



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 5, 2016 – 12:13 PM EDT

PDB ID : 5LDR
Title : Crystal structure of a cold-adapted dimeric beta-D-galactosidase from *Paracoccus* sp. 32d strain in complex with galactose
Authors : Rutkiewicz-Krotewicz, M.; Bujacz, A.; Pietrzyk, A.J.; Sekula, B.; Bujacz, G.
Deposited on : 2016-06-27
Resolution : 3.15 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027939
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) : rb-20027939

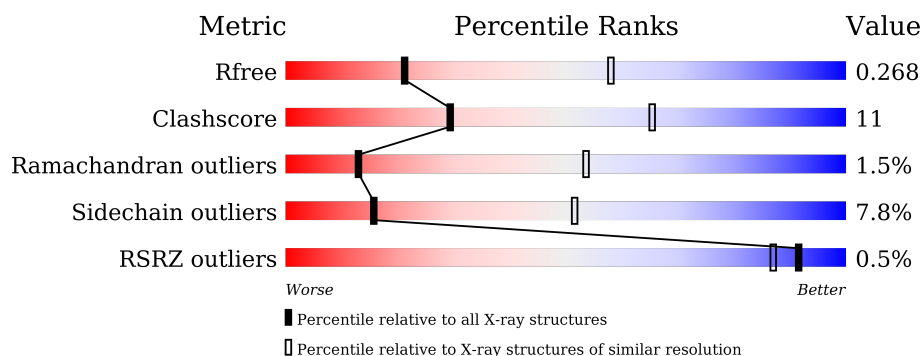
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.15 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	1112 (3.20-3.12)
Clashscore	102246	1249 (3.20-3.12)
Ramachandran outliers	100387	1222 (3.20-3.12)
Sidechain outliers	100360	1221 (3.20-3.12)
RSRZ outliers	91569	1117 (3.20-3.12)

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	732	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 26%, green 71%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 71% 26% . </div> </div>
2	B	731	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 26%, green 71%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 71% 26% . </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
5	PEG	A	803	-	-	-	X
5	PEG	B	802	-	-	-	X
5	PEG	B	805	-	-	X	X
7	ACT	B	808	-	-	-	X

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 11724 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-D-galactosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	731	Total	C	N	O	S	0	0	0
			5764	3632	1038	1068	26			

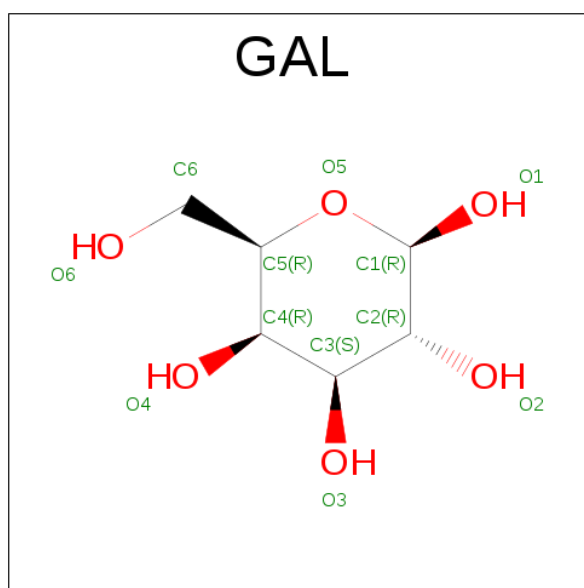
There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	ASN	-	expression tag	UNP D1LZK0

- Molecule 2 is a protein called Beta-D-galactosidase.

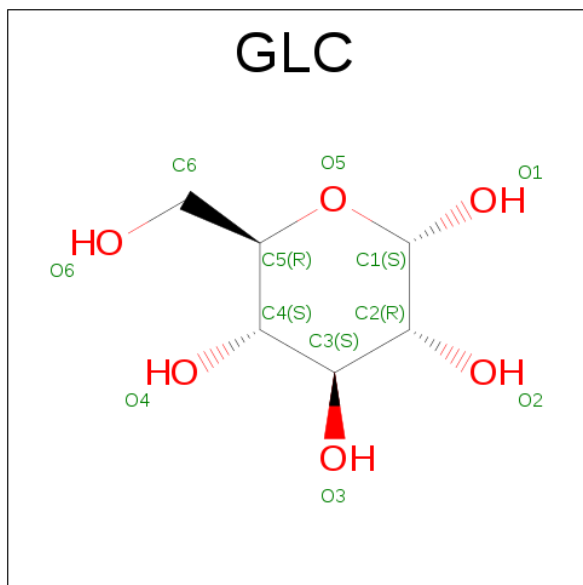
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	731	Total	C	N	O	S	0	0	0
			5763	3632	1038	1067	26			

- Molecule 3 is BETA-D-GALACTOSE (three-letter code: GAL) (formula: C₆H₁₂O₆).



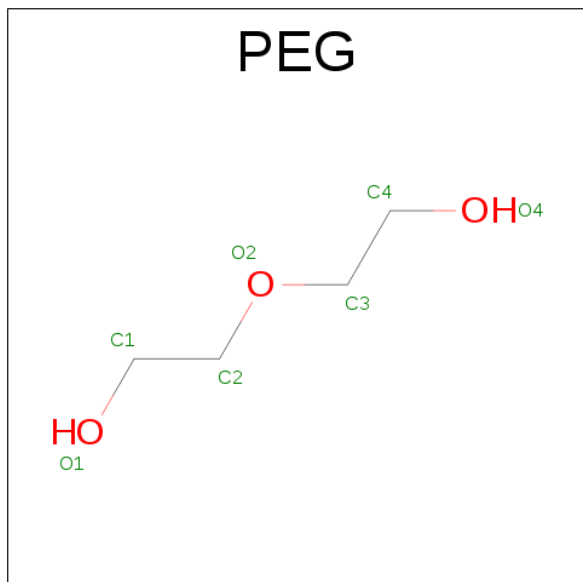
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			12	6	6		
3	B	1	Total	C	O	0	0
			12	6	6		

- Molecule 4 is ALPHA-D-GLUCOSE (three-letter code: GLC) (formula: C₆H₁₂O₆).



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

- Molecule 5 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).

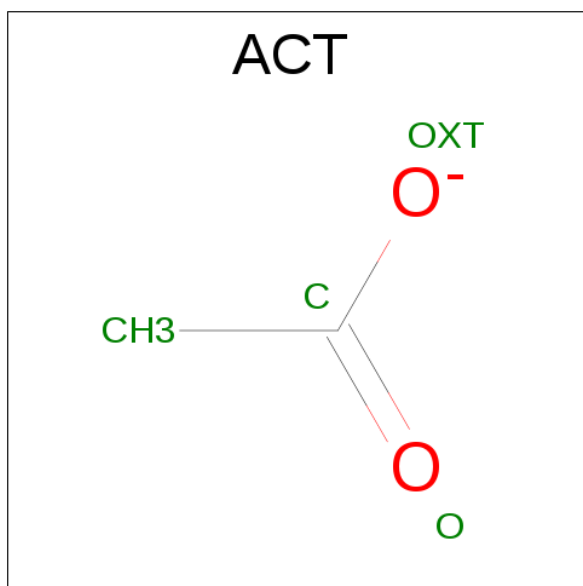


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 7 4 3	0	0
5	B	1	Total C O 7 4 3	0	0
5	B	1	Total C O 7 4 3	0	0
5	B	1	Total C O 7 4 3	0	0
5	B	1	Total C O 7 4 3	0	0

- Molecule 6 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

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

- Molecule 7 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	B	1	Total C O 4 2 2	0	0
7	B	1	Total C O 4 2 2	0	0

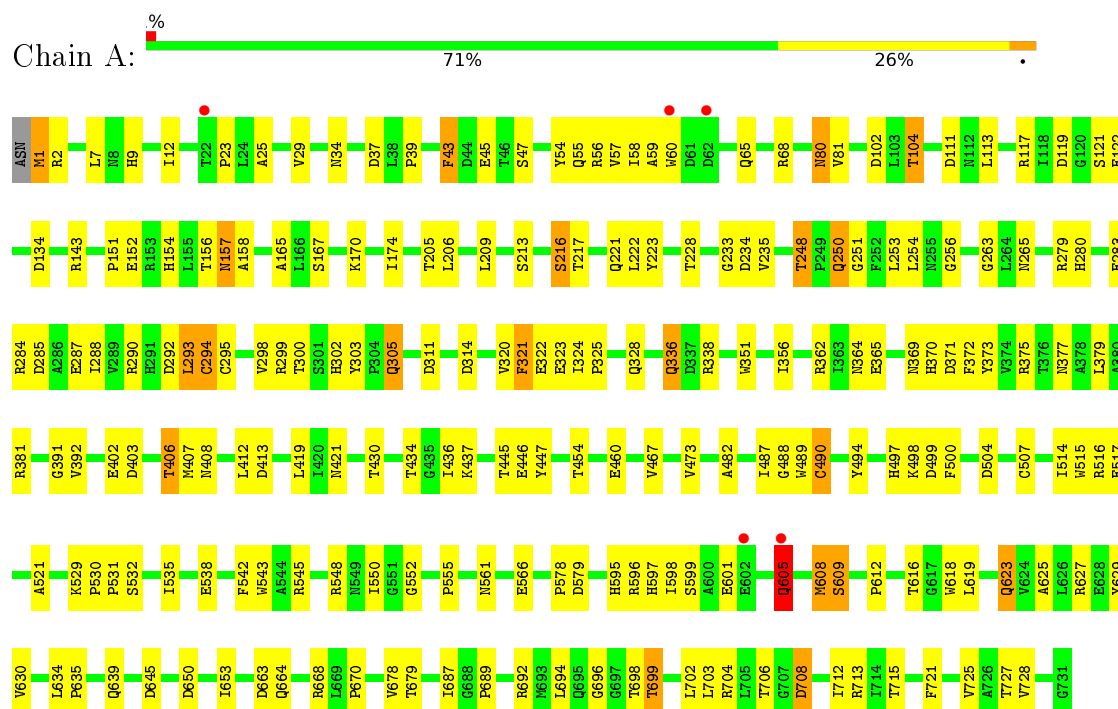
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	66	Total 66	O 66	0	0
8	B	48	Total 48	O 48	0	0

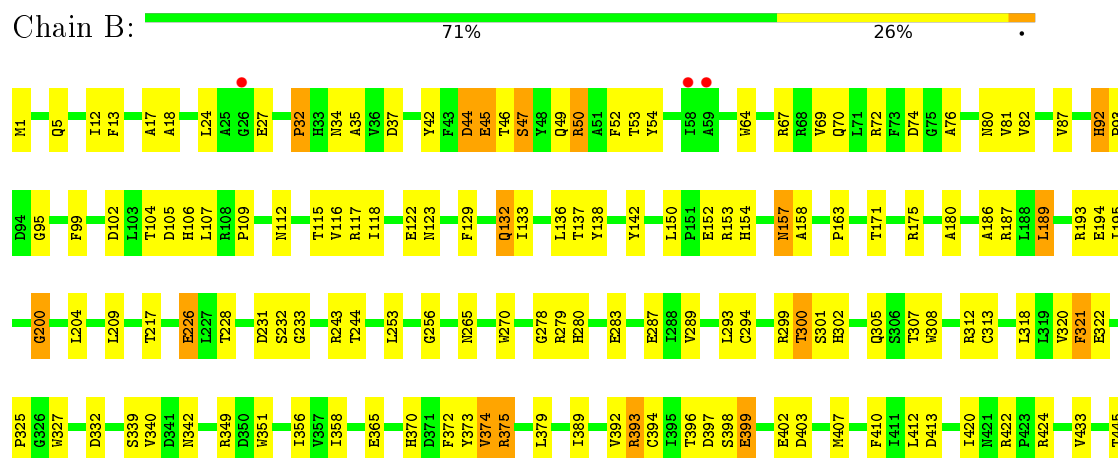
3 Residue-property plots

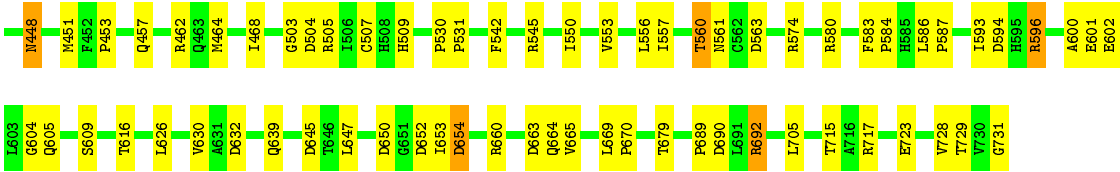
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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-D-galactosidase



• Molecule 2: Beta-D-galactosidase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	79.26Å 107.58Å 201.55Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.27 – 3.15 46.27 – 3.15	Depositor EDS
% Data completeness (in resolution range)	96.7 (46.27-3.15) 96.7 (46.27-3.15)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.91 (at 3.12Å)	Xtriage
Refinement program	REFMAC 5.8.0103	Depositor
R, R_{free}	0.184 , 0.267 0.195 , 0.268	Depositor DCC
R_{free} test set	1480 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	65.9	Xtriage
Anisotropy	0.069	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 62.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	11724	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.76% 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, 0.333 respectively for untwinned datasets, and 0.375, 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: ACT, GLC, PEG, GAL, 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	0.86	0/5919	1.04	0/8084
2	B	0.87	1/5918 (0.0%)	1.03	0/8084
All	All	0.86	1/11837 (0.0%)	1.04	0/16168

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	200	GLY	N-CA	5.59	1.54	1.46

There are no bond angle outliers.

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	5764	0	5591	139	0
2	B	5763	0	5591	122	1
3	A	12	0	12	3	0
3	B	12	0	12	0	0
4	A	12	0	12	0	0
5	A	7	0	10	0	0
5	B	28	0	40	5	0
6	A	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	2	0	0	1	0
7	B	8	0	6	0	0
8	A	66	0	0	4	0
8	B	48	0	0	4	0
All	All	11724	0	11274	259	1

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 11.

All (259) 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:678:VAL:HG23	1:A:715:THR:O	1.07	1.23
1:A:678:VAL:CG2	1:A:715:THR:O	2.02	1.07
1:A:157:ASN:HD22	1:A:158:ALA:H	1.19	0.89
1:A:460:GLU:OE2	1:A:548:ARG:NH1	2.12	0.81
1:A:165:ALA:O	1:A:170:LYS:NZ	2.15	0.79
2:B:107:LEU:HA	2:B:112:ASN:OD1	1.83	0.79
1:A:1:MET:HA	1:A:233:GLY:O	1.83	0.78
1:A:9:HIS:O	1:A:56:ARG:NH1	2.17	0.78
1:A:379:LEU:O	1:A:379:LEU:HD12	1.86	0.75
1:A:517:GLU:HG2	1:A:698:THR:HG21	1.68	0.74
1:A:1:MET:CE	1:A:235:VAL:HB	2.20	0.71
2:B:464:MET:O	2:B:468:ILE:HG12	1.90	0.71
1:A:369:ASN:OD1	1:A:372:PHE:HB3	1.91	0.70
1:A:157:ASN:HD22	1:A:158:ALA:N	1.90	0.70
1:A:601:GLU:HA	1:A:605:GLN:OE1	1.91	0.69
1:A:377:ASN:HD22	1:A:381:ARG:NH1	1.90	0.69
2:B:370:HIS:CD2	5:B:802:PEG:H32	2.28	0.69
1:A:706:THR:OG1	1:A:708:ASP:OD1	2.05	0.68
2:B:393:ARG:NH2	2:B:402:GLU:OE1	2.25	0.68
1:A:221:GLN:NE2	1:A:222:LEU:O	2.27	0.68
2:B:1:MET:N	2:B:233:GLY:O	2.26	0.68
1:A:595:HIS:HA	1:A:598:ILE:O	1.93	0.67
2:B:365:GLU:OE2	2:B:394:CYS:HB3	1.96	0.66
2:B:312:ARG:NH1	8:B:901:HOH:O	2.23	0.66
2:B:32:PRO:HB3	2:B:142:TYR:O	1.96	0.66
1:A:1:MET:HE2	1:A:235:VAL:HB	1.76	0.65
1:A:391:GLY:O	1:A:406:THR:HB	1.97	0.65
2:B:650:ASP:OD2	2:B:652:ASP:N	2.29	0.65
1:A:467:VAL:HG22	1:A:521:ALA:HA	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:157:ASN:ND2	1:A:158:ALA:H	1.93	0.63
1:A:635:PRO:HB3	1:A:721:PHE:CE1	2.34	0.63
1:A:702:LEU:C	1:A:703:LEU:HD12	2.20	0.62
1:A:447:TYR:CD1	1:A:473:VAL:HG11	2.36	0.61
1:A:579:ASP:OD2	1:A:596:ARG:NH2	2.32	0.61
2:B:12:ILE:O	2:B:54:TYR:HA	2.01	0.61
1:A:447:TYR:CE2	1:A:488:GLY:HA2	2.34	0.60
1:A:369:ASN:OD1	1:A:372:PHE:CB	2.50	0.60
2:B:586:LEU:HD23	2:B:587:PRO:HD2	1.84	0.60
2:B:601:GLU:O	2:B:604:GLY:N	2.34	0.60
2:B:115:THR:HG22	2:B:116:VAL:N	2.17	0.60
2:B:42:TYR:CE2	2:B:504:ASP:HB3	2.37	0.59
2:B:49:GLN:NE2	2:B:122:GLU:OE2	2.34	0.58
2:B:300:THR:OG1	2:B:322:GLU:HA	2.03	0.58
2:B:689:PRO:O	2:B:692:ARG:HD3	2.04	0.58
1:A:668:ARG:HG2	8:A:932:HOH:O	2.04	0.58
1:A:279:ARG:O	1:A:283:GLU:HG3	2.04	0.58
1:A:300:THR:HB	1:A:305:GLN:CG	2.33	0.58
1:A:668:ARG:HD2	1:A:696:GLY:O	2.04	0.58
2:B:340:VAL:CG1	2:B:379:LEU:HD22	2.34	0.58
1:A:377:ASN:HD22	1:A:381:ARG:HH12	1.51	0.57
2:B:17:ALA:HB1	2:B:117:ARG:NE	2.19	0.57
1:A:531:PRO:HG3	1:A:561:ASN:HB3	1.87	0.57
1:A:687:ILE:HD11	1:A:704:ARG:HB2	1.86	0.56
2:B:200:GLY:HA3	2:B:204:LEU:CD2	2.35	0.56
2:B:422:ARG:HD3	5:B:805:PEG:O1	2.04	0.56
1:A:362:ARG:HG2	1:A:373:TYR:CZ	2.39	0.56
1:A:601:GLU:CA	1:A:605:GLN:OE1	2.54	0.56
2:B:279:ARG:HD2	2:B:308:TRP:CE3	2.41	0.56
1:A:618:TRP:CD1	1:A:623:GLN:HA	2.41	0.55
1:A:548:ARG:NH2	1:A:555:PRO:O	2.40	0.55
1:A:619:LEU:HA	8:A:904:HOH:O	2.05	0.55
2:B:200:GLY:HA3	2:B:204:LEU:HD21	1.89	0.54
1:A:437:LYS:HA	8:A:933:HOH:O	2.06	0.54
2:B:580:ARG:HD3	8:B:929:HOH:O	2.07	0.54
2:B:542:PHE:CD1	2:B:670:PRO:HG3	2.41	0.54
2:B:340:VAL:HG12	2:B:379:LEU:HD22	1.89	0.54
1:A:365:GLU:OE2	3:A:801:GAL:O1	2.25	0.54
2:B:647:LEU:O	2:B:731:GLY:N	2.40	0.54
1:A:635:PRO:HB3	1:A:721:PHE:CZ	2.42	0.53
1:A:221:GLN:HG3	1:A:223:TYR:CE1	2.43	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:447:TYR:CE2	1:A:488:GLY:CA	2.91	0.53
2:B:76:ALA:O	2:B:95:GLY:HA2	2.08	0.53
1:A:538:GLU:OE1	1:A:627:ARG:NH1	2.42	0.53
1:A:290:ARG:O	1:A:294:CYS:HA	2.09	0.53
1:A:377:ASN:HD21	1:A:402:GLU:HG3	1.72	0.53
1:A:694:LEU:HD23	1:A:699:THR:HG22	1.91	0.53
1:A:284:ARG:O	1:A:285:ASP:C	2.47	0.53
1:A:498:LYS:HD2	2:B:632:ASP:HB3	1.90	0.52
1:A:708:ASP:N	1:A:708:ASP:OD1	2.37	0.52
1:A:156:THR:O	1:A:157:ASN:HB2	2.10	0.52
2:B:299:ARG:NH2	2:B:392:VAL:HG21	2.24	0.52
1:A:608:MET:O	1:A:609:SER:CB	2.57	0.52
2:B:35:ALA:HB1	2:B:52:PHE:CE2	2.45	0.52
1:A:678:VAL:HG22	1:A:679:THR:N	2.25	0.52
2:B:420:ILE:HG21	5:B:805:PEG:H22	1.92	0.52
1:A:12:ILE:O	1:A:54:TYR:HA	2.09	0.51
2:B:123:ASN:O	2:B:129:PHE:HB2	2.09	0.51
2:B:17:ALA:HB1	2:B:117:ARG:HE	1.75	0.51
2:B:189:LEU:HB2	2:B:193:ARG:O	2.10	0.51
2:B:705:LEU:HB3	8:B:909:HOH:O	2.11	0.51
1:A:494:TYR:CE2	1:A:507:CYS:HB2	2.45	0.51
2:B:293:LEU:O	2:B:294:CYS:C	2.48	0.51
2:B:64:TRP:CE3	2:B:69:VAL:HG11	2.46	0.51
2:B:92:HIS:ND1	2:B:99:PHE:HB3	2.24	0.51
2:B:339:SER:O	2:B:340:VAL:C	2.45	0.51
2:B:393:ARG:NH1	2:B:398:SER:OG	2.43	0.51
1:A:157:ASN:ND2	1:A:158:ALA:N	2.56	0.51
1:A:446:GLU:OE2	3:A:801:GAL:H1	2.11	0.50
1:A:645:ASP:HA	1:A:728:VAL:HG12	1.92	0.50
1:A:263:GLY:HA2	1:A:295:CYS:HB3	1.94	0.50
2:B:42:TYR:CE2	2:B:504:ASP:CB	2.95	0.50
2:B:647:LEU:N	2:B:729:THR:O	2.44	0.50
2:B:136:LEU:HB2	2:B:138:TYR:CZ	2.47	0.50
1:A:504:ASP:OD1	1:A:504:ASP:N	2.44	0.50
2:B:507:CYS:HB3	2:B:509:HIS:CE1	2.47	0.50
2:B:265:ASN:OD1	2:B:299:ARG:HD3	2.12	0.50
1:A:205:THR:C	1:A:206:LEU:HG	2.32	0.49
2:B:321:PHE:C	2:B:321:PHE:CD2	2.86	0.49
2:B:195:ILE:HG22	2:B:209:LEU:HD22	1.93	0.49
1:A:712:ILE:O	1:A:727:THR:HA	2.12	0.49
2:B:396:THR:HG22	2:B:433:VAL:HA	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:288:ILE:O	1:A:292:ASP:HB2	2.12	0.49
2:B:47:SER:O	2:B:50:ARG:NH1	2.45	0.49
1:A:248:THR:O	1:A:251:GLY:O	2.30	0.49
1:A:489:TRP:CD2	1:A:490:CYS:HB3	2.47	0.49
1:A:336:GLN:HG2	1:A:372:PHE:CG	2.48	0.49
1:A:336:GLN:HG2	1:A:372:PHE:CD1	2.48	0.49
1:A:545:ARG:NH2	1:A:608:MET:SD	2.85	0.48
1:A:634:LEU:O	1:A:663:ASP:HA	2.12	0.48
1:A:542:PHE:CD1	1:A:670:PRO:HG3	2.48	0.48
1:A:517:GLU:HG2	1:A:698:THR:CG2	2.41	0.48
2:B:132:GLN:C	2:B:133:ILE:HG13	2.34	0.48
1:A:381:ARG:NH2	1:A:403:ASP:OD1	2.45	0.48
2:B:81:VAL:O	2:B:116:VAL:HA	2.14	0.48
1:A:68:ARG:HA	1:A:104:THR:HG21	1.96	0.47
1:A:566:GLU:HB2	1:A:618:TRP:CZ3	2.49	0.47
1:A:1:MET:HE3	1:A:235:VAL:HB	1.94	0.47
2:B:115:THR:CG2	2:B:116:VAL:N	2.77	0.47
2:B:574:ARG:NH2	2:B:602:GLU:OE1	2.48	0.47
2:B:556:LEU:N	2:B:593:ILE:O	2.40	0.47
2:B:278:GLY:N	2:B:654:ASP:OD2	2.36	0.47
1:A:302:HIS:O	1:A:325:PRO:HA	2.15	0.47
2:B:424:ARG:HA	5:B:805:PEG:H41	1.97	0.47
1:A:338:ARG:CZ	1:A:338:ARG:HA	2.45	0.47
2:B:34:ASN:HB3	2:B:37:ASP:OD1	2.15	0.47
1:A:298:VAL:HG22	1:A:299:ARG:N	2.29	0.47
1:A:379:LEU:HD12	1:A:379:LEU:C	2.36	0.47
2:B:321:PHE:C	2:B:321:PHE:HD2	2.18	0.47
2:B:44:ASP:C	2:B:46:THR:H	2.19	0.47
1:A:499:ASP:O	2:B:545:ARG:HD3	2.15	0.46
2:B:370:HIS:O	2:B:374:VAL:HG12	2.15	0.46
1:A:548:ARG:HD3	1:A:552:GLY:O	2.16	0.46
2:B:299:ARG:CZ	2:B:392:VAL:HG21	2.46	0.46
1:A:300:THR:HB	1:A:305:GLN:CD	2.36	0.46
1:A:689:PRO:O	1:A:692:ARG:HD3	2.15	0.46
2:B:630:VAL:CG1	2:B:665:VAL:HG11	2.46	0.46
2:B:560:THR:HG22	2:B:561:ASN:N	2.31	0.46
2:B:457:GLN:NE2	2:B:670:PRO:O	2.49	0.46
1:A:412:LEU:HA	1:A:413:ASP:HA	1.81	0.45
1:A:663:ASP:C	1:A:663:ASP:OD1	2.52	0.45
1:A:515:TRP:HZ3	1:A:702:LEU:HD12	1.81	0.45
2:B:339:SER:O	2:B:342:ASN:N	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:349:ARG:HG2	2:B:349:ARG:O	2.16	0.45
2:B:542:PHE:CG	2:B:670:PRO:HG3	2.52	0.45
1:A:597:HIS:O	1:A:598:ILE:HG13	2.16	0.45
1:A:543:TRP:CG	1:A:629:TYR:HB3	2.52	0.45
1:A:102:ASP:OD1	1:A:104:THR:OG1	2.34	0.45
1:A:59:ALA:HA	1:A:111:ASP:HA	1.99	0.45
2:B:153:ARG:HB3	2:B:232:SER:HB3	1.99	0.45
2:B:645:ASP:O	2:B:728:VAL:HA	2.16	0.45
1:A:391:GLY:O	1:A:406:THR:CB	2.63	0.45
1:A:446:GLU:HA	1:A:487:ILE:O	2.17	0.45
1:A:407:MET:HG3	1:A:408:ASN:N	2.32	0.45
2:B:300:THR:HB	2:B:305:GLN:HG3	1.99	0.45
2:B:424:ARG:HA	5:B:805:PEG:C3	2.47	0.45
2:B:586:LEU:HD23	2:B:587:PRO:CD	2.47	0.44
1:A:369:ASN:O	1:A:369:ASN:OD1	2.36	0.44
1:A:369:ASN:O	1:A:370:HIS:C	2.55	0.44
1:A:447:TYR:CE1	1:A:473:VAL:HG11	2.52	0.44
1:A:250:GLN:HA	1:A:250:GLN:OE1	2.16	0.44
2:B:586:LEU:CD2	2:B:587:PRO:HD2	2.48	0.44
2:B:18:ALA:O	2:B:117:ARG:NH2	2.50	0.44
2:B:118:ILE:HG13	2:B:118:ILE:O	2.17	0.44
2:B:503:GLY:N	8:B:906:HOH:O	2.51	0.44
1:A:55:GLN:NE2	1:A:113:LEU:HD22	2.33	0.44
1:A:498:LYS:HG3	1:A:499:ASP:N	2.33	0.44
1:A:650:ASP:HB3	1:A:653:ILE:HD11	2.00	0.44
2:B:372:PHE:O	2:B:375:ARG:HB2	2.18	0.44
2:B:12:ILE:HG22	2:B:24:LEU:HD23	1.99	0.43
2:B:243:ARG:NH1	6:B:806:CL:CL	2.88	0.43
1:A:80:ASN:C	1:A:81:VAL:CG2	2.86	0.43
2:B:129:PHE:CE1	2:B:137:THR:HG21	2.53	0.43
2:B:253:LEU:HG	2:B:256:GLY:O	2.18	0.43
1:A:447:TYR:CD2	1:A:488:GLY:HA2	2.52	0.43
2:B:186:ALA:HB2	2:B:204:LEU:HD13	1.99	0.43
1:A:288:ILE:O	1:A:293:LEU:HD12	2.18	0.43
1:A:320:VAL:HB	1:A:356:ILE:HA	2.01	0.43
1:A:364:ASN:ND2	3:A:801:GAL:O2	2.52	0.43
1:A:430:THR:HG21	1:A:482:ALA:HB1	2.00	0.43
2:B:373:TYR:O	2:B:374:VAL:C	2.56	0.43
2:B:301:SER:HA	2:B:302:HIS:HA	1.85	0.43
2:B:187:ARG:NH1	2:B:226:GLU:OE1	2.51	0.43
2:B:410:PHE:HB3	2:B:451:MET:HE2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:468:ILE:HD11	2:B:557:ILE:HD11	2.00	0.43
2:B:530:PRO:O	2:B:531:PRO:C	2.57	0.43
1:A:529:LYS:O	1:A:561:ASN:ND2	2.52	0.43
2:B:105:ASP:N	2:B:105:ASP:OD1	2.52	0.43
2:B:163:PRO:HG2	2:B:171:THR:HB	2.01	0.43
2:B:13:PHE:HA	2:B:53:THR:O	2.19	0.43
1:A:284:ARG:O	1:A:287:GLU:N	2.52	0.43
2:B:302:HIS:O	2:B:325:PRO:HA	2.19	0.42
2:B:412:LEU:HA	2:B:413:ASP:HA	1.83	0.42
2:B:42:TYR:CZ	2:B:504:ASP:HB3	2.53	0.42
1:A:25:ALA:HB2	8:A:957:HOH:O	2.20	0.42
1:A:311:ASP:O	1:A:314:ASP:HB2	2.19	0.42
2:B:81:VAL:HG12	2:B:82:VAL:N	2.34	0.42
1:A:216:SER:OG	1:A:217:THR:N	2.52	0.42
1:A:1:MET:CG	1:A:2:ARG:H	2.33	0.42
1:A:362:ARG:HG2	1:A:373:TYR:CE1	2.55	0.42
2:B:321:PHE:CD1	2:B:358:ILE:CG2	3.03	0.42
1:A:122:GLU:OE2	1:A:497:HIS:NE2	2.51	0.42
1:A:324:ILE:HB	1:A:325:PRO:CD	2.50	0.42
1:A:39:PRO:HB2	1:A:43:PHE:HD2	1.84	0.42
1:A:57:VAL:HG13	1:A:111:ASP:HB2	2.01	0.42
2:B:70:GLN:NE2	2:B:102:ASP:OD2	2.52	0.42
1:A:514:ILE:O	1:A:514:ILE:CG1	2.67	0.42
1:A:616:THR:HA	1:A:625:ALA:O	2.20	0.42
2:B:150:LEU:HD23	2:B:154:HIS:CG	2.55	0.42
1:A:151:PRO:O	1:A:154:HIS:ND1	2.52	0.42
2:B:407:MET:HG2	2:B:433:VAL:HG11	2.02	0.42
2:B:453:PRO:O	2:B:462:ARG:NH2	2.53	0.42
2:B:553:VAL:HG23	2:B:553:VAL:O	2.18	0.42
2:B:663:ASP:O	2:B:664:GLN:C	2.58	0.42
1:A:228:THR:HA	1:A:234:ASP:O	2.20	0.41
1:A:299:ARG:NH2	1:A:392:VAL:HG21	2.35	0.41
2:B:583:PHE:O	2:B:584:PRO:C	2.58	0.41
2:B:653:ILE:HB	2:B:705:LEU:HD12	2.01	0.41
1:A:500:PHE:N	1:A:500:PHE:CD2	2.88	0.41
2:B:320:VAL:HB	2:B:356:ILE:HA	2.02	0.41
2:B:503:GLY:O	2:B:505:ARG:NH1	2.52	0.41
2:B:594:ASP:OD2	2:B:596:ARG:NH2	2.53	0.41
2:B:283:GLU:O	2:B:287:GLU:HG3	2.20	0.41
2:B:370:HIS:ND1	2:B:399:GLU:OE2	2.53	0.41
1:A:454:THR:HG22	1:A:454:THR:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:515:TRP:O	1:A:516:ARG:HB2	2.20	0.41
2:B:106:HIS:O	2:B:107:LEU:HD23	2.20	0.41
2:B:122:GLU:HA	2:B:129:PHE:CD1	2.56	0.41
2:B:157:ASN:HD22	2:B:158:ALA:H	1.69	0.41
1:A:170:LYS:HB2	1:A:209:LEU:HB2	2.02	0.41
1:A:265:ASN:ND2	1:A:487:ILE:HG22	2.35	0.41
1:A:174:ILE:HD12	1:A:174:ILE:N	2.36	0.41
1:A:320:VAL:HG12	1:A:321:PHE:N	2.36	0.41
1:A:300:THR:HB	1:A:305:GLN:HG3	2.01	0.41
2:B:153:ARG:NH2	2:B:180:ALA:O	2.54	0.41
1:A:134:ASP:OD1	1:A:494:TYR:OH	2.23	0.40
1:A:328:GLN:N	1:A:364:ASN:O	2.54	0.40
1:A:60:TRP:HZ2	1:A:65:GLN:HG3	1.86	0.40
2:B:313:CYS:SG	2:B:318:LEU:HD23	2.60	0.40
2:B:321:PHE:CD1	2:B:358:ILE:HG21	2.56	0.40
1:A:612:PRO:HB3	1:A:630:VAL:HA	2.03	0.40
1:A:254:LEU:O	1:A:256:GLY:N	2.54	0.40
1:A:280:HIS:O	1:A:283:GLU:N	2.54	0.40
1:A:372:PHE:O	1:A:375:ARG:HB2	2.21	0.40
1:A:447:TYR:CE2	1:A:488:GLY:N	2.90	0.40
2:B:279:ARG:O	2:B:280:HIS:C	2.58	0.40
2:B:389:ILE:H	2:B:403:ASP:HB2	1.85	0.40
2:B:639:GLN:HB3	2:B:660:ARG:HB2	2.02	0.40
2:B:321:PHE:CG	2:B:358:ILE:HG23	2.57	0.40
2:B:583:PHE:HB3	2:B:586:LEU:HD12	2.04	0.40

All (1) 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	Interatomic distance (Å)	Clash overlap (Å)
2:B:175:ARG:NH1	2:B:679:THR:OG1[3_745]	2.01	0.19

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	729/732 (100%)	621 (85%)	100 (14%)	8 (1%)	17	61
2	B	729/731 (100%)	631 (87%)	84 (12%)	14 (2%)	10	48
All	All	1458/1463 (100%)	1252 (86%)	184 (13%)	22 (2%)	13	53

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	609	SER
2	B	104	THR
2	B	332	ASP
1	A	532	SER
1	A	608	MET
2	B	550	ILE
1	A	152	GLU
1	A	623	GLN
2	B	669	LEU
1	A	605	GLN
2	B	45	GLU
2	B	399	GLU
2	B	600	ALA
2	B	109	PRO
2	B	375	ARG
2	B	448	ASN
1	A	550	ILE
2	B	300	THR
2	B	93	PRO
1	A	23	PRO
2	B	374	VAL
2	B	289	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	611/612 (100%)	561 (92%)	50 (8%)	14	48
2	B	611/611 (100%)	566 (93%)	45 (7%)	17	54
All	All	1222/1223 (100%)	1127 (92%)	95 (8%)	16	51

All (95) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	7	LEU
1	A	29	VAL
1	A	34	ASN
1	A	37	ASP
1	A	43	PHE
1	A	45	GLU
1	A	47	SER
1	A	58	ILE
1	A	80	ASN
1	A	104	THR
1	A	117	ARG
1	A	119	ASP
1	A	121	SER
1	A	143	ARG
1	A	157	ASN
1	A	167	SER
1	A	213	SER
1	A	216	SER
1	A	248	THR
1	A	250	GLN
1	A	253	LEU
1	A	293	LEU
1	A	294	CYS
1	A	303	TYR
1	A	305	GLN
1	A	321	PHE
1	A	322	GLU
1	A	323	GLU
1	A	336	GLN
1	A	351	TRP
1	A	371	ASP
1	A	406	THR
1	A	419	LEU
1	A	421	ASN

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Mol	Chain	Res	Type
1	A	434	THR
1	A	436	ILE
1	A	445	THR
1	A	490	CYS
1	A	530	PRO
1	A	535	ILE
1	A	578	PRO
1	A	599	SER
1	A	605	GLN
1	A	639	GLN
1	A	664	GLN
1	A	699	THR
1	A	708	ASP
1	A	713	ARG
1	A	725	VAL
2	B	5	GLN
2	B	27	GLU
2	B	32	PRO
2	B	44	ASP
2	B	45	GLU
2	B	47	SER
2	B	50	ARG
2	B	67	ARG
2	B	72	ARG
2	B	74	ASP
2	B	80	ASN
2	B	87	VAL
2	B	92	HIS
2	B	132	GLN
2	B	152	GLU
2	B	157	ASN
2	B	189	LEU
2	B	194	GLU
2	B	217	THR
2	B	226	GLU
2	B	228	THR
2	B	231	ASP
2	B	244	THR
2	B	270	TRP
2	B	307	THR
2	B	321	PHE
2	B	327	TRP

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Mol	Chain	Res	Type
2	B	351	TRP
2	B	393	ARG
2	B	397	ASP
2	B	445	THR
2	B	448	ASN
2	B	560	THR
2	B	563	ASP
2	B	596	ARG
2	B	605	GLN
2	B	609	SER
2	B	616	THR
2	B	626	LEU
2	B	654	ASP
2	B	690	ASP
2	B	692	ARG
2	B	715	THR
2	B	717	ARG
2	B	723	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	ASN
1	A	123	ASN
1	A	157	ASN
1	A	221	GLN
1	A	265	ASN
1	A	328	GLN
1	A	342	ASN
1	A	377	ASN
1	A	408	ASN
1	A	720	GLN
2	B	157	ASN
2	B	282	GLN
2	B	509	HIS
2	B	611	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 14 ligands modelled in this entry, 4 are monoatomic - leaving 10 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	GAL	A	801	-	12,12,12	0.96	0	17,17,17	1.81	4 (23%)
4	GLC	A	802	-	12,12,12	1.07	0	17,17,17	1.72	4 (23%)
5	PEG	A	803	-	6,6,6	0.57	0	5,5,5	0.16	0
3	GAL	B	801	-	12,12,12	1.41	3 (25%)	17,17,17	2.78	8 (47%)
5	PEG	B	802	-	6,6,6	0.55	0	5,5,5	0.43	0
5	PEG	B	803	-	6,6,6	0.70	0	5,5,5	0.91	0
5	PEG	B	804	-	6,6,6	0.42	0	5,5,5	0.44	0
5	PEG	B	805	-	6,6,6	0.47	0	5,5,5	0.98	0
7	ACT	B	808	-	0,3,3	0.00	-	0,3,3	0.00	-
7	ACT	B	809	-	0,3,3	0.00	-	0,3,3	0.00	-

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	GAL	A	801	-	-	0/2/22/22	0/1/1/1
4	GLC	A	802	-	-	0/2/22/22	0/1/1/1
5	PEG	A	803	-	-	0/4/4/4	0/0/0/0
3	GAL	B	801	-	-	0/2/22/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	PEG	B	802	-	-	0/4/4/4	0/0/0/0
5	PEG	B	803	-	-	0/4/4/4	0/0/0/0
5	PEG	B	804	-	-	0/4/4/4	0/0/0/0
5	PEG	B	805	-	-	0/4/4/4	0/0/0/0
7	ACT	B	808	-	-	0/0/0/0	0/0/0/0
7	ACT	B	809	-	-	0/0/0/0	0/0/0/0

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	801	GAL	O5-C1	2.06	1.47	1.43
3	B	801	GAL	C4-C5	2.31	1.58	1.53
3	B	801	GAL	C6-C5	2.35	1.60	1.51

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	801	GAL	O5-C5-C4	-2.50	104.89	109.67
3	B	801	GAL	O2-C2-C3	-2.31	105.14	110.36
3	B	801	GAL	O4-C4-C3	-2.27	105.25	110.36
4	A	802	GLC	C1-C2-C3	2.13	114.17	110.68
4	A	802	GLC	O5-C5-C6	2.63	113.20	106.38
3	A	801	GAL	O1-C1-C2	2.66	116.41	109.05
3	A	801	GAL	O5-C5-C6	2.73	113.45	106.38
4	A	802	GLC	O5-C1-C2	2.81	114.92	110.00
3	B	801	GAL	O2-C2-C1	2.94	116.18	109.74
3	B	801	GAL	O5-C1-C2	3.51	116.16	110.00
3	B	801	GAL	O6-C6-C5	3.70	123.67	111.30
4	A	802	GLC	C1-O5-C5	4.33	121.82	113.54
3	B	801	GAL	O5-C5-C6	4.51	118.05	106.38
3	A	801	GAL	O3-C3-C2	4.87	121.33	110.36
3	B	801	GAL	C3-C4-C5	5.23	119.55	110.23
3	B	801	GAL	C1-O5-C5	5.68	124.41	113.54

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	801	GAL	3	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	802	PEG	1	0
5	B	805	PEG	4	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	731/732 (99%)	-0.37	5 (0%)	89 82	31, 58, 84, 171	0
2	B	731/731 (100%)	-0.36	3 (0%)	93 90	36, 63, 89, 145	0
All	All	1462/1463 (99%)	-0.37	8 (0%)	91 87	31, 60, 87, 171	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	602	GLU	3.0
2	B	59	ALA	2.8
2	B	58	ILE	2.8
1	A	60	TRP	2.7
1	A	605	GLN	2.6
1	A	62	ASP	2.6
1	A	22	THR	2.0
2	B	26	GLY	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
5	PEG	B	805	7/7	0.79	0.69	4.89	113,117,121,122	0
5	PEG	A	803	7/7	0.85	0.27	3.16	54,66,80,82	0
7	ACT	B	808	4/4	0.93	0.32	2.57	65,66,70,71	0
5	PEG	B	802	7/7	0.86	0.25	2.50	59,74,80,80	0
3	GAL	A	801	12/12	0.93	0.21	1.59	66,75,79,79	0
4	GLC	A	802	12/12	0.82	0.28	0.74	72,91,102,103	0
3	GAL	B	801	12/12	0.94	0.17	0.42	49,56,59,59	0
6	CL	B	806	1/1	0.85	0.14	-0.42	42,42,42,42	0
5	PEG	B	804	7/7	0.92	0.17	-0.45	67,70,72,72	0
7	ACT	B	809	4/4	0.92	0.22	-	47,54,55,55	0
5	PEG	B	803	7/7	0.81	0.34	-	67,71,80,90	0
6	CL	A	805	1/1	0.93	0.24	-	61,61,61,61	0
6	CL	B	807	1/1	0.84	0.21	-	75,75,75,75	0
6	CL	A	804	1/1	0.86	0.21	-	57,57,57,57	0

6.5 Other polymers [i](#)

There are no such residues in this entry.