



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:39 PM GMT

PDB ID : 1GOI
Title : CRYSTAL STRUCTURE OF THE D140N MUTANT OF CHITINASE B
FROM SERRATIA MARCESCENS AT 1.45 Å RESOLUTION
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Deposited on : 2001-10-21
Resolution : 1.45 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

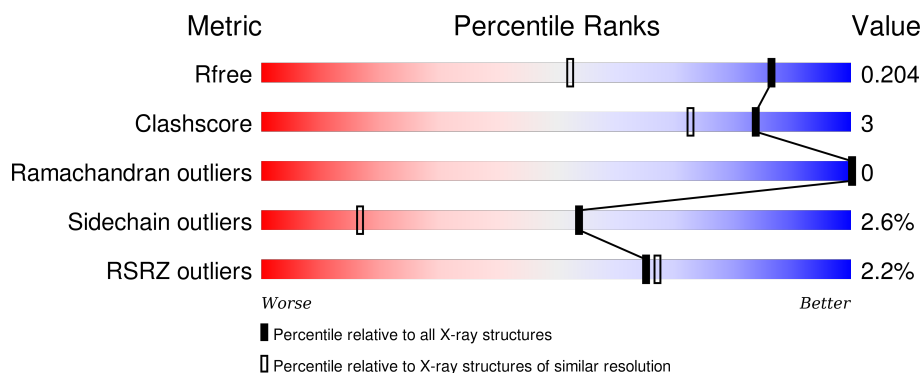
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.45 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	1278 (1.48-1.44)
Clashscore	102246	1336 (1.48-1.44)
Ramachandran outliers	100387	1320 (1.48-1.44)
Sidechain outliers	100360	1320 (1.48-1.44)
RSRZ outliers	91569	1279 (1.48-1.44)

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	499	<div> <div>3%</div> <div>84%</div> <div>14%</div> </div>
1	B	499	<div> <div>2%</div> <div>85%</div> <div>14%</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
2	GOL	A	1500	-	-	-	X
2	GOL	A	1501	-	-	-	X
2	GOL	A	1502	-	-	-	X
2	GOL	A	1503	-	-	-	X
2	GOL	A	1504	-	-	-	X
2	GOL	B	1500	-	-	-	X
2	GOL	B	1504	-	-	-	X
2	GOL	B	1505	-	-	-	X
3	SO4	B	1506	-	-	-	X
3	SO4	B	1508	-	-	-	X

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 8990 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 CHITINASE B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	497	Total	C	N	O	S	0	10	1
			3926	2511	660	741	14			
1	B	497	Total	C	N	O	S	0	14	0
			3972	2536	677	745	14			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	140	ASN	ASP	ENGINEERED MUTATION	UNP Q54276
B	140	ASN	ASP	ENGINEERED MUTATION	UNP Q54276

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



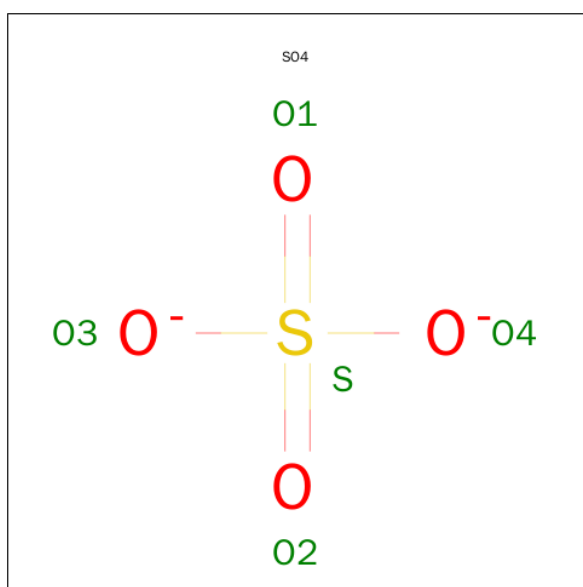
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		

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

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

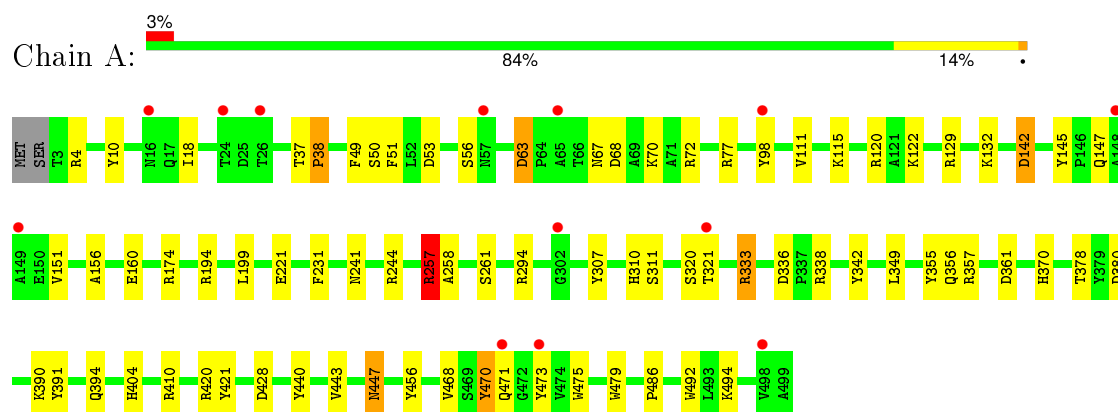
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	446	Total O 446 446	0	0
4	B	554	Total O 554 554	0	0

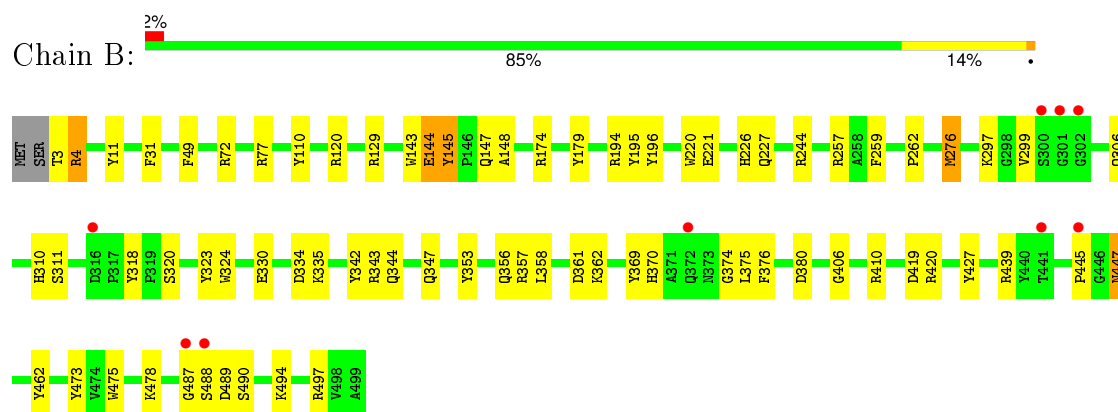
3 Residue-property plots

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: CHITINASE B



• Molecule 1: CHITINASE B



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	55.90Å 104.02Å 186.52Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	17.00 – 1.45 16.96 – 1.45	Depositor EDS
% Data completeness (in resolution range)	99.3 (17.00-1.45) 93.9 (16.96-1.45)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.15 (at 1.45Å)	Xtriage
Refinement program	SHELXH-97	Depositor
R, R_{free}	0.160 , 0.220 0.157 , 0.204	Depositor DCC
R_{free} test set	1921 reflections (1.07%)	DCC
Wilson B-factor (Å ²)	16.1	Xtriage
Anisotropy	0.293	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 64.2	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	0 of 191353 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	8990	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.17% of the height of the origin peak. No significant pseudotranslation is detected.*

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: GOL, SO4

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	0/4076	1.51	51/5558 (0.9%)
1	B	0.70	0/4142	1.47	56/5638 (1.0%)
All	All	0.70	0/8218	1.49	107/11196 (1.0%)

There are no bond length outliers.

All (107) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	257	ARG	NE-CZ-NH2	17.60	129.10	120.30
1	B	343	ARG	NE-CZ-NH1	-13.79	113.40	120.30
1	A	257	ARG	NE-CZ-NH2	-13.00	113.80	120.30
1	A	194	ARG	NE-CZ-NH2	12.64	126.62	120.30
1	A	447	ASN	CB-CG-OD1	11.42	144.44	121.60
1	B	497	ARG	NE-CZ-NH2	-10.77	114.92	120.30
1	A	72	ARG	NE-CZ-NH1	10.76	125.68	120.30
1	A	4	ARG	NE-CZ-NH2	9.96	125.28	120.30
1	B	194	ARG	NE-CZ-NH2	-9.62	115.49	120.30
1	A	72	ARG	NE-CZ-NH2	-9.59	115.50	120.30
1	A	120	ARG	NE-CZ-NH1	9.36	124.98	120.30
1	B	410	ARG	NE-CZ-NH2	-9.21	115.70	120.30
1	A	311	SER	C-N-CA	9.04	144.30	121.70
1	A	77	ARG	NE-CZ-NH2	-8.92	115.84	120.30
1	A	443	VAL	C-N-CA	8.92	141.03	122.30
1	A	333	ARG	NE-CZ-NH1	8.66	124.63	120.30
1	B	497	ARG	NE-CZ-NH1	8.66	124.63	120.30
1	A	410	ARG	NE-CZ-NH1	8.57	124.59	120.30
1	A	145	TYR	CB-CG-CD2	8.55	126.13	121.00
1	A	174	ARG	NE-CZ-NH2	-8.40	116.10	120.30
1	B	439	ARG	NE-CZ-NH1	-8.20	116.20	120.30
1	B	174	ARG	NE-CZ-NH1	8.20	124.40	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	129	ARG	CD-NE-CZ	7.93	134.70	123.60
1	A	447	ASN	OD1-CG-ND2	-7.92	103.68	121.90
1	A	244	ARG	NE-CZ-NH1	-7.88	116.36	120.30
1	B	194	ARG	NE-CZ-NH1	7.85	124.22	120.30
1	A	311	SER	O-C-N	-7.78	110.25	122.70
1	B	330	GLU	OE1-CD-OE2	7.74	132.59	123.30
1	A	142	ASP	CB-CG-OD2	-7.72	111.35	118.30
1	B	276	MET	CA-CB-CG	7.55	126.14	113.30
1	A	470	TYR	CB-CG-CD2	7.34	125.40	121.00
1	B	77	ARG	NE-CZ-NH1	7.25	123.92	120.30
1	A	129	ARG	NE-CZ-NH2	-7.24	116.68	120.30
1	B	72	ARG	NE-CZ-NH1	7.22	123.91	120.30
1	A	63	ASP	CB-CG-OD1	7.20	124.78	118.30
1	B	420[A]	ARG	CD-NE-CZ	6.82	133.15	123.60
1	B	420[B]	ARG	CD-NE-CZ	6.82	133.15	123.60
1	B	120	ARG	NE-CZ-NH1	6.81	123.70	120.30
1	B	120	ARG	NE-CZ-NH2	-6.61	117.00	120.30
1	A	456	TYR	CB-CG-CD2	6.58	124.94	121.00
1	B	342	TYR	CB-CG-CD1	-6.48	117.11	121.00
1	A	370	HIS	CG-ND1-CE1	6.48	117.27	108.20
1	B	257	ARG	NE-CZ-NH1	-6.43	117.08	120.30
1	A	456	TYR	CB-CG-CD1	-6.35	117.19	121.00
1	B	110	TYR	CB-CG-CD2	6.33	124.80	121.00
1	A	420	ARG	NE-CZ-NH1	6.26	123.43	120.30
1	A	142	ASP	CB-CG-OD1	6.17	123.85	118.30
1	B	77	ARG	CG-CD-NE	-6.15	98.88	111.80
1	B	129[A]	ARG	NE-CZ-NH1	6.15	123.38	120.30
1	B	129[B]	ARG	NE-CZ-NH1	6.15	123.38	120.30
1	B	462	TYR	CG-CD1-CE1	6.13	126.21	121.30
1	A	355	TYR	CB-CG-CD2	-6.11	117.33	121.00
1	A	294	ARG	NE-CZ-NH2	-6.10	117.25	120.30
1	A	194	ARG	CD-NE-CZ	6.08	132.11	123.60
1	A	391	TYR	CB-CG-CD2	6.05	124.63	121.00
1	B	244[A]	ARG	NE-CZ-NH2	-5.97	117.32	120.30
1	B	244[B]	ARG	NE-CZ-NH2	-5.97	117.32	120.30
1	B	488	SER	N-CA-CB	5.94	119.41	110.50
1	B	174	ARG	NE-CZ-NH2	-5.92	117.34	120.30
1	A	77	ARG	NE-CZ-NH1	5.89	123.25	120.30
1	A	404	HIS	CG-ND1-CE1	5.88	116.43	108.20
1	B	369	TYR	CB-CG-CD1	5.88	124.53	121.00
1	A	428	ASP	CB-CG-OD1	5.86	123.57	118.30
1	B	357	ARG	NE-CZ-NH2	-5.85	117.38	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	4	ARG	NE-CZ-NH1	-5.84	117.38	120.30
1	A	336	ASP	CB-CG-OD2	-5.82	113.07	118.30
1	A	49	PHE	CB-CG-CD1	-5.75	116.77	120.80
1	A	231	PHE	CB-CG-CD2	-5.75	116.78	120.80
1	B	143	TRP	CG-CD1-NE1	5.70	115.80	110.10
1	A	342	TYR	CB-CG-CD2	5.70	124.42	121.00
1	B	49	PHE	CB-CG-CD2	5.61	124.72	120.80
1	B	427	TYR	CB-CG-CD1	-5.59	117.64	121.00
1	B	144	GLU	C-N-CA	5.57	135.62	121.70
1	B	318	TYR	CB-CG-CD2	-5.57	117.66	121.00
1	A	244	ARG	NH1-CZ-NH2	5.53	125.48	119.40
1	B	361	ASP	CB-CG-OD2	5.52	123.27	118.30
1	B	344	GLN	CG-CD-OE1	5.48	132.57	121.60
1	A	68	ASP	CB-CG-OD2	-5.48	113.37	118.30
1	A	440	TYR	CB-CG-CD2	5.47	124.28	121.00
1	A	257	ARG	NH1-CZ-NH2	5.46	125.41	119.40
1	B	179	TYR	CG-CD2-CE2	-5.41	116.98	121.30
1	B	195	TYR	CB-CG-CD1	5.39	124.24	121.00
1	B	311	SER	C-N-CA	5.38	135.15	121.70
1	B	488	SER	O-C-N	5.38	131.31	122.70
1	A	361	ASP	CB-CG-OD1	5.37	123.13	118.30
1	B	343	ARG	NE-CZ-NH2	5.33	122.96	120.30
1	B	145	TYR	CB-CG-CD2	5.33	124.20	121.00
1	B	323	TYR	CB-CG-CD2	5.32	124.19	121.00
1	B	196	TYR	CB-CG-CD2	5.31	124.18	121.00
1	B	489	ASP	CB-CG-OD1	5.30	123.07	118.30
1	B	179	TYR	CZ-CE2-CD2	5.30	124.57	119.80
1	B	244[A]	ARG	NE-CZ-NH1	5.29	122.94	120.30
1	B	244[B]	ARG	NE-CZ-NH1	5.29	122.94	120.30
1	B	376	PHE	CB-CG-CD1	5.28	124.49	120.80
1	B	419	ASP	CB-CG-OD2	-5.21	113.61	118.30
1	A	98	TYR	CB-CG-CD1	-5.16	117.91	121.00
1	B	4	ARG	NE-CZ-NH2	-5.15	117.72	120.30
1	B	11	TYR	CB-CG-CD1	5.15	124.09	121.00
1	A	421	TYR	CG-CD1-CE1	5.14	125.41	121.30
1	A	307	TYR	CB-CG-CD1	5.08	124.05	121.00
1	A	174	ARG	NE-CZ-NH1	5.08	122.84	120.30
1	B	257	ARG	NH1-CZ-NH2	-5.08	113.82	119.40
1	A	338	ARG	NE-CZ-NH1	-5.07	117.77	120.30
1	A	53	ASP	CB-CG-OD1	5.06	122.86	118.30
1	B	143	TRP	CD1-CG-CD2	-5.04	102.27	106.30
1	A	357	ARG	NE-CZ-NH1	-5.04	117.78	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	347	GLN	CB-CG-CD	5.01	124.63	111.60

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	3926	0	3747	25	0
1	B	3972	0	3794	21	0
2	A	36	0	48	5	0
2	B	36	0	47	3	0
3	A	5	0	0	0	0
3	B	15	0	0	0	0
4	A	446	0	0	4	0
4	B	554	0	0	5	0
All	All	8990	0	7636	47	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 3.

All (47) 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:479:TRP:HE1	2:B:1505:GOL:H2	1.42	0.85
1:A:261:SER:H	2:A:1501:GOL:H31	1.41	0.84
2:B:1500:GOL:H31	4:B:2536:HOH:O	1.89	0.72
1:A:111:VAL:O	1:A:115:LYS:HG3	1.95	0.66
1:B:4:ARG:HD3	2:B:1504:GOL:O2	1.95	0.66
1:A:261:SER:N	2:A:1501:GOL:H31	2.10	0.63
1:B:356[A]:GLN:NE2	1:B:358:LEU:HD21	2.17	0.58
1:A:221:GLU:O	1:A:310:HIS:HE1	1.86	0.58
1:B:221:GLU:O	1:B:310:HIS:HE1	1.87	0.57
1:B:147[B]:GLN:HG3	1:B:148:ALA:N	2.20	0.57
1:B:475:TRP:CZ3	1:B:487:GLY:HA3	2.41	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:356[B]:GLN:HG3	4:B:2401:HOH:O	2.05	0.55
1:B:226:HIS:HD2	4:B:2418:HOH:O	1.88	0.55
1:B:478:LYS:HE3	1:B:490:SER:O	2.08	0.54
1:B:226:HIS:HE1	1:B:306:GLN:OE1	1.90	0.54
1:B:297:LYS:HD3	1:B:324:TRP:CE2	2.43	0.53
1:A:122[B]:LYS:HE2	4:A:2114:HOH:O	2.11	0.51
1:B:445:PRO:HD3	4:B:2487:HOH:O	2.11	0.50
1:B:353:TYR:O	1:B:370:HIS:HE1	1.95	0.49
1:B:362[B]:LYS:HD2	4:B:2309:HOH:O	2.12	0.49
1:A:122[B]:LYS:HB3	1:A:122[B]:LYS:HE2	1.53	0.48
1:B:276:MET:O	1:B:447:ASN:ND2	2.45	0.48
1:A:37[B]:THR:OG1	1:A:38:PRO:HD2	2.13	0.48
1:A:142:ASP:OD2	2:A:1504:GOL:O1	2.29	0.47
1:A:18:ILE:O	1:A:70:LYS:HE3	2.15	0.46
1:B:299:VAL:HG22	1:B:374:GLY:C	2.36	0.46
1:A:475:TRP:CD2	1:A:486:PRO:HB3	2.51	0.45
1:A:147:GLN:O	1:A:151:VAL:HG23	2.16	0.45
1:A:51:PHE:HZ	2:A:1504:GOL:O2	1.99	0.45
1:A:10:TYR:HD2	1:A:50:SER:HG	1.61	0.45
1:A:390:LYS:O	1:A:394:GLN:HG3	2.16	0.45
1:A:473:TYR:CD1	1:A:494:LYS:HD3	2.52	0.45
1:A:257:ARG:NH2	4:A:2235:HOH:O	2.50	0.44
1:A:356:GLN:HG3	4:A:2309:HOH:O	2.18	0.44
1:B:220:TRP:CD1	1:B:221:GLU:HG2	2.53	0.43
1:A:67:ASN:OD1	1:A:70:LYS:HG2	2.19	0.43
1:A:470:TYR:CD2	1:A:471:GLN:HG3	2.54	0.43
1:B:144:GLU:HA	1:B:145:TYR:CG	2.54	0.42
1:A:156:ALA:O	1:A:160:GLU:HG2	2.19	0.42
1:B:31:PHE:CG	1:B:406:GLY:HA2	2.54	0.42
1:B:473:TYR:CE2	1:B:494:LYS:HD3	2.54	0.42
1:A:257:ARG:HG3	1:A:258:ALA:N	2.33	0.42
1:B:334:ASP:O	1:B:335:LYS:HB2	2.20	0.42
1:B:259:PHE:CE1	1:B:276:MET:HE1	2.56	0.41
1:A:468:VAL:HG22	1:A:492:TRP:CZ3	2.56	0.41
1:A:356:GLN:NE2	4:A:2309:HOH:O	2.51	0.41
1:A:261:SER:H	2:A:1501:GOL:C3	2.23	0.41

There are no symmetry-related clashes.

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	505/499 (101%)	493 (98%)	12 (2%)	0	100	100
1	B	509/499 (102%)	502 (99%)	7 (1%)	0	100	100
All	All	1014/998 (102%)	995 (98%)	19 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	411/405 (102%)	395 (96%)	16 (4%)	39	7
1	B	414/405 (102%)	407 (98%)	7 (2%)	68	32
All	All	825/810 (102%)	802 (97%)	23 (3%)	54	14

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	PRO
1	A	56[A]	SER
1	A	56[B]	SER
1	A	63	ASP
1	A	132	LYS
1	A	199	LEU
1	A	241	ASN
1	A	257	ARG

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Mol	Chain	Res	Type
1	A	320	SER
1	A	321[A]	THR
1	A	321[B]	THR
1	A	333	ARG
1	A	349	LEU
1	A	378	THR
1	A	380	ASP
1	A	447	ASN
1	B	3	THR
1	B	227	GLN
1	B	262	PRO
1	B	320	SER
1	B	375	LEU
1	B	380	ASP
1	B	447	ASN

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	147	GLN
1	A	241	ASN
1	A	310	HIS
1	A	447	ASN
1	B	112	ASN
1	B	180	GLN
1	B	226	HIS
1	B	310	HIS
1	B	344	GLN
1	B	370	HIS
1	B	447	ASN

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

16 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	GOL	A	1499	-	5,5,5	0.52	0	5,5,5	2.02	1 (20%)
2	GOL	A	1500	-	5,5,5	0.53	0	5,5,5	3.24	4 (80%)
2	GOL	A	1501	-	5,5,5	0.36	0	5,5,5	4.45	4 (80%)
2	GOL	A	1502	-	5,5,5	0.71	0	5,5,5	3.51	3 (60%)
2	GOL	A	1503	-	5,5,5	0.29	0	5,5,5	3.34	3 (60%)
2	GOL	A	1504	-	5,5,5	0.31	0	5,5,5	4.03	3 (60%)
3	SO4	A	1505	-	4,4,4	0.69	0	6,6,6	0.52	0
2	GOL	B	1500	-	5,5,5	0.87	0	5,5,5	5.39	3 (60%)
2	GOL	B	1501	-	5,5,5	0.41	0	5,5,5	3.94	2 (40%)
2	GOL	B	1502	-	5,5,5	0.65	0	5,5,5	2.95	3 (60%)
2	GOL	B	1503	-	5,5,5	0.62	0	5,5,5	3.98	4 (80%)
2	GOL	B	1504	-	5,5,5	0.54	0	5,5,5	3.60	3 (60%)
2	GOL	B	1505	-	5,5,5	0.28	0	5,5,5	4.29	4 (80%)
3	SO4	B	1506	-	4,4,4	0.89	0	6,6,6	0.59	0
3	SO4	B	1507	-	4,4,4	0.71	0	6,6,6	0.35	0
3	SO4	B	1508	-	4,4,4	0.78	0	6,6,6	0.26	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	GOL	A	1499	-	-	0/4/4/4	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	A	1500	-	-	0/4/4/4	0/0/0/0
2	GOL	A	1501	-	-	0/4/4/4	0/0/0/0
2	GOL	A	1502	-	-	0/4/4/4	0/0/0/0
2	GOL	A	1503	-	-	0/4/4/4	0/0/0/0
2	GOL	A	1504	-	-	0/4/4/4	0/0/0/0
3	SO4	A	1505	-	-	0/0/0/0	0/0/0/0
2	GOL	B	1500	-	-	0/4/4/4	0/0/0/0
2	GOL	B	1501	-	-	0/4/4/4	0/0/0/0
2	GOL	B	1502	-	-	0/4/4/4	0/0/0/0
2	GOL	B	1503	-	-	0/4/4/4	0/0/0/0
2	GOL	B	1504	-	-	0/4/4/4	0/0/0/0
2	GOL	B	1505	-	-	0/4/4/4	0/0/0/0
3	SO4	B	1506	-	-	0/0/0/0	0/0/0/0
3	SO4	B	1507	-	-	0/0/0/0	0/0/0/0
3	SO4	B	1508	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1500	GOL	C3-C2-C1	-4.56	93.25	111.12
2	A	1502	GOL	C3-C2-C1	-4.17	94.75	111.12
2	A	1504	GOL	C3-C2-C1	-3.64	96.86	111.12
2	B	1502	GOL	C3-C2-C1	-3.58	97.10	111.12
2	A	1500	GOL	C3-C2-C1	-3.50	97.38	111.12
2	B	1503	GOL	C3-C2-C1	-3.25	98.39	111.12
2	B	1504	GOL	C3-C2-C1	-2.83	100.01	111.12
2	B	1505	GOL	C3-C2-C1	-2.75	100.33	111.12
2	A	1503	GOL	C3-C2-C1	-2.72	100.44	111.12
2	B	1503	GOL	O2-C2-C1	-2.49	97.24	108.65
2	A	1501	GOL	C3-C2-C1	-2.45	101.52	111.12
2	B	1500	GOL	O2-C2-C3	2.23	118.88	108.65
2	A	1500	GOL	O1-C1-C2	2.36	121.65	110.18
2	B	1503	GOL	O2-C2-C3	2.67	120.90	108.65
2	A	1501	GOL	O2-C2-C3	2.77	121.33	108.65
2	B	1502	GOL	O2-C2-C3	3.03	122.56	108.65
2	A	1500	GOL	O2-C2-C3	3.17	123.16	108.65
2	A	1503	GOL	O2-C2-C3	3.47	124.56	108.65
2	B	1505	GOL	O2-C2-C3	3.68	125.54	108.65
2	A	1499	GOL	O3-C3-C2	3.69	128.08	110.18
2	B	1505	GOL	O1-C1-C2	3.87	128.95	110.18
2	A	1501	GOL	O1-C1-C2	4.07	129.90	110.18

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1502	GOL	O3-C3-C2	4.11	130.12	110.18
2	A	1502	GOL	O2-C2-C3	4.26	128.17	108.65
2	B	1504	GOL	O2-C2-C3	4.39	128.80	108.65
2	A	1500	GOL	O3-C3-C2	4.93	134.09	110.18
2	A	1502	GOL	O3-C3-C2	5.08	134.80	110.18
2	A	1504	GOL	O2-C2-C3	5.22	132.60	108.65
2	B	1501	GOL	O2-C2-C3	5.35	133.19	108.65
2	A	1503	GOL	O3-C3-C2	5.86	138.62	110.18
2	B	1504	GOL	O3-C3-C2	5.98	139.20	110.18
2	A	1504	GOL	O3-C3-C2	6.16	140.06	110.18
2	B	1501	GOL	O3-C3-C2	6.77	143.04	110.18
2	B	1503	GOL	O3-C3-C2	7.45	146.29	110.18
2	B	1505	GOL	O3-C3-C2	7.48	146.48	110.18
2	A	1501	GOL	O3-C3-C2	8.30	150.45	110.18
2	B	1500	GOL	O3-C3-C2	10.77	162.39	110.18

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1501	GOL	3	0
2	A	1504	GOL	2	0
2	B	1500	GOL	1	0
2	B	1504	GOL	1	0
2	B	1505	GOL	1	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, 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	497/499 (99%)	-0.16	13 (2%) 59 60	11, 22, 38, 48	0
1	B	497/499 (99%)	-0.31	9 (1%) 71 72	13, 19, 31, 47	0
All	All	994/998 (99%)	-0.24	22 (2%) 65 67	11, 20, 36, 48	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	488	SER	5.4
1	B	301	GLY	4.9
1	A	57	ASN	3.8
1	A	24	THR	3.3
1	B	302	GLY	3.1
1	B	487	GLY	2.9
1	A	149	ALA	2.9
1	A	321[A]	THR	2.8
1	B	441	THR	2.6
1	A	16	ASN	2.6
1	A	98	TYR	2.4
1	B	316	ASP	2.3
1	A	302	GLY	2.3
1	B	445	PRO	2.3
1	B	300	SER	2.3
1	A	471	GLN	2.2
1	A	473	TYR	2.2
1	A	26[A]	THR	2.1
1	A	65	ALA	2.1
1	A	498	VAL	2.1
1	B	372	GLN	2.1
1	A	148	ALA	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 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, 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
2	GOL	A	1504	6/6	0.65	0.28	15.18	61,65,67,68	0
2	GOL	A	1503	6/6	0.73	0.19	11.81	50,54,54,59	0
2	GOL	B	1500	6/6	0.82	0.24	9.87	55,58,59,59	0
2	GOL	B	1505	6/6	0.80	0.32	8.81	69,69,70,72	0
2	GOL	A	1501	6/6	0.69	0.22	8.07	38,45,49,53	0
2	GOL	A	1502	6/6	0.86	0.13	7.48	38,40,43,51	0
3	SO4	B	1506	5/5	0.97	0.17	4.92	35,36,37,40	0
3	SO4	B	1508	5/5	0.97	0.17	3.57	61,66,67,67	0
2	GOL	B	1504	6/6	0.87	0.16	3.55	23,36,38,43	0
2	GOL	A	1500	6/6	0.91	0.12	3.05	22,31,34,35	0
2	GOL	B	1501	6/6	0.90	0.12	1.96	31,37,39,40	0
3	SO4	A	1505	5/5	0.98	0.09	1.78	36,39,43,45	0
2	GOL	A	1499	6/6	0.92	0.08	1.28	31,33,35,40	0
2	GOL	B	1502	6/6	0.90	0.12	0.89	23,33,37,40	0
2	GOL	B	1503	6/6	0.88	0.10	0.19	30,35,38,39	0
3	SO4	B	1507	5/5	0.96	0.21	-	48,52,54,57	0

6.5 Other polymers [i](#)

There are no such residues in this entry.