



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:02 PM GMT

PDB ID : 4UOZ
Title : beta-(1,6)-galactosidase from Bifidobacterium animalis subsp. lactis Bl-04 nucleophile mutant E324A in complex with galactose
Authors : Viborg, A.H.; Fredslund, F.; Katayama, T.; Nielsen, S.K.; Svensson, B.; Kitaoka, M.; Lo Leggio, L.; Abou Hachem, M.
Deposited on : 2014-06-11
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

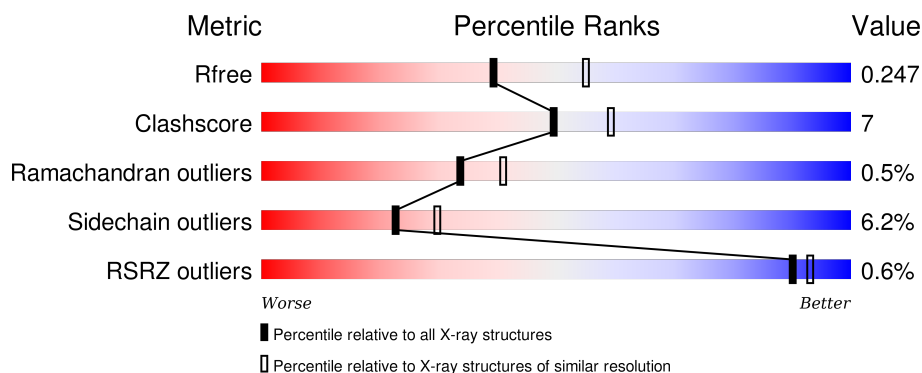
1 Overall quality at a glance

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	695	<div> <div style="width: 100%; height: 10px; background-color: red; position: relative;"> % </div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> 81% 16% .. </div> </div>
1	B	695	<div> <div style="width: 100%; height: 10px; background-color: green; position: relative;"> </div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> 83% 14% .. </div> </div>
1	C	695	<div> <div style="width: 100%; height: 10px; background-color: red; position: relative;"> % </div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> 82% 15% .. </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	GLA	A	701	-	-	-	X
4	PGE	A	1697	-	-	-	X
4	PGE	A	1698	-	-	-	X
4	PGE	B	1696	-	-	-	X
4	PGE	C	1696	-	-	-	X
4	PGE	C	1697	-	-	-	X

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 17041 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.

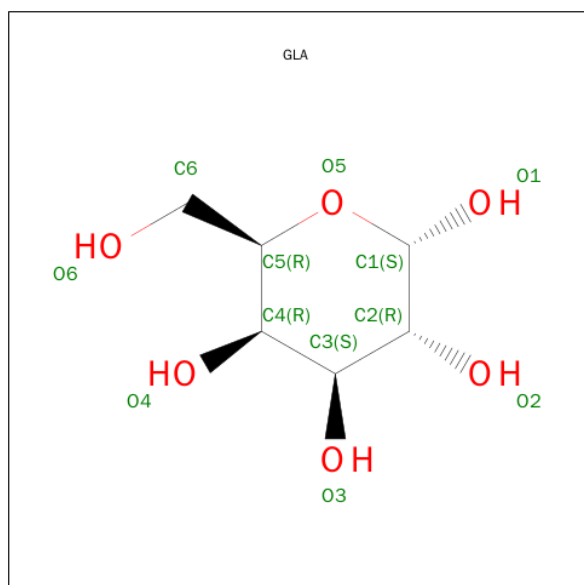
- Molecule 1 is a protein called BETA-GALACTOSIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	688	Total	C	N	O	S	0	1	0
			5474	3460	954	1037	23			
1	B	688	Total	C	N	O	S	0	1	0
			5474	3460	955	1036	23			
1	C	688	Total	C	N	O	S	0	0	0
			5469	3457	954	1035	23			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	324	ALA	GLU	ENGINEERED MUTATION	UNP C6A6W5
B	324	ALA	GLU	ENGINEERED MUTATION	UNP C6A6W5
C	324	ALA	GLU	ENGINEERED MUTATION	UNP C6A6W5

- Molecule 2 is SUGAR (ALPHA D-GALACTOSE) (three-letter code: GLA) (formula: C₆H₁₂O₆).

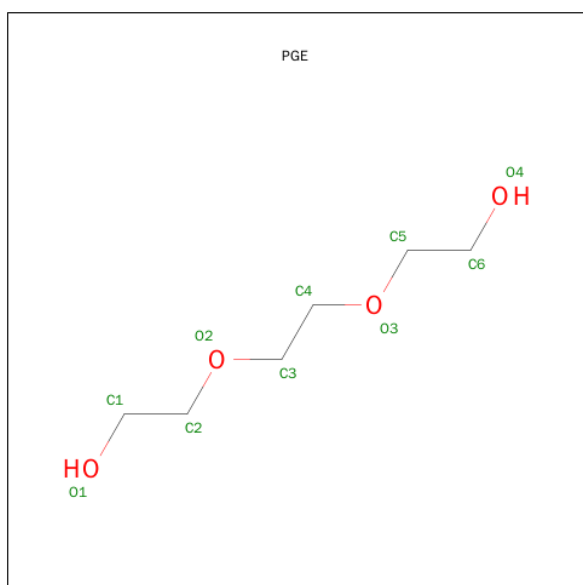


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			12	6	6		
2	B	1	Total	C	O	0	0
			12	6	6		
2	C	1	Total	C	O	0	0
			12	6	6		

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		

- Molecule 4 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: C₆H₁₄O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			10	6	4		
4	A	1	Total	C	O	0	0
			10	6	4		
4	B	1	Total	C	O	0	0
			10	6	4		
4	B	1	Total	C	O	0	0
			10	6	4		
4	C	1	Total	C	O	0	0
			10	6	4		
4	C	1	Total	C	O	0	0
			10	6	4		

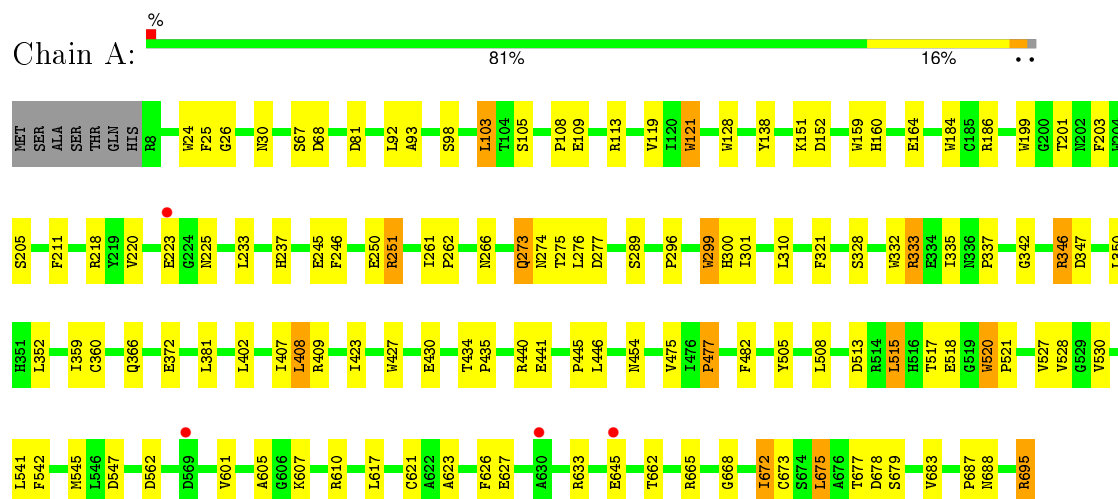
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	185	Total 185	O 185	0	0
5	B	200	Total 200	O 200	0	0
5	C	142	Total 142	O 142	0	0

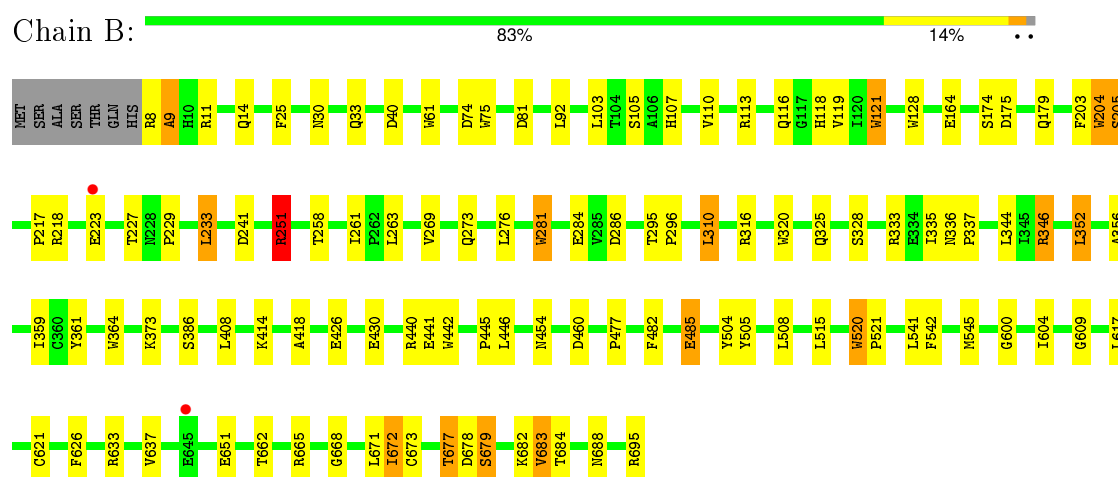
3 Residue-property plots

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.

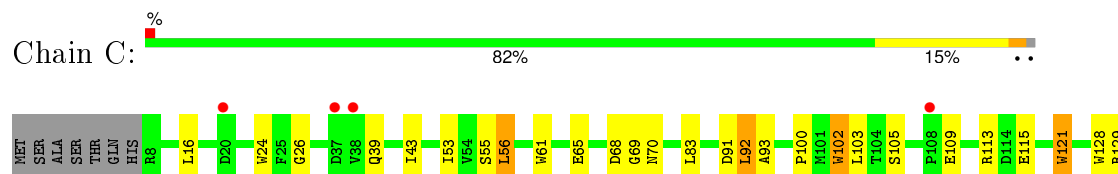
• Molecule 1: BETA-GALACTOSIDASE

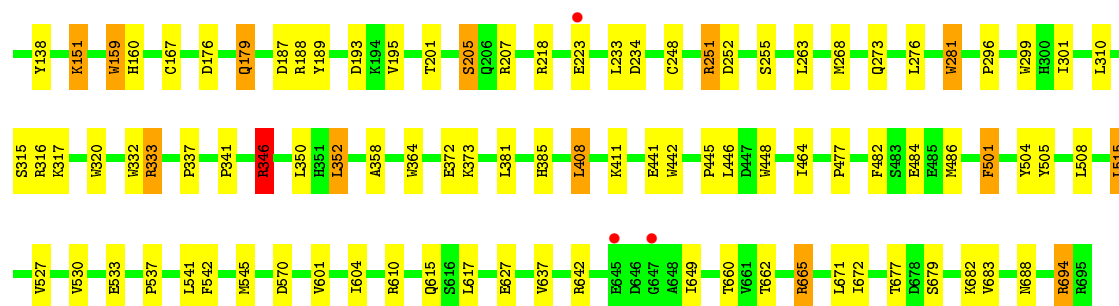


• Molecule 1: BETA-GALACTOSIDASE



• Molecule 1: BETA-GALACTOSIDASE





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	134.20Å 167.97Å 108.58Å 90.00° 115.64° 90.00°	Depositor
Resolution (Å)	28.68 – 2.30 28.68 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.9 (28.68-2.30) 99.9 (28.68-2.30)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.02 (at 2.31Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.186 , 0.248 0.186 , 0.247	Depositor DCC
R_{free} test set	4795 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	22.3	Xtriage
Anisotropy	0.063	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 31.1	EDS
Estimated twinning fraction	0.085 for -1/2*h+1/2*k-l,1/2*h-1/2*k-l,-1/2 *h-1/2*k 0.045 for -1/2*h-1/2*k-l,-1/2*h-1/2*k+l,-1/ 2*h+1/2*k	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	1 of 95831 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	17041	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 2.98% 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: GLA, ZN, PGE

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	0.91	9/5631 (0.2%)	0.93	8/7677 (0.1%)
1	B	0.95	10/5631 (0.2%)	0.96	12/7678 (0.2%)
1	C	0.90	10/5623 (0.2%)	0.93	7/7666 (0.1%)
All	All	0.92	29/16885 (0.2%)	0.94	27/23021 (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	B	0	1

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	281	TRP	CD2-CE2	6.95	1.49	1.41
1	A	121	TRP	CD2-CE2	6.55	1.49	1.41
1	A	184	TRP	CD2-CE2	6.48	1.49	1.41
1	C	320	TRP	CD2-CE2	6.47	1.49	1.41
1	C	281	TRP	CD2-CE2	6.47	1.49	1.41
1	C	332	TRP	CD2-CE2	6.31	1.49	1.41
1	B	121	TRP	CD2-CE2	6.29	1.48	1.41
1	C	448	TRP	CD2-CE2	6.25	1.48	1.41
1	B	204	TRP	CD2-CE2	6.18	1.48	1.41
1	C	102	TRP	CD2-CE2	5.97	1.48	1.41
1	B	520	TRP	CD2-CE2	5.97	1.48	1.41
1	B	128	TRP	CD2-CE2	5.93	1.48	1.41
1	A	24	TRP	CD2-CE2	5.82	1.48	1.41
1	A	332	TRP	CD2-CE2	5.81	1.48	1.41
1	C	128	TRP	CD2-CE2	5.78	1.48	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	299	TRP	CD2-CE2	5.78	1.48	1.41
1	A	520	TRP	CD2-CE2	5.72	1.48	1.41
1	B	442	TRP	CD2-CE2	5.62	1.48	1.41
1	C	442	TRP	CD2-CE2	5.60	1.48	1.41
1	A	128	TRP	CD2-CE2	5.57	1.48	1.41
1	B	609	GLY	N-CA	5.37	1.54	1.46
1	A	199	TRP	CD2-CE2	5.34	1.47	1.41
1	C	364	TRP	CD2-CE2	5.25	1.47	1.41
1	B	364	TRP	CD2-CE2	5.25	1.47	1.41
1	C	121	TRP	CD2-CE2	5.24	1.47	1.41
1	A	159	TRP	CD2-CE2	5.14	1.47	1.41
1	C	159	TRP	CD2-CE2	5.13	1.47	1.41
1	B	61	TRP	CD2-CE2	5.06	1.47	1.41
1	B	75	TRP	CD2-CE2	5.00	1.47	1.41

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	346	ARG	NE-CZ-NH2	-18.14	111.23	120.30
1	A	346	ARG	NE-CZ-NH2	-13.80	113.40	120.30
1	B	346	ARG	NE-CZ-NH1	10.84	125.72	120.30
1	C	346	ARG	NE-CZ-NH2	-10.69	114.95	120.30
1	A	440	ARG	NE-CZ-NH1	8.23	124.41	120.30
1	B	346	ARG	CG-CD-NE	-8.20	94.57	111.80
1	C	346	ARG	NE-CZ-NH1	7.93	124.26	120.30
1	A	440	ARG	NE-CZ-NH2	-7.37	116.62	120.30
1	C	346	ARG	CG-CD-NE	-7.10	96.89	111.80
1	B	251	ARG	NE-CZ-NH1	7.04	123.82	120.30
1	A	346	ARG	NE-CZ-NH1	6.92	123.76	120.30
1	A	251	ARG	NE-CZ-NH1	6.59	123.59	120.30
1	C	234	ASP	CB-CG-OD1	6.55	124.20	118.30
1	B	440	ARG	NE-CZ-NH2	-6.47	117.06	120.30
1	C	333	ARG	NE-CZ-NH1	-6.39	117.10	120.30
1	C	316	ARG	NE-CZ-NH1	6.06	123.33	120.30
1	B	241	ASP	CB-CG-OD1	5.98	123.68	118.30
1	A	333	ARG	NE-CZ-NH1	-5.66	117.47	120.30
1	A	346	ARG	CG-CD-NE	-5.56	100.12	111.80
1	C	408	LEU	CA-CB-CG	5.41	127.74	115.30
1	B	352	LEU	CA-CB-CG	5.38	127.67	115.30
1	B	346	ARG	CD-NE-CZ	5.31	131.03	123.60
1	B	40	ASP	CB-CG-OD1	5.30	123.07	118.30
1	B	251	ARG	NE-CZ-NH2	-5.18	117.71	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	683	VAL	CG1-CB-CG2	5.17	119.17	110.90
1	A	277	ASP	CB-CG-OD1	5.10	122.89	118.30
1	B	233	LEU	CA-CB-CG	5.01	126.82	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	361	TYR	Peptide

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5474	0	5179	76	0
1	B	5474	0	5181	81	0
1	C	5469	0	5175	82	0
2	A	12	0	12	1	0
2	B	12	0	12	1	0
2	C	12	0	12	0	0
3	A	1	0	0	0	0
4	A	20	0	28	1	0
4	B	20	0	28	1	0
4	C	20	0	28	1	0
5	A	185	0	0	17	0
5	B	200	0	0	21	0
5	C	142	0	0	16	0
All	All	17041	0	15655	229	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 7.

All (229) 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:B:9:ALA:HB1	5:B:2001:HOH:O	1.20	1.32

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:615:GLN:HB2	5:C:2136:HOH:O	1.21	1.31
1:B:118:HIS:HD2	5:B:2052:HOH:O	1.13	1.30
1:C:251:ARG:HD3	5:C:2053:HOH:O	1.27	1.27
1:C:352:LEU:HG	5:C:2085:HOH:O	1.09	1.25
1:B:673:CYS:HB2	5:B:2197:HOH:O	1.41	1.21
1:A:673:CYS:HB2	5:A:2182:HOH:O	1.41	1.19
1:C:251:ARG:HB2	5:C:2053:HOH:O	1.41	1.19
1:C:542:PHE:H	1:C:545:MET:HE3	1.05	1.14
1:C:545:MET:SD	5:C:2120:HOH:O	2.10	1.09
1:B:118:HIS:CD2	5:B:2052:HOH:O	1.93	1.08
1:A:30:ASN:HB2	5:A:2017:HOH:O	1.54	1.04
1:B:673:CYS:CB	5:B:2197:HOH:O	2.02	0.98
1:C:542:PHE:H	1:C:545:MET:CE	1.76	0.96
1:B:673:CYS:SG	5:B:2197:HOH:O	2.22	0.95
1:C:542:PHE:N	1:C:545:MET:HE3	1.85	0.91
1:B:113:ARG:HH21	1:B:119:VAL:HG22	1.35	0.88
1:B:8:ARG:O	1:B:9:ALA:HB3	1.73	0.87
1:B:30[B]:ASN:HD22	1:B:33:GLN:NE2	1.73	0.86
1:B:333:ARG:HD2	5:B:2114:HOH:O	1.74	0.86
1:B:542:PHE:H	1:B:545:MET:HE3	1.40	0.86
1:A:542:PHE:H	1:A:545:MET:HE3	1.40	0.85
1:C:542:PHE:HB2	1:C:545:MET:CE	2.09	0.83
1:A:623:ALA:HA	4:A:1698:PGE:H5	1.62	0.82
1:C:56:LEU:HB3	5:C:2012:HOH:O	1.80	0.81
1:B:542:PHE:N	1:B:545:MET:HE3	1.95	0.81
1:C:251:ARG:CD	5:C:2053:HOH:O	2.02	0.80
1:C:660:THR:O	5:C:2141:HOH:O	1.92	0.80
1:C:56:LEU:CB	5:C:2012:HOH:O	2.31	0.79
1:A:346:ARG:NH2	1:A:688:ASN:OD1	2.17	0.78
1:C:694:ARG:HH11	1:C:694:ARG:HG2	1.49	0.77
1:A:542:PHE:HB2	1:A:545:MET:CE	2.15	0.77
1:A:542:PHE:HD2	1:A:545:MET:HE1	1.50	0.76
1:C:56:LEU:HD12	1:C:92:LEU:HD23	1.66	0.75
1:B:203:PHE:HB3	5:B:2071:HOH:O	1.85	0.74
1:C:542:PHE:N	1:C:545:MET:CE	2.47	0.73
1:A:665:ARG:NH2	1:A:668:GLY:O	2.22	0.73
1:B:8:ARG:O	1:B:9:ALA:CB	2.36	0.73
1:C:333:ARG:HD3	5:C:2077:HOH:O	1.89	0.72
1:A:542:PHE:N	1:A:545:MET:HE3	2.05	0.72
1:B:333:ARG:CD	5:B:2114:HOH:O	2.35	0.72
5:A:2116:HOH:O	1:B:229:PRO:HG2	1.88	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:409:ARG:HG2	5:A:2001:HOH:O	1.88	0.71
1:B:30[B]:ASN:HD22	1:B:33:GLN:CD	1.94	0.69
1:C:542:PHE:HB2	1:C:545:MET:HE2	1.73	0.69
1:C:276:LEU:H	1:C:276:LEU:HD12	1.57	0.68
1:C:248:CYS:O	1:C:251:ARG:HG3	1.94	0.68
1:B:665:ARG:NH2	1:B:668:GLY:O	2.27	0.67
1:B:346:ARG:NH2	1:B:688:ASN:OD1	2.27	0.66
1:C:542:PHE:HD2	1:C:545:MET:HE1	1.59	0.66
1:B:333:ARG:CG	5:B:2114:HOH:O	2.44	0.66
1:C:205:SER:HA	5:C:2042:HOH:O	1.94	0.66
1:B:205:SER:HA	5:B:2072:HOH:O	1.96	0.65
1:C:299:TRP:CD1	4:C:1696:PGE:H42	2.34	0.63
1:C:56:LEU:CD1	1:C:92:LEU:HD23	2.28	0.63
1:B:542:PHE:H	1:B:545:MET:CE	2.10	0.62
1:B:454:ASN:CG	5:B:2165:HOH:O	2.36	0.62
1:B:454:ASN:HB3	5:B:2165:HOH:O	1.98	0.62
1:B:542:PHE:HD2	1:B:545:MET:HE1	1.64	0.62
1:B:671:LEU:O	1:B:672:ILE:HD12	1.99	0.62
1:C:694:ARG:NH1	1:C:694:ARG:HG2	2.14	0.61
1:B:665:ARG:NH1	1:B:678:ASP:OD2	2.33	0.61
1:B:30[B]:ASN:ND2	1:B:33:GLN:CD	2.54	0.61
1:B:333:ARG:HG3	5:B:2114:HOH:O	2.01	0.61
1:C:296:PRO:HB3	1:C:337:PRO:HG2	1.81	0.60
1:A:610:ARG:HD2	5:A:2147:HOH:O	2.02	0.60
1:C:542:PHE:CD2	1:C:545:MET:HE1	2.37	0.60
1:B:8:ARG:CG	1:B:8:ARG:O	2.50	0.60
1:A:372:GLU:OE1	1:B:203:PHE:HZ	1.85	0.59
1:C:179:GLN:HA	1:C:179:GLN:HE21	1.68	0.59
1:C:255:SER:HB2	5:C:2001:HOH:O	2.02	0.59
1:B:276:LEU:HD23	1:B:281:TRP:CZ2	2.38	0.59
1:B:542:PHE:HB2	1:B:545:MET:HE2	1.86	0.58
1:A:542:PHE:CD2	1:A:545:MET:HE1	2.36	0.58
1:C:346:ARG:NH2	1:C:688:ASN:OD1	2.35	0.58
1:B:121:TRP:CD1	1:C:218:ARG:HB3	2.39	0.57
1:A:333:ARG:NH1	5:A:2085:HOH:O	2.36	0.57
1:A:151:LYS:O	1:A:152:ASP:HB2	2.04	0.57
1:A:673:CYS:CB	5:A:2182:HOH:O	2.19	0.56
1:C:55:SER:HA	1:C:91:ASP:O	2.05	0.56
1:C:542:PHE:CB	1:C:545:MET:CE	2.83	0.56
1:A:346:ARG:NH1	1:A:347:ASP:OD1	2.33	0.56
1:A:409:ARG:NH1	5:A:2007:HOH:O	2.39	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:8:ARG:HG2	1:B:8:ARG:O	2.06	0.55
1:C:679:SER:HB3	1:C:682:LYS:O	2.07	0.55
1:A:695:ARG:HH11	1:A:695:ARG:HG2	1.70	0.55
1:B:269:VAL:HG13	1:B:310:LEU:HD13	1.89	0.55
1:B:107:HIS:O	1:B:110:VAL:HG22	2.07	0.55
1:C:352:LEU:HD13	1:C:642:ARG:NH1	2.22	0.54
1:A:109:GLU:O	5:A:2033:HOH:O	2.18	0.54
1:A:513:ASP:OD2	1:C:373:LYS:NZ	2.38	0.54
1:A:515:LEU:HB3	1:C:381:LEU:HD13	1.90	0.54
1:A:203:PHE:HB3	5:A:2057:HOH:O	2.07	0.54
1:B:542:PHE:HB2	1:B:545:MET:CE	2.38	0.54
1:C:352:LEU:CD1	1:C:642:ARG:NH1	2.70	0.54
1:B:333:ARG:HD3	5:B:2113:HOH:O	2.08	0.54
1:C:251:ARG:CB	5:C:2053:HOH:O	2.20	0.54
1:A:93:ALA:HA	1:A:160:HIS:O	2.09	0.53
1:C:570:ASP:OD1	1:C:570:ASP:N	2.39	0.53
1:C:103:LEU:HD11	1:C:138:TYR:CG	2.43	0.53
1:A:203:PHE:HZ	1:C:372:GLU:OE1	1.92	0.52
1:A:607:LYS:HE2	5:A:2142:HOH:O	2.09	0.52
1:B:542:PHE:HD2	1:B:545:MET:CE	2.23	0.51
1:B:386:SER:HA	1:C:533:GLU:OE2	2.10	0.51
1:C:201:THR:OG1	1:C:207:ARG:HA	2.09	0.51
1:C:604:ILE:N	1:C:604:ILE:HD12	2.25	0.51
1:C:39:GLN:O	1:C:43:ILE:HG12	2.11	0.51
1:B:269:VAL:CG1	1:B:310:LEU:HD13	2.40	0.51
1:A:113:ARG:HH21	1:A:119:VAL:CG2	2.23	0.51
1:B:454:ASN:CB	5:B:2165:HOH:O	2.55	0.51
1:B:276:LEU:H	1:B:276:LEU:HD12	1.76	0.50
1:B:373:LYS:HG3	1:C:515:LEU:HD22	1.92	0.50
1:A:520:TRP:HA	1:A:521:PRO:C	2.32	0.50
1:A:621:CYS:HB3	1:A:626:PHE:O	2.12	0.50
1:B:520:TRP:HA	1:B:521:PRO:C	2.31	0.50
1:B:113:ARG:HG3	1:B:174:SER:HB3	1.94	0.50
1:B:651:GLU:OE1	1:B:695:ARG:NH1	2.45	0.50
1:B:103:LEU:HD23	1:B:103:LEU:O	2.11	0.50
1:A:121:TRP:CD1	1:B:218:ARG:HB3	2.47	0.49
1:C:43:ILE:HD13	1:C:83:LEU:HD23	1.95	0.49
1:B:11:ARG:HB2	1:B:286:ASP:HB3	1.95	0.49
1:B:316:ARG:HA	5:B:2102:HOH:O	2.12	0.49
1:C:189:TYR:CD2	1:C:195:VAL:HG22	2.48	0.49
1:A:335:ILE:O	1:A:337:PRO:HD3	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:335:ILE:O	1:B:337:PRO:HD3	2.13	0.49
1:B:251:ARG:HB2	1:B:263:LEU:HD22	1.96	0.48
1:A:275:THR:O	1:A:427:TRP:HZ2	1.97	0.48
1:C:56:LEU:CD2	5:C:2012:HOH:O	2.60	0.48
1:C:109:GLU:O	5:C:2025:HOH:O	2.19	0.48
1:B:679:SER:HB2	1:B:682:LYS:O	2.13	0.48
1:A:186:ARG:HA	1:A:211:PHE:CE2	2.49	0.48
1:B:118:HIS:HA	5:B:2052:HOH:O	2.14	0.48
1:C:527:VAL:HG12	1:C:601:VAL:HG21	1.96	0.48
1:A:164:GLU:OE2	2:A:701:GLA:H1	2.14	0.47
1:C:151:LYS:NZ	1:C:151:LYS:HB2	2.28	0.47
1:A:25:PHE:HA	1:A:359:ILE:O	2.15	0.47
1:C:43:ILE:CD1	1:C:83:LEU:HD23	2.45	0.47
1:A:299:TRP:CH2	1:A:610:ARG:HD3	2.50	0.47
1:A:300:HIS:HE1	1:A:347:ASP:OD2	1.97	0.46
1:C:276:LEU:HD12	1:C:276:LEU:N	2.29	0.46
1:B:296:PRO:HB3	1:B:337:PRO:HG2	1.97	0.46
1:A:273:GLN:HE22	1:A:423:ILE:HD13	1.81	0.46
1:B:203:PHE:CE2	1:B:204:TRP:CE2	3.04	0.46
1:A:113:ARG:HH21	1:A:119:VAL:HG22	1.80	0.46
1:B:164:GLU:OE2	2:B:701:GLA:H1	2.15	0.46
1:A:26:GLY:O	1:A:360:CYS:HA	2.16	0.46
1:A:276:LEU:H	1:A:276:LEU:HD12	1.81	0.46
1:C:93:ALA:HA	1:C:160:HIS:HB3	1.98	0.46
1:A:673:CYS:SG	5:A:2182:HOH:O	2.61	0.45
1:A:108:PRO:HD2	1:A:109:GLU:OE1	2.16	0.45
1:A:237:HIS:HE1	5:A:2076:HOH:O	1.99	0.45
1:C:26:GLY:HA3	1:C:53:ILE:O	2.17	0.45
1:A:562:ASP:HB2	1:A:605:ALA:HB1	1.99	0.45
1:A:218:ARG:HB3	1:C:121:TRP:CD1	2.52	0.45
1:A:237:HIS:HD2	5:A:2074:HOH:O	2.00	0.45
1:C:24:TRP:HB2	1:C:358:ALA:CB	2.46	0.45
1:B:621:CYS:HB3	1:B:626:PHE:O	2.17	0.45
1:A:246:PHE:HE1	1:A:250:GLU:OE2	1.98	0.45
1:A:342:GLY:HA3	1:A:675:LEU:HD22	1.98	0.45
1:C:542:PHE:N	1:C:545:MET:HE1	2.31	0.45
1:C:464:ILE:HD11	1:C:486:MET:SD	2.56	0.45
1:A:542:PHE:HB2	1:A:545:MET:HE3	1.93	0.45
1:B:344:LEU:HD23	1:B:344:LEU:C	2.36	0.45
1:B:346:ARG:HD2	1:B:346:ARG:C	2.38	0.44
1:C:103:LEU:O	1:C:103:LEU:HD23	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:671:LEU:HA	1:C:671:LEU:HD23	1.67	0.44
1:A:205:SER:HA	5:A:2058:HOH:O	2.18	0.44
1:C:68:ASP:HA	1:C:102:TRP:CH2	2.53	0.44
1:A:273:GLN:NE2	1:A:423:ILE:HD13	2.33	0.44
1:C:276:LEU:HD23	1:C:281:TRP:CZ2	2.53	0.44
1:C:501:PHE:CD1	1:C:501:PHE:N	2.86	0.44
1:B:251:ARG:NH2	1:B:284:GLU:OE1	2.51	0.43
1:A:289:SER:HB3	1:A:321:PHE:HB3	1.99	0.43
1:A:296:PRO:HB3	1:A:337:PRO:HG2	2.00	0.43
1:A:301:ILE:HD11	1:A:346:ARG:NH2	2.34	0.43
1:C:317:LYS:O	1:C:411:LYS:HG2	2.18	0.43
1:A:434:THR:HB	1:A:435:PRO:HD2	2.01	0.43
1:C:92:LEU:HD13	1:C:159:TRP:CZ3	2.53	0.43
1:A:475:VAL:O	1:A:477:PRO:HD3	2.19	0.43
1:A:517:THR:C	1:A:518:GLU:HG2	2.39	0.43
1:A:266:ASN:OD1	1:A:266:ASN:N	2.43	0.43
1:A:527:VAL:HG12	1:A:601:VAL:HG21	2.01	0.43
1:C:649:ILE:O	1:C:694:ARG:HA	2.19	0.43
1:B:604:ILE:N	1:B:604:ILE:HD12	2.33	0.43
1:A:103:LEU:HD21	1:A:138:TYR:CD2	2.53	0.43
1:B:74:ASP:HB2	5:B:2040:HOH:O	2.19	0.43
1:B:441:GLU:O	1:B:445:PRO:HD2	2.18	0.43
1:C:255:SER:HB3	1:C:263:LEU:HD11	2.00	0.43
1:A:151:LYS:O	1:A:152:ASP:CB	2.61	0.42
1:C:167:CYS:HA	1:C:268:MET:HE1	2.01	0.42
1:A:441:GLU:O	1:A:445:PRO:HD2	2.20	0.42
1:B:677:THR:HB	1:B:684:THR:HB	2.00	0.42
1:C:129:ARG:NH1	1:C:176:ASP:OD2	2.45	0.42
1:B:320:TRP:O	1:B:356:ALA:HA	2.20	0.42
1:B:113:ARG:NH2	1:B:119:VAL:HG22	2.18	0.42
1:B:203:PHE:CZ	1:B:204:TRP:NE1	2.87	0.42
1:A:408:LEU:O	1:A:409:ARG:HB2	2.18	0.42
1:A:547:ASP:OD1	1:A:547:ASP:C	2.58	0.42
1:C:65:GLU:HA	1:C:70:ASN:O	2.20	0.42
1:A:408:LEU:HB2	5:A:2124:HOH:O	2.20	0.42
1:A:261:ILE:HA	1:A:262:PRO:HD3	1.93	0.42
1:A:687:PRO:O	1:A:688:ASN:HB2	2.19	0.42
1:A:67:SER:O	1:A:68:ASP:C	2.58	0.42
1:B:14:GLN:HB2	5:B:2011:HOH:O	2.19	0.42
1:B:295:THR:HG21	4:B:1696:PGE:H2	2.01	0.41
1:A:665:ARG:NH1	1:A:678:ASP:OD1	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:665:ARG:HD2	1:C:679:SER:O	2.20	0.41
1:A:402:LEU:O	1:A:407:ILE:HG22	2.20	0.41
1:B:258:THR:HB	1:B:261:ILE:HD12	2.02	0.41
1:C:251:ARG:HD2	1:C:252:ASP:OD1	2.21	0.41
1:C:61:TRP:CG	1:C:100:PRO:HD3	2.55	0.41
1:B:426:GLU:OE2	1:B:430:GLU:OE2	2.38	0.41
1:C:56:LEU:HB2	5:C:2012:HOH:O	2.10	0.41
1:C:441:GLU:O	1:C:445:PRO:HD2	2.21	0.41
1:A:672:ILE:HD12	1:A:672:ILE:HA	1.81	0.41
1:A:274:ASN:HB2	1:A:430:GLU:OE2	2.21	0.41
1:B:116:GLN:HB2	1:B:118:HIS:CE1	2.55	0.41
1:C:301:ILE:HD11	1:C:346:ARG:NH2	2.35	0.41
1:C:346:ARG:C	1:C:346:ARG:HD2	2.41	0.41
1:B:25:PHE:HA	1:B:359:ILE:O	2.21	0.41
1:B:485:GLU:CD	1:B:485:GLU:H	2.23	0.41
1:B:418:ALA:HA	1:B:460:ASP:O	2.21	0.41
1:B:414:LYS:NZ	5:B:2149:HOH:O	2.48	0.40
5:A:2175:HOH:O	1:C:341:PRO:HG3	2.20	0.40
1:B:336:ASN:O	1:C:537:PRO:HG2	2.22	0.40
1:A:289:SER:HA	1:A:321:PHE:O	2.21	0.40
1:A:366:GLN:HB3	1:A:381:LEU:HG	2.02	0.40
1:A:528:VAL:HG12	1:A:530:VAL:HB	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	687/695 (99%)	659 (96%)	26 (4%)	2 (0%)	46 57
1	B	687/695 (99%)	654 (95%)	27 (4%)	6 (1%)	21 24
1	C	686/695 (99%)	657 (96%)	26 (4%)	3 (0%)	39 48

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	2060/2085 (99%)	1970 (96%)	79 (4%)	11 (0%)	34	41

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	9	ALA
1	B	217	PRO
1	B	477	PRO
1	B	504	TYR
1	C	504	TYR
1	A	201	THR
1	B	600	GLY
1	B	328	SER
1	C	477	PRO
1	C	69	GLY
1	A	477	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	575/580 (99%)	540 (94%)	35 (6%)	23	30
1	B	575/580 (99%)	545 (95%)	30 (5%)	29	38
1	C	574/580 (99%)	532 (93%)	42 (7%)	17	22
All	All	1724/1740 (99%)	1617 (94%)	107 (6%)	23	30

All (107) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	81	ASP
1	A	92	LEU
1	A	98	SER
1	A	103	LEU
1	A	105	SER

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Mol	Chain	Res	Type
1	A	220	VAL
1	A	223	GLU
1	A	225	ASN
1	A	233	LEU
1	A	245	GLU
1	A	251	ARG
1	A	273	GLN
1	A	310	LEU
1	A	328	SER
1	A	350	LEU
1	A	352	LEU
1	A	408	LEU
1	A	446	LEU
1	A	454	ASN
1	A	482	PHE
1	A	505	TYR
1	A	508	LEU
1	A	515	LEU
1	A	541	LEU
1	A	617	LEU
1	A	627	GLU
1	A	633	ARG
1	A	645	GLU
1	A	662	THR
1	A	672	ILE
1	A	675	LEU
1	A	677	THR
1	A	679	SER
1	A	683	VAL
1	A	695	ARG
1	B	81	ASP
1	B	92	LEU
1	B	105	SER
1	B	175	ASP
1	B	179	GLN
1	B	205	SER
1	B	223	GLU
1	B	227	THR
1	B	233	LEU
1	B	251	ARG
1	B	273	GLN
1	B	310	LEU

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Mol	Chain	Res	Type
1	B	325	GLN
1	B	352	LEU
1	B	408	LEU
1	B	446	LEU
1	B	482	PHE
1	B	485	GLU
1	B	505	TYR
1	B	508	LEU
1	B	515	LEU
1	B	541	LEU
1	B	617	LEU
1	B	633	ARG
1	B	637	VAL
1	B	662	THR
1	B	672	ILE
1	B	677	THR
1	B	679	SER
1	B	683	VAL
1	C	16	LEU
1	C	56	LEU
1	C	92	LEU
1	C	105	SER
1	C	113	ARG
1	C	115	GLU
1	C	151	LYS
1	C	179	GLN
1	C	187	ASP
1	C	188	ARG
1	C	193	ASP
1	C	205	SER
1	C	223	GLU
1	C	233	LEU
1	C	251	ARG
1	C	273	GLN
1	C	310	LEU
1	C	315	SER
1	C	346	ARG
1	C	350	LEU
1	C	352	LEU
1	C	385	HIS
1	C	408	LEU
1	C	446	LEU

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Mol	Chain	Res	Type
1	C	482	PHE
1	C	484	GLU
1	C	501	PHE
1	C	505	TYR
1	C	508	LEU
1	C	515	LEU
1	C	530	VAL
1	C	541	LEU
1	C	610	ARG
1	C	617	LEU
1	C	627	GLU
1	C	637	VAL
1	C	662	THR
1	C	665	ARG
1	C	672	ILE
1	C	677	THR
1	C	683	VAL
1	C	694	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	237	HIS
1	A	273	GLN
1	A	300	HIS
1	A	325	GLN
1	A	454	ASN
1	A	465	HIS
1	B	33	GLN
1	B	273	GLN
1	B	325	GLN
1	B	465	HIS
1	C	179	GLN
1	C	237	HIS

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 10 ligands modelled in this entry, 1 is monoatomic - leaving 9 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	$\# Z > 2$
4	PGE	A	1697	-	9,9,9	0.51	0	8,8,8	0.47	0
4	PGE	A	1698	-	9,9,9	0.86	0	8,8,8	0.99	0
2	GLA	A	701	-	12,12,12	0.82	0	17,17,17	1.52	3 (17%)
4	PGE	B	1696	-	9,9,9	0.54	0	8,8,8	0.73	0
4	PGE	B	1697	-	9,9,9	0.83	0	8,8,8	0.87	1 (12%)
2	GLA	B	701	-	12,12,12	0.79	0	17,17,17	1.74	5 (29%)
4	PGE	C	1696	-	9,9,9	0.53	0	8,8,8	0.72	0
4	PGE	C	1697	-	9,9,9	0.93	0	8,8,8	0.85	0
2	GLA	C	701	-	12,12,12	1.04	1 (8%)	17,17,17	1.84	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
4	PGE	A	1697	-	-	0/7/7/7	0/0/0/0
4	PGE	A	1698	-	-	0/7/7/7	0/0/0/0
2	GLA	A	701	-	-	0/2/22/22	0/1/1/1
4	PGE	B	1696	-	-	0/7/7/7	0/0/0/0
4	PGE	B	1697	-	-	0/7/7/7	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GLA	B	701	-	-	0/2/22/22	0/1/1/1
4	PGE	C	1696	-	-	0/7/7/7	0/0/0/0
4	PGE	C	1697	-	-	0/7/7/7	0/0/0/0
2	GLA	C	701	-	-	0/2/22/22	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	701	GLA	C4-C5	2.78	1.59	1.53

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	701	GLA	C1-C2-C3	-4.07	104.37	110.43
2	B	701	GLA	O3-C3-C2	-2.94	103.72	110.34
2	B	701	GLA	C1-C2-C3	-2.18	107.18	110.43
2	C	701	GLA	C6-C5-C4	2.05	118.07	113.02
4	B	1697	PGE	O2-C3-C4	2.09	119.67	110.36
2	B	701	GLA	O6-C6-C5	2.11	118.32	111.33
2	A	701	GLA	O3-C3-C4	2.33	115.57	110.34
2	A	701	GLA	C1-O5-C5	2.37	117.85	113.47
2	C	701	GLA	O5-C1-C2	2.58	113.92	109.80
2	B	701	GLA	O3-C3-C4	2.70	116.41	110.34
2	C	701	GLA	C1-O5-C5	3.35	119.66	113.47
2	B	701	GLA	C1-O5-C5	3.60	120.12	113.47
2	C	701	GLA	O3-C3-C4	3.86	119.03	110.34

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1698	PGE	1	0
2	A	701	GLA	1	0
4	B	1696	PGE	1	0
2	B	701	GLA	1	0
4	C	1696	PGE	1	0

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 [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	688/695 (98%)	-0.29	4 (0%) 90 93	11, 22, 38, 52	0
1	B	688/695 (98%)	-0.34	2 (0%) 94 96	8, 18, 36, 51	0
1	C	688/695 (98%)	-0.22	7 (1%) 84 88	10, 23, 41, 62	0
All	All	2064/2085 (98%)	-0.28	13 (0%) 90 93	8, 21, 39, 62	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	645	GLU	3.3
1	C	223	GLU	2.9
1	C	108	PRO	2.9
1	A	630	ALA	2.9
1	A	569	ASP	2.8
1	C	38	VAL	2.7
1	C	647	GLY	2.5
1	C	645	GLU	2.5
1	B	645	GLU	2.4
1	C	20	ASP	2.3
1	C	37	ASP	2.2
1	A	223	GLU	2.1
1	B	223	GLU	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
4	PGE	C	1697	10/10	0.78	0.23	4.15	28,37,44,45	0
4	PGE	C	1696	10/10	0.91	0.17	3.80	28,31,47,51	0
4	PGE	B	1696	10/10	0.88	0.16	3.73	31,33,44,45	0
4	PGE	A	1697	10/10	0.90	0.17	3.70	32,36,51,52	0
4	PGE	A	1698	10/10	0.88	0.17	2.20	28,30,39,39	0
2	GLA	A	701	12/12	0.90	0.18	2.05	28,33,35,37	0
4	PGE	B	1697	10/10	0.77	0.16	1.80	36,40,44,49	0
2	GLA	B	701	12/12	0.95	0.12	0.52	23,26,28,28	0
2	GLA	C	701	12/12	0.94	0.12	0.28	27,30,33,33	0
3	ZN	A	1696	1/1	0.99	0.05	-3.78	43,43,43,43	0

6.5 Other polymers [i](#)

There are no such residues in this entry.