	ASP	ASN	CONFLICT	UNP Q0TR53

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Cl	0	0
			2	2		
2	A	2	Total	Cl	0	0
			2	2		

- Molecule 3 is N-[(5R,6R,7R,8S)-6,7-DIHYDROXY-5-(HYDROXYMETHYL)-2-(2-PHENYLETHYL)-1,5,6,7,8,8A-HEXAHYDROIMIDAZO[1,2-A]PYRIDIN-8-YL]-2-METHYLPROPANAMIDE (three-letter code: GSZ) (formula: C<sub>20</sub>H<sub>28</sub>N<sub>3</sub>O<sub>4</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			27	20	3	4		
3	B	1	Total	C	N	O	0	0
			27	20	3	4		

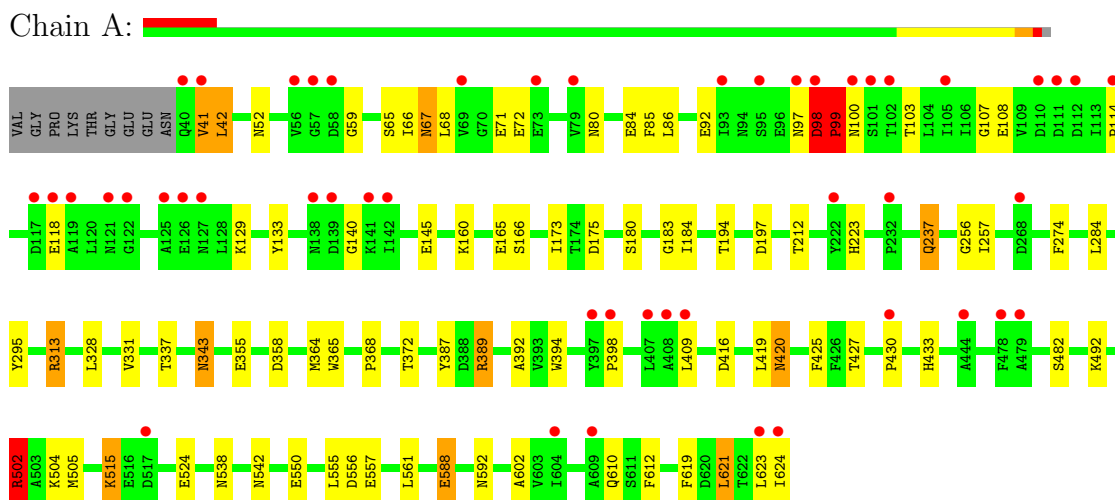
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	198	Total	O	0	0
			198	198		
4	B	261	Total	O	0	0
			261	261		

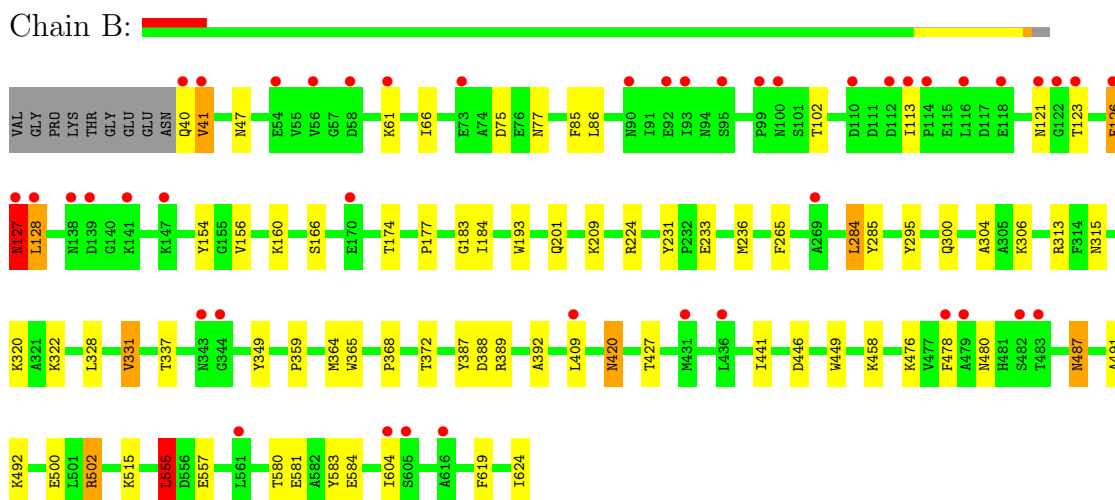
### 3 Residue-property plots

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: HYALURONIDASE



#### • Molecule 1: HYALURONIDASE



## 4 Data and refinement statistics

Property	Value	Source
Space group	I 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	130.25Å 145.92Å 152.57Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.26 20.00 – 2.26	Depositor EDS
% Data completeness (in resolution range)	99.6 (20.00-2.26) 99.6 (20.00-2.26)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.80 (at 2.26Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.179 , 0.219 0.179 , 0.217	Depositor DCC
$R_{free}$ test set	1372 reflections (2.06%)	DCC
Wilson B-factor (Å <sup>2</sup> )	40.9	Xtriage
Anisotropy	0.051	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 30.9	EDS
Estimated twinning fraction	0.007 for -h,-l,-k	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 67920 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	9774	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.45% of the height of the origin peak. No significant pseudotranslation is detected.*

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

## 5 Model quality

### 5.1 Standard geometry

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

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	0.69	4/4724 (0.1%)	0.68	3/6413 (0.0%)
1	B	0.64	0/4725	0.70	6/6415 (0.1%)
All	All	0.67	4/9449 (0.0%)	0.69	9/12828 (0.1%)

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	A	0	1
1	B	0	2
All	All	0	3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	140	GLY	C-O	14.29	1.46	1.23
1	A	140	GLY	C-N	8.56	1.53	1.34
1	A	67	ASN	CG-ND2	7.71	1.52	1.32
1	A	67	ASN	CG-OD1	6.18	1.37	1.24

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	502	ARG	NE-CZ-NH1	9.11	124.85	120.30
1	B	584	GLU	N-CA-C	-7.61	90.45	111.00
1	B	502	ARG	NE-CZ-NH2	-7.04	116.78	120.30
1	A	502	ARG	NE-CZ-NH2	-6.53	117.03	120.30
1	A	98	ASP	C-N-CD	-5.47	108.57	120.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	502	ARG	NE-CZ-NH1	5.44	123.02	120.30
1	B	284	LEU	CA-CB-CG	-5.41	102.86	115.30
1	B	555	LEU	CA-CB-CG	5.22	127.31	115.30
1	B	126	GLU	C-N-CA	5.19	134.67	121.70

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	99	PRO	Peptide
1	B	127	ASN	Peptide
1	B	583	TYR	Peptide

## 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4628	0	4416	60	0
1	B	4629	0	4418	46	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	27	0	28	0	0
3	B	27	0	28	0	0
4	A	198	0	0	4	0
4	B	261	0	0	1	2
All	All	9774	0	8890	104	2

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 6.

All (104) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:99:PRO:CB	1:A:100:ASN:HB2	1.73	1.18
1:A:99:PRO:HB2	1:A:100:ASN:HB2	1.22	1.11
1:A:99:PRO:CB	1:A:100:ASN:CB	2.49	0.89

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:502:ARG:HD3	1:B:502:ARG:O	1.74	0.87
1:A:99:PRO:HB3	1:A:100:ASN:HB2	1.56	0.85
1:A:99:PRO:HB2	1:A:100:ASN:CB	2.04	0.82
1:A:515:LYS:HG2	1:A:515:LYS:O	1.77	0.81
1:A:99:PRO:HB3	1:A:100:ASN:CB	2.14	0.77
1:A:237:GLN:HG3	4:A:2046:HOH:O	1.84	0.76
1:B:47:ASN:HD21	1:B:446:ASP:HB2	1.53	0.72
1:B:359:PRO:HA	1:B:389:ARG:HH21	1.56	0.71
1:A:108:GLU:OE1	4:A:2009:HOH:O	2.08	0.69
1:A:41:VAL:HG13	1:A:59:GLY:HA3	1.78	0.65
1:A:502:ARG:O	1:A:502:ARG:HD3	1.97	0.64
1:A:99:PRO:HB3	1:A:100:ASN:CG	2.17	0.63
1:B:487:ASN:HD21	1:B:491:ALA:H	1.46	0.62
1:B:156:VAL:O	1:B:160:LYS:HG3	2.01	0.60
1:B:502:ARG:C	1:B:502:ARG:HD3	2.18	0.60
1:A:619:PHE:HD2	1:A:621:LEU:HD22	1.66	0.59
1:A:502:ARG:C	1:A:502:ARG:HD3	2.22	0.59
1:A:68:LEU:HG	1:A:71:GLU:HG3	1.82	0.59
1:A:619:PHE:CD2	1:A:621:LEU:CD2	2.86	0.58
1:B:47:ASN:ND2	1:B:446:ASP:HB2	2.20	0.57
1:A:99:PRO:HB3	1:A:100:ASN:OD1	2.05	0.56
1:A:80:ASN:O	1:A:84:GLU:HG3	2.04	0.56
1:A:107:GLY:O	1:A:145:GLU:HA	2.06	0.56
1:B:285:TYR:CZ	1:B:322:LYS:HD3	2.41	0.56
1:A:420:ASN:H	1:A:420:ASN:HD22	1.54	0.55
1:A:619:PHE:CD2	1:A:621:LEU:HD22	2.42	0.55
1:B:420:ASN:H	1:B:420:ASN:HD22	1.56	0.54
1:B:500:GLU:CD	1:B:500:GLU:H	2.11	0.54
1:A:538:ASN:OD1	1:A:542:ASN:ND2	2.41	0.54
1:A:52:ASN:O	1:A:173:ILE:HA	2.07	0.54
1:A:343:ASN:HB3	4:A:2076:HOH:O	2.08	0.54
1:A:387:TYR:HB3	1:A:389:ARG:HD2	1.89	0.53
1:B:368:PRO:HD2	1:B:372:THR:HG21	1.90	0.53
1:B:66:ILE:HG22	1:B:102:THR:HB	1.90	0.53
1:A:67:ASN:HB2	1:A:103:THR:HG23	1.91	0.53
1:A:42:LEU:H	1:A:42:LEU:HD12	1.72	0.53
1:B:285:TYR:CE1	1:B:322:LYS:HD3	2.44	0.52
1:B:126:GLU:N	1:B:127:ASN:HB3	2.25	0.52
1:A:504:LYS:HE2	1:A:524:GLU:OE1	2.10	0.52
1:B:126:GLU:N	1:B:127:ASN:CB	2.73	0.52
1:A:619:PHE:CD2	1:A:621:LEU:HD21	2.45	0.52
1:B:487:ASN:ND2	1:B:491:ALA:H	2.07	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:331:VAL:HG22	1:A:364:MET:HB2	1.92	0.51
1:A:392:ALA:HA	1:A:425:PHE:HB3	1.92	0.51
1:A:624:ILE:HD12	1:B:224:ARG:HH21	1.76	0.51
1:A:133:TYR:CE2	1:A:175:ASP:HB3	2.45	0.50
1:B:295:TYR:CD2	1:B:331:VAL:HG13	2.46	0.50
1:B:557:GLU:HG2	1:B:604:ILE:HG22	1.93	0.50
1:A:398:PRO:HD2	1:A:430:PRO:HA	1.93	0.50
1:A:65:SER:HB3	1:A:92:GLU:HB3	1.93	0.49
1:B:85:PHE:HB2	1:B:160:LYS:HD2	1.94	0.49
1:A:183:GLY:C	1:A:184:ILE:HD12	2.32	0.49
1:B:359:PRO:HA	1:B:389:ARG:NH2	2.23	0.49
1:B:183:GLY:O	1:B:427:THR:HA	2.13	0.49
1:B:364:MET:HA	1:B:392:ALA:O	2.12	0.48
1:B:315:ASN:O	1:B:320:LYS:HG3	2.13	0.48
1:A:183:GLY:O	1:A:427:THR:HA	2.13	0.48
1:A:295:TYR:CE2	1:A:331:VAL:HG23	2.48	0.48
1:B:295:TYR:CD2	1:B:331:VAL:CG1	2.98	0.47
1:B:387:TYR:HB3	1:B:389:ARG:HG2	1.97	0.46
1:B:387:TYR:HB3	1:B:389:ARG:HH11	1.80	0.46
1:A:85:PHE:CD1	1:A:160:LYS:HG2	2.51	0.46
1:B:154:TYR:CE1	1:B:209:LYS:HA	2.51	0.45
1:A:557:GLU:HG2	1:A:602:ALA:HB3	1.96	0.45
1:A:68:LEU:HG	1:A:71:GLU:CG	2.45	0.45
1:B:184:ILE:HD11	1:B:441:ILE:HG23	1.98	0.45
1:B:193:TRP:CZ3	1:B:201:GLN:HG3	2.51	0.45
1:A:505:MET:HG2	1:A:612:PHE:CD1	2.52	0.44
1:B:75:ASP:OD1	1:B:77:ASN:HB2	2.17	0.44
1:B:61:LYS:HE2	1:B:166:SER:HB2	2.00	0.44
1:A:99:PRO:HB2	1:A:100:ASN:CA	2.48	0.44
1:A:194:THR:O	1:A:197:ASP:HB2	2.17	0.44
1:A:98:ASP:HA	1:A:99:PRO:O	2.18	0.43
1:B:265:PHE:CD2	1:B:306:LYS:HD3	2.52	0.43
1:B:177:PRO:HB3	1:B:449:TRP:CE3	2.53	0.43
1:A:515:LYS:HG3	1:B:300:GLN:OE1	2.19	0.43
1:A:588:GLU:HG2	4:A:2048:HOH:O	2.17	0.43
1:B:337:THR:HB	1:B:368:PRO:HA	2.01	0.43
1:B:387:TYR:HD2	1:B:389:ARG:NH1	2.16	0.43
1:A:416:ASP:HB3	1:A:419:LEU:HD13	2.01	0.42
1:B:304:ALA:HB2	1:B:349:TYR:CD1	2.54	0.42
1:A:223:HIS:HD2	1:A:256:GLY:O	2.02	0.42
1:B:154:TYR:CD1	1:B:209:LYS:HG2	2.54	0.42
1:B:231:TYR:HB2	1:B:236:MET:SD	2.59	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:98:ASP:OD1	1:A:99:PRO:O	2.38	0.42
1:A:274:PHE:CE2	1:A:313:ARG:HD2	2.55	0.42
1:B:209:LYS:HD3	1:B:449:TRP:CH2	2.55	0.42
1:A:41:VAL:CG1	1:A:59:GLY:HA3	2.49	0.42
1:A:274:PHE:CZ	1:A:313:ARG:HD2	2.56	0.41
1:B:478:PHE:CE1	1:B:555:LEU:HD13	2.55	0.41
1:A:183:GLY:HA3	1:A:212:THR:O	2.21	0.41
1:A:337:THR:HB	1:A:368:PRO:HA	2.01	0.41
1:B:476:LYS:O	1:B:480:ASN:HB2	2.21	0.41
1:A:433:HIS:HB3	1:A:550:GLU:OE1	2.20	0.41
1:A:97:ASN:OD1	1:A:98:ASP:N	2.54	0.41
1:B:420:ASN:N	1:B:420:ASN:HD22	2.19	0.41
1:A:561:LEU:HD22	1:A:610:GLN:HA	2.04	0.40
1:A:114:PRO:O	1:A:118:GLU:HG2	2.21	0.40
1:A:165:GLU:O	1:A:166:SER:HB2	2.21	0.40
1:B:47:ASN:ND2	4:B:2004:HOH:O	2.54	0.40
1:B:557:GLU:CG	1:B:604:ILE:HG22	2.51	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	Distance(Å)	Clash(Å)
4:B:2241:HOH:O	4:B:2241:HOH:O[6_555]	1.97	0.23
4:B:2237:HOH:O	4:B:2241:HOH:O[6_555]	2.13	0.07

## 5.3 Torsion angles

### 5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	583/594 (98%)	559 (96%)	21 (4%)	3 (0%)	38	38
1	B	583/594 (98%)	558 (96%)	21 (4%)	4 (1%)	30	29
All	All	1166/1188 (98%)	1117 (96%)	42 (4%)	7 (1%)	33	33

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	99	PRO
1	B	128	LEU
1	B	127	ASN
1	B	388	ASP
1	A	98	ASP
1	A	343	ASN
1	B	41	VAL

### 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	500/507 (99%)	469 (94%)	31 (6%)	26	24
1	B	500/507 (99%)	475 (95%)	25 (5%)	34	36
All	All	1000/1014 (99%)	944 (94%)	56 (6%)	30	30

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	41	VAL
1	A	42	LEU
1	A	66	ILE
1	A	72	GLU
1	A	86	LEU
1	A	98	ASP
1	A	129	LYS
1	A	180	SER
1	A	237	GLN
1	A	257	ILE
1	A	284	LEU
1	A	313	ARG
1	A	328	LEU
1	A	355	GLU
1	A	358	ASP
1	A	365	TRP
1	A	372	THR
1	A	389	ARG
1	A	394	TRP

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Mol	Chain	Res	Type
1	A	409	LEU
1	A	420	ASN
1	A	482	SER
1	A	492	LYS
1	A	502	ARG
1	A	515	LYS
1	A	555	LEU
1	A	556	ASP
1	A	588	GLU
1	A	592	ASN
1	A	621	LEU
1	A	623	LEU
1	B	40	GLN
1	B	41	VAL
1	B	86	LEU
1	B	113	ILE
1	B	121	ASN
1	B	123	THR
1	B	128	LEU
1	B	174	THR
1	B	233	GLU
1	B	284	LEU
1	B	313	ARG
1	B	328	LEU
1	B	331	VAL
1	B	365	TRP
1	B	409	LEU
1	B	420	ASN
1	B	458	LYS
1	B	487	ASN
1	B	492	LYS
1	B	515	LYS
1	B	555	LEU
1	B	580	THR
1	B	581	GLU
1	B	619	PHE
1	B	624	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	121	ASN

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Mol	Chain	Res	Type
1	A	201	GLN
1	A	223	HIS
1	A	390	ASN
1	A	420	ASN
1	A	610	GLN
1	B	47	ASN
1	B	77	ASN
1	B	121	ASN
1	B	420	ASN
1	B	487	ASN
1	B	610	GLN
1	B	614	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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	GSZ	A	1626	-	29,29,29	1.51	4 (13%)	37,41,41	1.69	6 (16%)
3	GSZ	B	1627	-	29,29,29	1.47	5 (17%)	37,41,41	1.69	5 (13%)

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
3	GSZ	A	1626	-	-	0/15/35/35	0/1/3/3
3	GSZ	B	1627	-	-	0/15/35/35	0/1/3/3

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1626	GSZ	CAR-CAP	3.68	1.55	1.50
3	A	1626	GSZ	CAD-NAE	3.63	1.42	1.34
3	B	1627	GSZ	CAD-NAE	3.27	1.41	1.34
3	B	1627	GSZ	CAR-CAP	3.21	1.54	1.50
3	B	1627	GSZ	CAQ-NAE	-3.03	1.33	1.38
3	A	1626	GSZ	CAQ-NAE	-2.79	1.33	1.38
3	A	1626	GSZ	CAC-NAK	2.65	1.50	1.46
3	B	1627	GSZ	CAQ-CAP	-2.11	1.34	1.36
3	B	1627	GSZ	CAC-NAK	2.06	1.49	1.46

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1626	GSZ	CAC-CAD-NAO	5.57	130.80	119.48
3	B	1627	GSZ	CAC-CAD-NAO	5.44	130.54	119.48
3	B	1627	GSZ	CAQ-NAE-CAD	-4.98	106.21	109.50
3	A	1626	GSZ	CAQ-NAE-CAD	-4.97	106.22	109.50
3	B	1627	GSZ	CAR-CAP-CAQ	-3.05	124.60	129.50
3	B	1627	GSZ	CAN-CAL-NAK	-2.79	114.58	116.40
3	A	1626	GSZ	CAN-CAL-NAK	-2.79	114.58	116.40
3	A	1626	GSZ	CAR-CAP-CAQ	-2.56	125.39	129.50
3	A	1626	GSZ	CAB-CAA-CAF	-2.43	107.76	111.62
3	B	1627	GSZ	CAQ-CAP-NAO	2.24	108.75	106.15
3	A	1626	GSZ	CAQ-CAP-NAO	2.10	108.58	106.15

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

There are no chain breaks in this entry.



## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	585/594 (98%)	0.50	49 (8%)	11 13	35, 44, 54, 62	10 (1%)
1	B	585/594 (98%)	0.49	44 (7%)	14 17	32, 44, 55, 66	8 (1%)
All	All	1170/1188 (98%)	0.49	93 (7%)	13 15	32, 44, 54, 66	18 (1%)

All (93) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	41	VAL	12.1
1	B	122	GLY	5.7
1	A	112	ASP	5.6
1	A	100	ASN	5.2
1	B	100	ASN	5.1
1	B	99	PRO	5.0
1	A	57	GLY	4.8
1	A	58	ASP	4.6
1	A	624	ILE	4.6
1	B	121	ASN	4.4
1	A	93	ILE	4.3
1	A	101	SER	4.3
1	B	118	GLU	4.2
1	B	58	ASP	3.9
1	B	112	ASP	3.9
1	B	95	SER	3.7
1	A	98	ASP	3.6
1	A	110	ASP	3.6
1	A	40	GLN	3.5
1	B	93	ILE	3.4
1	A	138	ASN	3.4
1	A	69	VAL	3.4
1	A	409	LEU	3.4
1	A	139	ASP	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	409	LEU	3.3
1	B	56	VAL	3.3
1	B	113	ILE	3.3
1	A	95	SER	3.3
1	B	114	PRO	3.1
1	A	125	ALA	3.1
1	B	478	PHE	3.1
1	A	141	LYS	3.0
1	A	268	ASP	3.0
1	A	117	ASP	2.9
1	B	73	GLU	2.9
1	B	482	SER	2.9
1	B	147	LYS	2.9
1	B	269	ALA	2.9
1	A	118	GLU	2.8
1	A	232	PRO	2.8
1	B	110	ASP	2.8
1	A	478	PHE	2.7
1	B	127	ASN	2.7
1	A	407	LEU	2.7
1	A	517	ASP	2.7
1	B	139	ASP	2.7
1	B	138	ASN	2.7
1	A	73	GLU	2.6
1	B	90	ASN	2.6
1	B	483	THR	2.6
1	A	397	TYR	2.6
1	B	123	THR	2.5
1	A	114	PRO	2.5
1	B	561	LEU	2.5
1	B	479	ALA	2.5
1	B	126	GLU	2.4
1	B	616	ALA	2.4
1	A	222	TYR	2.4
1	A	127	ASN	2.4
1	A	623	LEU	2.4
1	A	122	GLY	2.4
1	B	604	ILE	2.4
1	B	344	GLY	2.4
1	A	111	ASP	2.3
1	A	41	VAL	2.3
1	B	170	GLU	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	408	ALA	2.3
1	B	54	GLU	2.3
1	B	141	LYS	2.3
1	B	128	LEU	2.3
1	A	126	GLU	2.3
1	A	609	ALA	2.3
1	A	398	PRO	2.2
1	A	444	ALA	2.2
1	A	142	ILE	2.2
1	B	431	MET	2.2
1	A	79	VAL	2.2
1	A	604	ILE	2.2
1	A	97	ASN	2.2
1	A	56	VAL	2.2
1	B	61	LYS	2.1
1	A	430	PRO	2.1
1	A	121	ASN	2.1
1	B	116	LEU	2.1
1	B	92	GLU	2.1
1	B	343	ASN	2.1
1	A	479	ALA	2.1
1	B	436	LEU	2.1
1	A	102	THR	2.1
1	A	105	ILE	2.1
1	A	119	ALA	2.1
1	B	605	SER	2.0
1	B	40	GLN	2.0

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

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

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	CL	B	1626	1/1	0.23	5.82	67,67,67,67	0
3	GSZ	B	1627	27/27	0.20	1.17	37,44,62,64	0
3	GSZ	A	1626	27/27	0.16	0.47	37,42,53,54	0
2	CL	A	1627	1/1	0.13	-1.03	67,67,67,67	0
2	CL	B	1625	1/1	0.08	-4.25	67,67,67,67	0
2	CL	A	1625	1/1	0.05	-7.19	75,75,75,75	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.