



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:21 PM GMT

PDB ID : 4EK7
Title : High speed X-ray analysis of plant enzymes at room temperature
Authors : Xia, L.; Rajendran, C.; Ruppert, M.; Panjikar, S.; Wang, M.; Stoeckigt, J.
Deposited on : 2012-04-09
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
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

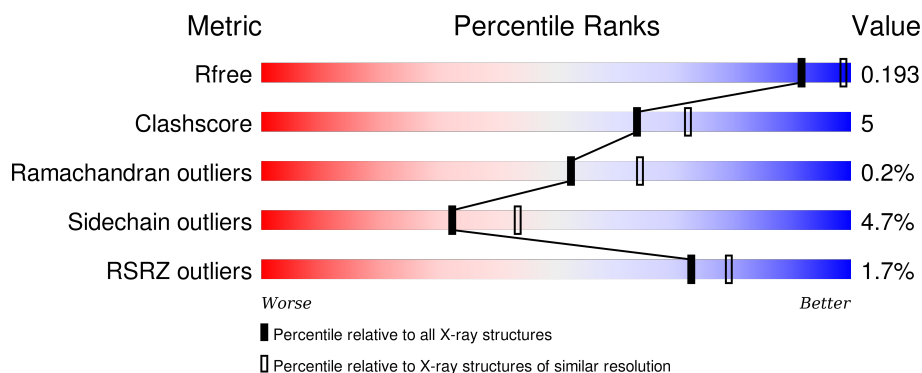
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

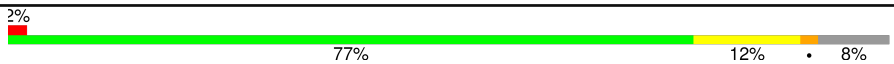

The reported resolution of this entry is 2.30 Å.

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 ≥ 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	513	 2% 77% 12% • 8%
1	B	513	 % 76% 14% • 8%

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	BGC	A	601	-	-	-	X
3	CL	B	601	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7792 atoms, of which 12 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.

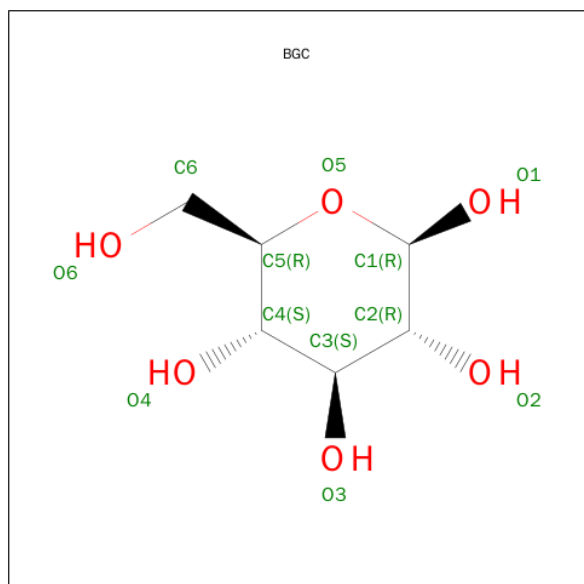
- Molecule 1 is a protein called Raucaffricine-O-beta-D-glucosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	470	Total	C	N	O	S	0	0	0
			3766	2410	640	703	13			
1	B	470	Total	C	N	O	S	0	0	0
			3766	2410	640	703	13			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	186	GLN	GLU	ENGINEERED MUTATION	UNP Q9SPP9
B	186	GLN	GLU	ENGINEERED MUTATION	UNP Q9SPP9

- Molecule 2 is SUGAR (BETA-D-GLUCOSE) (three-letter code: BGC) (formula: C₆H₁₂O₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	H	O	0	0
			24	6	12	6		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	C	O	0	0
			12	6	6		

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

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

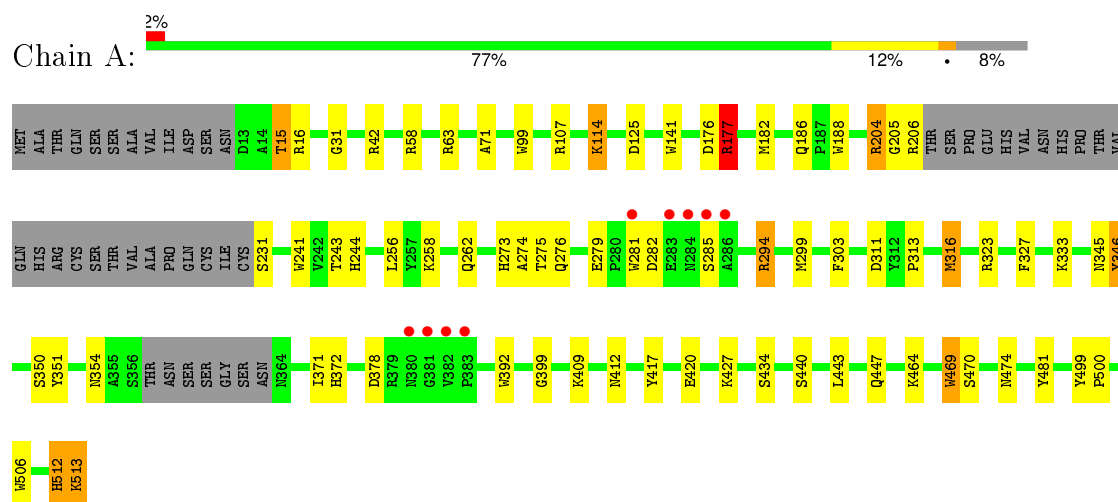
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	109	Total	O	0	0
			109	109		
4	B	113	Total	O	0	0
			113	113		

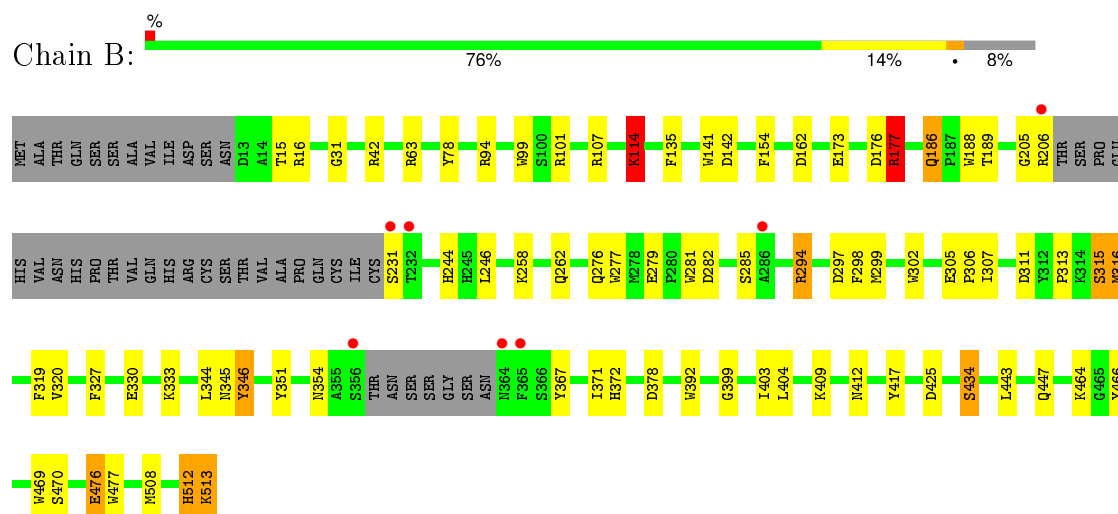
3 Residue-property plots

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 ($\text{RSRZ} > 2$). 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: Raucaffricine-O-beta-D-glucosidase



• Molecule 1: Raucaffricine-O-beta-D-glucosidase



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	105.56Å 129.52Å 216.64Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.75 – 2.30 47.68 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.3 (47.75-2.30) 99.3 (47.68-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.91 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.151 , 0.192 0.156 , 0.193	Depositor DCC
R_{free} test set	3282 reflections (5.24%)	DCC
Wilson B-factor (Å ²)	41.4	Xtriage
Anisotropy	0.056	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 41.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 66091 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	7792	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

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

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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: BGC, 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	1.18	15/3877 (0.4%)	1.05	14/5263 (0.3%)
1	B	1.17	17/3877 (0.4%)	1.04	12/5263 (0.2%)
All	All	1.17	32/7754 (0.4%)	1.04	26/10526 (0.2%)

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	1
All	All	0	2

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	188	TRP	CD2-CE2	7.87	1.50	1.41
1	A	417	TYR	CG-CD1	-7.81	1.28	1.39
1	A	417	TYR	CE1-CZ	-7.21	1.29	1.38
1	A	204	ARG	CZ-NH1	-6.62	1.24	1.33
1	B	346	TYR	CE1-CZ	-6.47	1.30	1.38
1	A	42	ARG	CZ-NH1	6.29	1.41	1.33
1	A	346	TYR	CE1-CZ	-6.26	1.30	1.38
1	B	78	TYR	CE1-CZ	-6.08	1.30	1.38
1	A	141	TRP	CD2-CE2	6.06	1.48	1.41
1	B	141	TRP	CD2-CE2	6.05	1.48	1.41
1	B	466	TYR	CE1-CZ	-5.99	1.30	1.38
1	B	281	TRP	CD2-CE2	5.94	1.48	1.41
1	B	417	TYR	CG-CD1	-5.93	1.31	1.39
1	A	99	TRP	CD2-CE2	5.90	1.48	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	281	TRP	CD2-CE2	5.77	1.48	1.41
1	B	476	GLU	CD-OE1	5.72	1.31	1.25
1	A	141	TRP	CG-CD1	5.66	1.44	1.36
1	B	99	TRP	CD2-CE2	5.57	1.48	1.41
1	B	477	TRP	CD2-CE2	5.56	1.48	1.41
1	A	506	TRP	CD2-CE2	5.56	1.48	1.41
1	B	392	TRP	CD2-CE2	5.51	1.48	1.41
1	B	188	TRP	CD2-CE2	5.51	1.48	1.41
1	A	469	TRP	CD2-CE2	5.45	1.47	1.41
1	B	173	GLU	CG-CD	5.32	1.59	1.51
1	B	277	TRP	CD2-CE2	5.31	1.47	1.41
1	B	302	TRP	CD2-CE2	5.24	1.47	1.41
1	B	417	TYR	CE2-CZ	-5.22	1.31	1.38
1	A	241	TRP	CD2-CE2	5.21	1.47	1.41
1	A	392	TRP	CD2-CE2	5.15	1.47	1.41
1	A	350	SER	CB-OG	-5.15	1.35	1.42
1	B	466	TYR	CG-CD2	-5.13	1.32	1.39
1	B	42	ARG	CZ-NH1	5.08	1.39	1.33

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	42	ARG	NE-CZ-NH2	-10.54	115.03	120.30
1	B	42	ARG	NE-CZ-NH2	-9.23	115.68	120.30
1	A	177	ARG	NE-CZ-NH1	9.14	124.87	120.30
1	A	177	ARG	NE-CZ-NH2	-9.13	115.73	120.30
1	B	294	ARG	NE-CZ-NH2	-8.06	116.27	120.30
1	A	42	ARG	NE-CZ-NH1	7.77	124.19	120.30
1	A	294	ARG	NE-CZ-NH2	-7.14	116.73	120.30
1	B	42	ARG	NE-CZ-NH1	6.66	123.63	120.30
1	B	316	MET	CG-SD-CE	-6.60	89.64	100.20
1	B	311	ASP	CB-CG-OD1	6.24	123.92	118.30
1	B	425	ASP	CB-CG-OD1	6.17	123.86	118.30
1	A	58	ARG	NE-CZ-NH1	5.82	123.21	120.30
1	A	443	LEU	CB-CG-CD2	-5.80	101.13	111.00
1	B	101	ARG	NE-CZ-NH1	5.78	123.19	120.30
1	B	177	ARG	NE-CZ-NH1	5.76	123.18	120.30
1	B	162	ASP	CB-CG-OD2	5.68	123.41	118.30
1	B	177	ARG	NE-CZ-NH2	-5.62	117.49	120.30
1	A	282	ASP	CB-CG-OD2	-5.45	113.40	118.30
1	A	311	ASP	CB-CG-OD1	5.40	123.16	118.30
1	A	323	ARG	NE-CZ-NH1	5.29	122.95	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	114	LYS	CB-CA-C	-5.27	99.86	110.40
1	A	125	ASP	CB-CG-OD1	5.19	122.97	118.30
1	A	316	MET	CG-SD-CE	-5.16	91.94	100.20
1	A	58	ARG	NE-CZ-NH2	-5.13	117.74	120.30
1	A	177	ARG	CD-NE-CZ	5.07	130.70	123.60
1	B	282	ASP	CB-CG-OD1	5.04	122.84	118.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	512	HIS	Peptide
1	B	512	HIS	Peptide

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	3766	0	3566	35	0
1	B	3766	0	3566	39	0
2	A	12	12	12	0	0
2	B	12	0	11	2	0
3	B	2	0	0	3	0
4	A	109	0	0	0	0
4	B	113	0	0	3	0
All	All	7780	12	7155	76	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 5.

All (76) 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)
3:B:602:CL:CL	4:B:795:HOH:O	2.15	1.00
3:B:601:CL:CL	4:B:779:HOH:O	2.21	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:276:GLN:N	1:A:299:MET:HE1	1.88	0.89
1:A:313:PRO:HD2	1:A:316:MET:HE3	1.56	0.85
1:B:276:GLN:N	1:B:299:MET:HE1	1.97	0.79
1:B:313:PRO:HD2	1:B:316:MET:HE3	1.64	0.79
1:A:276:GLN:H	1:A:299:MET:HE1	1.45	0.79
1:B:276:GLN:H	1:B:299:MET:HE1	1.51	0.75
1:B:434:SER:OG	3:B:601:CL:CL	2.42	0.74
1:A:313:PRO:CD	1:A:316:MET:HE3	2.20	0.72
1:A:313:PRO:HD2	1:A:316:MET:CE	2.21	0.69
1:A:15:THR:HA	1:A:447:GLN:OE1	1.95	0.67
1:A:512:HIS:O	1:A:513:LYS:HB2	1.99	0.63
1:B:294:ARG:HD2	1:B:371:ILE:O	2.00	0.62
1:B:205:GLY:O	1:B:206:ARG:CB	2.48	0.62
1:B:279:GLU:O	1:B:351:TYR:HA	2.00	0.61
1:B:469:TRP:CZ2	2:B:603:BGC:H3	2.38	0.59
1:A:256:LEU:C	1:A:256:LEU:HD23	2.23	0.58
1:A:279:GLU:O	1:A:351:TYR:HA	2.03	0.58
1:A:294:ARG:HD2	1:A:371:ILE:O	2.02	0.58
1:B:313:PRO:HD2	1:B:316:MET:CE	2.34	0.57
1:B:313:PRO:CD	1:B:316:MET:HE3	2.35	0.57
1:A:313:PRO:CD	1:A:316:MET:CE	2.83	0.56
1:A:186:GLN:HE21	1:A:274:ALA:HB2	1.70	0.55
1:A:469:TRP:CD2	1:A:470:SER:HB3	2.41	0.55
1:A:256:LEU:O	1:A:256:LEU:HD23	2.06	0.55
1:B:297:ASP:OD2	1:B:315:SER:OG	2.24	0.54
1:B:298:PHE:O	1:B:367:TYR:OH	2.24	0.52
1:B:512:HIS:O	1:B:513:LYS:HB2	2.10	0.52
1:B:186:GLN:HG2	1:B:345:ASN:HD22	1.75	0.52
1:B:244:HIS:CE1	1:B:327:PHE:CE1	2.98	0.52
1:A:499:TYR:CD1	1:A:500:PRO:HD2	2.44	0.52
1:B:476:GLU:OE1	2:B:603:BGC:O4	2.24	0.51
1:B:294:ARG:NH1	1:B:372:HIS:O	2.44	0.50
1:B:316:MET:O	1:B:320:VAL:HG23	2.12	0.49
1:A:31:GLY:O	1:A:470:SER:HB2	2.12	0.49
1:A:205:GLY:O	1:A:206:ARG:CB	2.59	0.49
1:A:346:TYR:OH	1:A:399:GLY:HA3	2.13	0.48
1:A:273:HIS:CE1	1:A:303:PHE:HB3	2.49	0.47
1:B:258:LYS:HG2	1:B:262:GLN:OE1	2.14	0.47
1:A:345:ASN:HB3	1:A:420:GLU:HB2	1.96	0.47
1:A:469:TRP:CE2	1:A:470:SER:HB3	2.49	0.47
1:A:346:TYR:CD1	1:A:346:TYR:C	2.88	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:469:TRP:HA	1:A:470:SER:HA	1.62	0.46
1:A:294:ARG:NH1	1:A:372:HIS:O	2.49	0.46
1:B:508:MET:O	1:B:512:HIS:HB2	2.16	0.45
1:A:258:LYS:HG2	1:A:262:GLN:OE1	2.15	0.45
1:B:469:TRP:CD2	1:B:470:SER:HB3	2.51	0.45
1:B:114:LYS:HB3	1:B:114:LYS:HE2	1.77	0.45
1:A:244:HIS:CE1	1:A:327:PHE:CE1	3.05	0.45
1:A:313:PRO:HG2	1:A:316:MET:CE	2.47	0.44
1:B:464:LYS:HA	1:B:464:LYS:HD2	1.75	0.43
1:B:186:GLN:NE2	1:B:189:THR:OG1	2.51	0.43
1:A:186:GLN:HG3	1:A:345:ASN:HD22	1.84	0.43
1:B:114:LYS:NZ	4:B:810:HOH:O	2.51	0.43
1:B:403:ILE:HD12	1:B:403:ILE:HA	1.89	0.43
1:B:305:GLU:HB3	1:B:306:PRO:HD3	1.99	0.43
1:B:443:LEU:HD11	1:B:447:GLN:HE21	1.83	0.42
1:B:344:LEU:HD12	1:B:404:LEU:HD23	2.02	0.42
1:B:94:ARG:HA	1:B:135:PHE:O	2.19	0.42
1:B:346:TYR:OH	1:B:399:GLY:HA3	2.20	0.42
1:B:154:PHE:CE1	1:B:246:LEU:HD23	2.55	0.42
1:B:469:TRP:CE2	1:B:470:SER:HB3	2.55	0.42
1:A:313:PRO:HG2	1:A:316:MET:HE2	2.02	0.41
1:B:315:SER:O	1:B:319:PHE:HD1	2.02	0.41
1:A:176:ASP:OD2	1:A:177:ARG:HD3	2.21	0.41
1:B:142:ASP:OD1	1:B:142:ASP:N	2.49	0.41
1:A:275:THR:C	1:A:299:MET:HE1	2.41	0.41
1:A:71:ALA:HA	1:A:481:TYR:OH	2.21	0.41
1:B:186:GLN:CG	1:B:345:ASN:HD22	2.32	0.41
1:A:114:LYS:HE2	1:A:114:LYS:HB3	1.31	0.41
1:B:31:GLY:O	1:B:470:SER:HB2	2.21	0.40
1:A:243:THR:HG23	1:A:303:PHE:CE1	2.56	0.40
1:B:176:ASP:OD2	1:B:177:ARG:HD3	2.21	0.40
1:B:469:TRP:HA	1:B:470:SER:HA	1.57	0.40
1:A:464:LYS:HA	1:A:464:LYS:HD2	1.83	0.40

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	464/513 (90%)	445 (96%)	18 (4%)	1 (0%)	52	64
1	B	464/513 (90%)	444 (96%)	19 (4%)	1 (0%)	52	64
All	All	928/1026 (90%)	889 (96%)	37 (4%)	2 (0%)	52	64

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	285	SER
1	B	285	SER

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	395/435 (91%)	376 (95%)	19 (5%)	31	42
1	B	395/435 (91%)	377 (95%)	18 (5%)	33	44
All	All	790/870 (91%)	753 (95%)	37 (5%)	32	43

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	THR
1	A	16	ARG
1	A	63	ARG
1	A	107	ARG

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Mol	Chain	Res	Type
1	A	114	LYS
1	A	177	ARG
1	A	182	MET
1	A	204	ARG
1	A	231	SER
1	A	333	LYS
1	A	354	ASN
1	A	378	ASP
1	A	409	LYS
1	A	412	ASN
1	A	427	LYS
1	A	434	SER
1	A	440	SER
1	A	474	ASN
1	A	513	LYS
1	B	15	THR
1	B	16	ARG
1	B	63	ARG
1	B	107	ARG
1	B	114	LYS
1	B	177	ARG
1	B	186	GLN
1	B	231	SER
1	B	307	ILE
1	B	315	SER
1	B	330	GLU
1	B	333	LYS
1	B	354	ASN
1	B	378	ASP
1	B	409	LYS
1	B	412	ASN
1	B	434	SER
1	B	513	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	474	ASN
1	B	186	GLN
1	B	447	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 ⓘ

Of 4 ligands modelled in this entry, 2 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$
2	BGC	A	601	-	12,12,12	1.11	2 (16%)	17,17,17	2.08	6 (35%)
2	BGC	B	603	-	12,12,12	0.42	0	17,17,17	3.34	4 (23%)

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	BGC	A	601	-	-	0/2/22/22	0/1/1/1
2	BGC	B	603	-	-	0/2/22/22	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	BGC	O5-C1	-2.27	1.38	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	BGC	O5-C5	-2.10	1.39	1.44

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	603	BGC	C4-C3-C2	-4.72	101.98	110.79
2	A	601	BGC	O1-C1-O5	-4.27	98.57	110.25
2	B	603	BGC	O3-C3-C2	-2.99	103.62	110.34
2	A	601	BGC	O2-C2-C1	-2.55	104.19	109.82
2	A	601	BGC	O5-C5-C6	-2.48	100.08	106.36
2	A	601	BGC	C1-C2-C3	2.29	113.83	110.43
2	A	601	BGC	C1-O5-C5	2.43	117.97	113.47
2	B	603	BGC	C6-C5-C4	3.01	120.44	113.02
2	A	601	BGC	O5-C1-C2	4.39	116.80	109.80
2	B	603	BGC	C3-C4-C5	11.58	130.39	110.20

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	603	BGC	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	470/513 (91%)	-0.43	9 (1%) 70 76	26, 40, 74, 107	0
1	B	470/513 (91%)	-0.55	7 (1%) 76 81	27, 42, 73, 118	0
All	All	940/1026 (91%)	-0.49	16 (1%) 73 79	26, 41, 74, 118	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	380	ASN	5.0
1	A	382	VAL	4.2
1	A	284	ASN	4.0
1	B	364	ASN	3.9
1	A	381	GLY	3.4
1	B	231	SER	3.3
1	B	356	SER	3.3
1	B	365	PHE	2.9
1	B	206	ARG	2.6
1	A	281	TRP	2.6
1	B	286	ALA	2.5
1	A	283	GLU	2.3
1	A	286	ALA	2.3
1	A	383	PRO	2.2
1	A	285	SER	2.1
1	B	232	THR	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(\AA^2)	Q<0.9
2	BGC	A	601	12/12	0.82	0.25	2.83	20,20,20,20	0
2	BGC	B	603	12/12	0.90	0.16	1.37	36,55,58,60	0
3	CL	B	602	1/1	0.99	0.03	-	58,58,58,58	0
3	CL	B	601	1/1	0.99	0.05	-	55,55,55,55	0

6.5 Other polymers [i](#)

There are no such residues in this entry.