



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:45 PM UTC

PDB ID : 1HJU / pdb_00001hju
Title : Structure of two fungal beta-1,4-galactanases: searching for the basis for temperature and pH optimum.
Authors : Le Nours, J.; Ryttersgaard, C.; Lo Leggio, L.; Ostergaard, P.R.; Borchert, T.V.; Christensen, L.L.H.; Larsen, S.
Deposited on : 2003-02-27
Resolution : 2.15 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

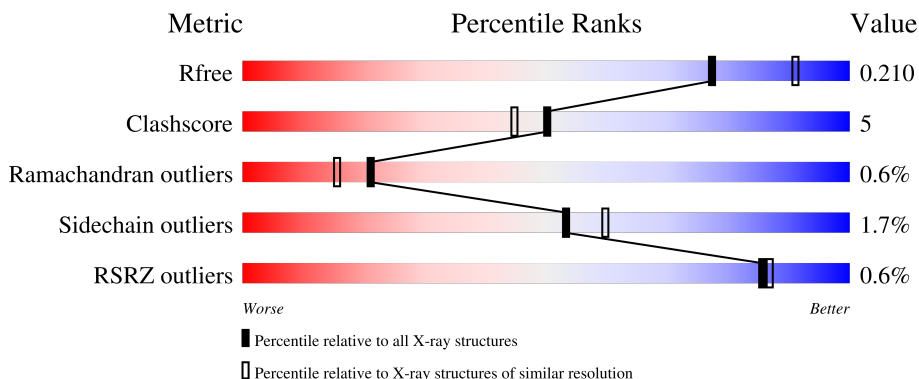
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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}	180053	2057 (2.16-2.16)
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	332	% 86% 14%
1	B	332	% 86% 13%
1	C	332	 83% 16%
1	D	332	% 85% 14%

2 Entry composition [i](#)

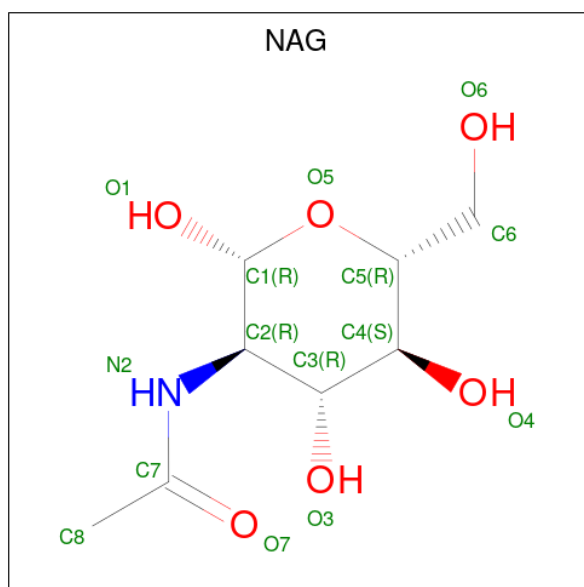
There are 6 unique types of molecules in this entry. The entry contains 11439 atoms, of which 0 are hydrogens and 0 are deuteriums.

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-1,4-GALACTANASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	332	Total	C	N	O	S	0	4	0
			2625	1672	448	497	8			
1	B	332	Total	C	N	O	S	0	1	0
			2609	1661	446	494	8			
1	C	332	Total	C	N	O	S	0	1	0
			2609	1663	445	493	8			
1	D	332	Total	C	N	O	S	0	2	0
			2614	1666	446	494	8			

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



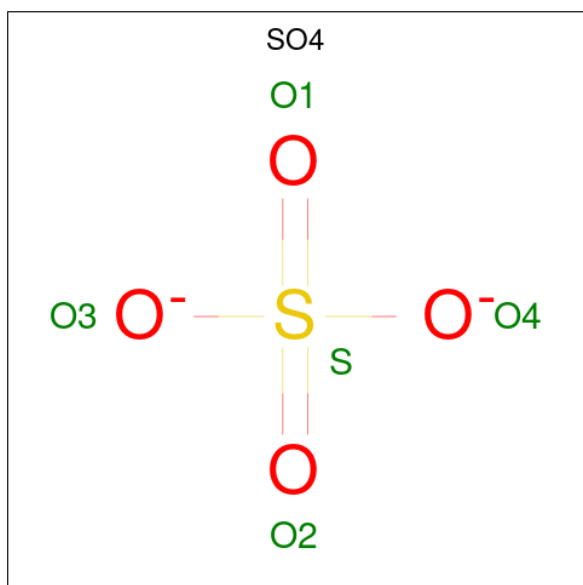
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	C	1	Total	C	N	O	0	0
			14	8	1	5		
2	D	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



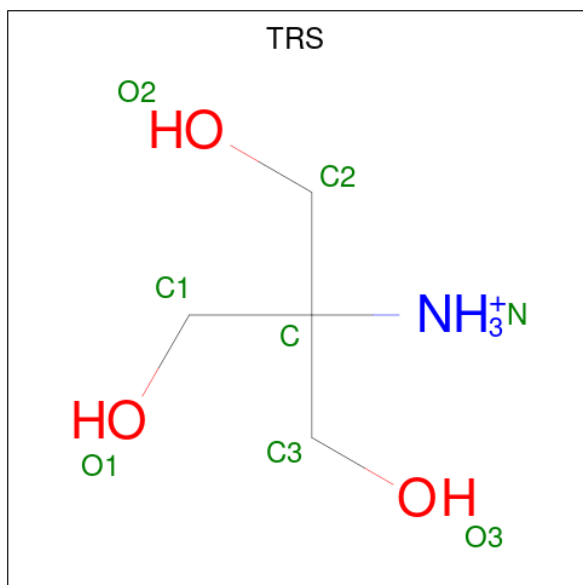
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total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

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

- Molecule 4 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (CCD ID: TRS) (formula: $C_4H_{12}NO_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			8	4	1	3		
4	D	1	Total	C	N	O	0	0
			8	4	1	3		

- Molecule 5 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



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

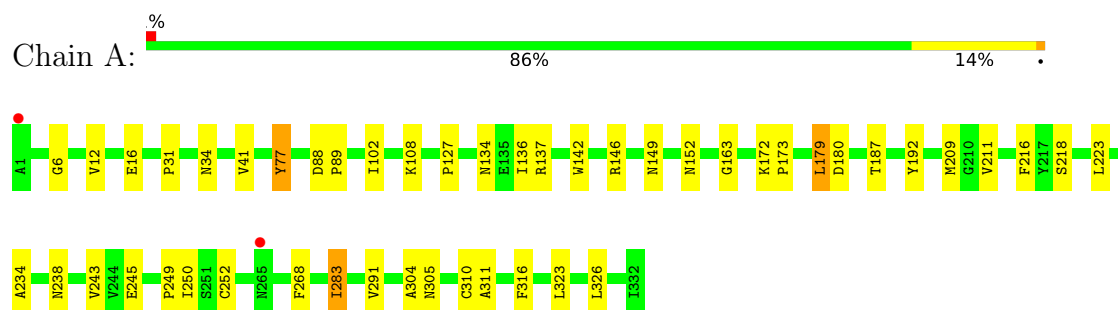
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	237	Total	O	0	0
			237	237		
6	B	206	Total	O	0	0
			206	206		
6	C	202	Total	O	0	1
			202	202		
6	D	187	Total	O	0	0
			187	187		

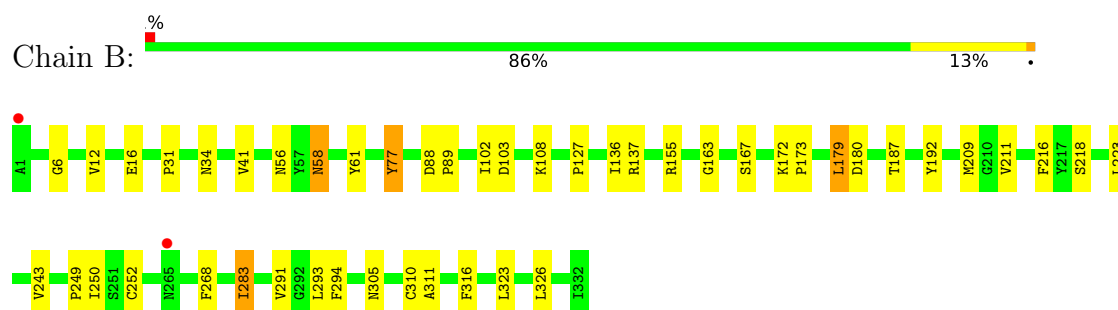
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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.

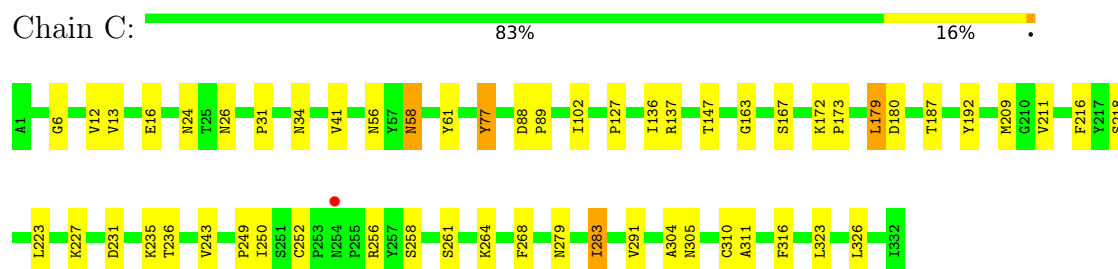
• Molecule 1: BETA-1,4-GALACTANASE



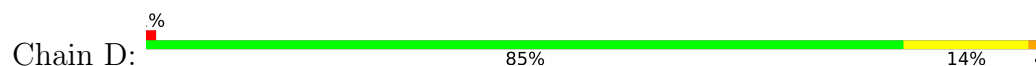
• Molecule 1: BETA-1,4-GALACTANASE

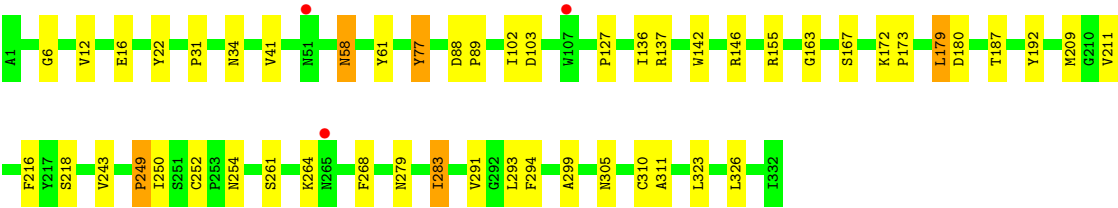


• Molecule 1: BETA-1,4-GALACTANASE



• Molecule 1: BETA-1,4-GALACTANASE





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	107.99Å 136.59Å 110.88Å 90.00° 118.92° 90.00°	Depositor
Resolution (Å)	18.01 – 2.15 18.01 – 2.15	Depositor EDS
% Data completeness (in resolution range)	98.2 (18.01-2.15) 98.0 (18.01-2.15)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	8.47 (at 2.15Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.187 , 0.212 0.187 , 0.210	Depositor DCC
R_{free} test set	7713 reflections (10.09%)	wwPDB-VP
Wilson B-factor (Å ²)	22.2	Xtriage
Anisotropy	0.570	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 34.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.014 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11439	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.93% 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: PEG, SO4, NAG, TRS

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	0.41	0/2710	0.92	11/3707 (0.3%)
1	B	0.38	0/2688	0.93	10/3678 (0.3%)
1	C	0.38	0/2688	0.93	13/3677 (0.4%)
1	D	0.37	0/2696	0.91	11/3689 (0.3%)
All	All	0.39	0/10782	0.92	45/14751 (0.3%)

There are no bond length outliers.

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	250	ILE	N-CA-C	-10.13	102.80	112.83
1	D	250	ILE	N-CA-C	-10.08	102.85	112.83
1	C	250	ILE	N-CA-C	-9.98	102.94	112.83
1	A	250	ILE	N-CA-C	-9.92	103.01	112.83
1	D	291	VAL	N-CA-C	7.60	121.14	113.47
1	B	291	VAL	N-CA-C	7.53	121.07	113.47
1	A	291	VAL	N-CA-C	7.45	121.00	113.47
1	C	291	VAL	N-CA-C	7.35	120.89	113.47
1	C	311	ALA	N-CA-C	6.50	118.44	111.36
1	C	252	CYS	CA-C-N	6.42	125.64	118.97
1	C	252	CYS	C-N-CA	6.42	125.64	118.97
1	C	304	ALA	N-CA-C	6.35	119.01	111.33
1	B	311	ALA	N-CA-C	6.33	118.26	111.36
1	A	304	ALA	N-CA-C	6.29	118.95	111.33
1	A	311	ALA	N-CA-C	6.17	118.08	111.36
1	D	311	ALA	N-CA-C	6.13	118.04	111.36
1	B	283	ILE	N-CA-C	-5.97	104.69	110.72
1	A	252	CYS	CA-C-N	5.93	125.40	119.24
1	A	252	CYS	C-N-CA	5.93	125.40	119.24
1	C	102	ILE	N-CA-C	5.91	116.57	110.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	102	ILE	N-CA-C	5.91	116.56	110.36
1	C	216	PHE	N-CA-C	5.82	119.95	112.26
1	D	252	CYS	CA-C-N	5.82	125.29	119.24
1	D	252	CYS	C-N-CA	5.82	125.29	119.24
1	A	102	ILE	N-CA-C	5.78	116.43	110.36
1	B	102	ILE	N-CA-C	5.77	116.42	110.36
1	C	147	THR	N-CA-C	-5.75	102.40	110.50
1	A	216	PHE	N-CA-C	5.70	119.79	112.26
1	B	216	PHE	N-CA-C	5.66	119.73	112.26
1	B	252	CYS	CA-C-N	5.63	125.09	119.24
1	B	252	CYS	C-N-CA	5.63	125.09	119.24
1	D	216	PHE	N-CA-C	5.62	119.67	112.26
1	C	283[A]	ILE	N-CA-C	-5.57	105.10	110.72
1	C	283[B]	ILE	N-CA-C	-5.57	105.10	110.72
1	D	163	GLY	N-CA-C	-5.51	105.99	113.37
1	D	283	ILE	N-CA-C	-5.47	105.19	110.72
1	B	163	GLY	N-CA-C	-5.46	106.05	113.37
1	B	243	VAL	N-CA-C	-5.38	99.33	107.73
1	C	163	GLY	N-CA-C	-5.38	106.16	113.37
1	A	283	ILE	N-CA-C	-5.28	105.39	110.72
1	A	163	GLY	N-CA-C	-5.19	106.42	113.37
1	C	243	VAL	N-CA-C	-5.19	99.64	107.73
1	D	243	VAL	N-CA-C	-5.12	99.74	107.73
1	A	243	VAL	N-CA-C	-5.09	99.79	107.73
1	D	249	PRO	N-CA-C	5.05	119.96	111.32

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	2625	0	2522	25	0
1	B	2609	0	2501	29	0
1	C	2609	0	2506	30	0
1	D	2614	0	2512	25	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	14	0	13	0	0
2	B	14	0	13	0	0
2	C	14	0	13	0	0
2	D	14	0	13	0	0
3	A	15	0	0	0	0
3	B	10	0	0	0	0
3	C	10	0	0	0	0
3	D	15	0	0	0	0
4	A	8	0	12	3	0
4	D	8	0	12	0	0
5	A	7	0	10	0	0
5	B	7	0	10	0	0
5	C	7	0	10	0	0
5	D	7	0	10	0	0
6	A	237	0	0	2	0
6	B	206	0	0	1	0
6	C	202	0	0	2	0
6	D	187	0	0	1	0
All	All	11439	0	10157	110	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 (110) 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:149[A]:ASN:ND2	6:A:2128:HOH:O	1.88	1.06
1:C:167:SER:O	1:C:172:LYS:NZ	2.15	0.80
1:B:167:SER:O	1:B:172:LYS:NZ	2.16	0.78
1:D:167:SER:O	1:D:172:LYS:NZ	2.17	0.77
1:C:223:LEU:HB3	1:C:283[B]:ILE:HD11	1.68	0.75
1:A:223:LEU:HB3	1:A:283:ILE:HD11	1.69	0.74
1:B:223:LEU:HB3	1:B:283:ILE:HD11	1.73	0.71
1:C:58:ASN:C	1:C:58:ASN:HD22	2.01	0.69
1:B:293:LEU:HD23	1:B:294:PHE:N	2.08	0.68
1:D:58:ASN:HD22	1:D:58:ASN:C	2.02	0.67
1:B:293:LEU:HD23	1:B:293:LEU:C	2.20	0.67
1:A:134:ASN:HD21	4:A:704:TRS:H12	1.61	0.66
1:B:58:ASN:C	1:B:58:ASN:HD22	2.02	0.65
1:A:149[B]:ASN:ND2	1:A:152:ASN:CG	2.55	0.65
1:D:58:ASN:ND2	1:D:61:TYR:H	1.94	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:137:ARG:HD3	1:B:180:ASP:OD1	1.97	0.64
1:C:58:ASN:ND2	1:C:61:TYR:H	1.95	0.64
1:C:137:ARG:HD3	1:C:180:ASP:OD1	1.97	0.64
1:B:58:ASN:ND2	1:B:61:TYR:H	1.96	0.64
1:C:261:SER:HA	1:C:264:LYS:HE3	1.80	0.64
1:D:137:ARG:HD3	1:D:180:ASP:OD1	1.97	0.63
1:A:137:ARG:HD3	1:A:180:ASP:OD1	1.99	0.63
1:A:245:GLU:OE2	4:A:704:TRS:O1	2.18	0.61
1:A:149[B]:ASN:ND2	1:A:152:ASN:ND2	2.49	0.61
1:B:103:ASP:OD1	1:B:155:ARG:NH2	2.31	0.59
1:A:172:LYS:HG3	1:A:173:PRO:HD2	1.85	0.58
4:A:704:TRS:H11	6:A:2083:HOH:O	2.05	0.57
1:D:293:LEU:C	1:D:293:LEU:HD23	2.33	0.54
1:A:223:LEU:HB3	1:A:283:ILE:CD1	2.37	0.54
1:D:249:PRO:HG2	1:D:268:PHE:CD1	2.43	0.54
1:A:249:PRO:HG2	1:A:268:PHE:CD1	2.43	0.53
1:C:279:ASN:O	1:C:283[A]:ILE:HG12	2.08	0.53
1:C:249:PRO:HG2	1:C:268:PHE:CD1	2.43	0.53
1:C:223:LEU:HB3	1:C:283[B]:ILE:CD1	2.37	0.53
1:B:249:PRO:HG2	1:B:268:PHE:CD1	2.43	0.52
1:B:88:ASP:HB2	1:B:89:PRO:CD	2.39	0.52
1:C:256:ARG:NH1	6:C:2174[A]:HOH:O	2.43	0.52
1:C:192:TYR:CE1	1:C:209:MET:HE2	2.45	0.52
1:A:142:TRP:CE2	1:A:146:ARG:HD3	2.44	0.51
1:C:88:ASP:HB2	1:C:89:PRO:CD	2.41	0.51
1:D:12:VAL:O	1:D:16:GLU:HG3	2.12	0.50
1:C:137:ARG:HD2	1:C:187:THR:HG21	1.94	0.50
1:D:279:ASN:O	1:D:283:ILE:HG12	2.12	0.50
1:B:58:ASN:C	1:B:58:ASN:ND2	2.68	0.49
1:B:192:TYR:CE1	1:B:209:MET:HE2	2.47	0.49
1:C:24:ASN:OD1	1:C:26:ASN:HB2	2.12	0.49
1:D:88:ASP:HB2	1:D:89:PRO:CD	2.42	0.49
1:A:137:ARG:HD2	1:A:187:THR:HG21	1.94	0.49
1:A:192:TYR:CE1	1:A:209:MET:HE2	2.47	0.49
1:C:305:ASN:HB2	1:C:310:CYS:O	2.12	0.49
1:B:137:ARG:HD2	1:B:187:THR:HG21	1.94	0.49
1:C:12:VAL:O	1:C:16:GLU:HG3	2.11	0.49
1:B:223:LEU:HB3	1:B:283:ILE:CD1	2.41	0.49
1:D:192:TYR:CE1	1:D:209:MET:HE2	2.47	0.49
1:D:261:SER:HA	1:D:264:LYS:HE3	1.93	0.49
1:A:88:ASP:HB2	1:A:89:PRO:CD	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:77:TYR:CD1	1:B:77:TYR:C	2.91	0.49
1:B:305:ASN:HB2	1:B:310:CYS:O	2.13	0.49
1:C:77:TYR:C	1:C:77:TYR:CD1	2.90	0.49
1:D:77:TYR:C	1:D:77:TYR:CD1	2.91	0.49
1:B:6:GLY:HA2	1:B:41:VAL:HG13	1.95	0.48
1:D:137:ARG:HD2	1:D:187:THR:HG21	1.95	0.48
1:D:127:PRO:O	1:D:173:PRO:HB3	2.13	0.48
1:C:31:PRO:HB2	1:C:34:ASN:ND2	2.28	0.48
1:D:305:ASN:HB2	1:D:310:CYS:O	2.13	0.48
1:A:77:TYR:C	1:A:77:TYR:CD1	2.92	0.48
1:D:31:PRO:HB2	1:D:34:ASN:ND2	2.29	0.48
1:A:6:GLY:HA2	1:A:41:VAL:HG13	1.96	0.48
1:A:305:ASN:HB2	1:A:310:CYS:O	2.14	0.47
1:B:31:PRO:HD2	1:B:34:ASN:HD22	1.79	0.47
1:B:293:LEU:C	1:B:293:LEU:CD2	2.86	0.47
1:D:142:TRP:CE2	1:D:146:ARG:HD3	2.49	0.47
1:B:12:VAL:O	1:B:16:GLU:HG3	2.13	0.47
1:A:12:VAL:O	1:A:16:GLU:HG3	2.14	0.47
1:A:31:PRO:HD2	1:A:34:ASN:HD22	1.80	0.47
1:B:316:PHE:HA	1:B:323:LEU:HD13	1.97	0.46
1:D:6:GLY:HA2	1:D:41:VAL:HG13	1.96	0.46
1:A:31:PRO:HB2	1:A:34:ASN:ND2	2.29	0.46
1:B:127:PRO:O	1:B:173:PRO:HB3	2.15	0.46
1:C:6:GLY:HA2	1:C:41:VAL:HG13	1.97	0.46
1:B:31:PRO:HB2	1:B:34:ASN:ND2	2.30	0.46
1:D:58:ASN:C	1:D:58:ASN:ND2	2.68	0.46
1:A:234:ALA:O	1:A:238[A]:ASN:ND2	2.49	0.46
1:A:127:PRO:O	1:A:173:PRO:HB3	2.16	0.45
1:D:31:PRO:HD2	1:D:34:ASN:HD22	1.81	0.45
1:C:31:PRO:HD2	1:C:34:ASN:HD22	1.81	0.45
1:B:179:LEU:O	1:B:211:VAL:HA	2.17	0.44
1:D:293:LEU:HD23	1:D:294:PHE:N	2.32	0.44
1:C:127:PRO:O	1:C:173:PRO:HB3	2.16	0.44
1:C:316:PHE:HA	1:C:323:LEU:HD13	1.99	0.43
1:A:179:LEU:O	1:A:211:VAL:HA	2.18	0.43
1:C:179:LEU:O	1:C:211:VAL:HA	2.19	0.43
1:D:103:ASP:OD1	1:D:155:ARG:NH2	2.45	0.43
1:D:179:LEU:O	1:D:211:VAL:HA	2.19	0.43
1:D:254[A]:ASN:OD1	6:D:2145:HOH:O	2.21	0.43
1:C:58:ASN:C	1:C:58:ASN:ND2	2.67	0.42
1:B:316:PHE:CA	1:B:323:LEU:HD13	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:58:ASN:HD21	1:B:61:TYR:H	1.67	0.42
1:B:77:TYR:C	1:B:77:TYR:HD1	2.28	0.41
1:C:58:ASN:HD21	1:C:61:TYR:H	1.67	0.41
1:C:227:LYS:HE3	1:C:231:ASP:OD2	2.20	0.41
1:C:12:VAL:HG13	1:C:13:VAL:N	2.35	0.41
1:A:316:PHE:HA	1:A:323[B]:LEU:HD13	2.01	0.41
1:C:77:TYR:C	1:C:77:TYR:HD1	2.28	0.41
1:A:149[B]:ASN:HD21	1:A:152:ASN:CG	2.28	0.41
1:C:235:LYS:HG3	1:C:236:THR:N	2.36	0.41
1:B:56:ASN:HB2	6:B:2052:HOH:O	2.21	0.40
1:B:88:ASP:HB2	1:B:89:PRO:HD2	2.04	0.40
1:C:56:ASN:HB2	6:C:2056:HOH:O	2.21	0.40
1:D:22:TYR:CE1	1:D:299:ALA:HB1	2.56	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	334/332 (101%)	321 (96%)	11 (3%)	2 (1%)	21	15
1	B	331/332 (100%)	316 (96%)	13 (4%)	2 (1%)	21	15
1	C	331/332 (100%)	318 (96%)	11 (3%)	2 (1%)	21	15
1	D	332/332 (100%)	317 (96%)	13 (4%)	2 (1%)	21	15
All	All	1328/1328 (100%)	1272 (96%)	48 (4%)	8 (1%)	21	15

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	218	SER
1	B	218	SER

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Mol	Chain	Res	Type
1	C	218	SER
1	D	218	SER
1	A	136	ILE
1	B	136	ILE
1	C	136	ILE
1	D	136	ILE

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	283/279 (101%)	279 (99%)	4 (1%)	59	66
1	B	280/279 (100%)	275 (98%)	5 (2%)	51	58
1	C	280/279 (100%)	275 (98%)	5 (2%)	51	58
1	D	281/279 (101%)	276 (98%)	5 (2%)	51	58
All	All	1124/1116 (101%)	1105 (98%)	19 (2%)	53	60

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	77	TYR
1	A	108	LYS
1	A	179	LEU
1	A	326	LEU
1	B	58	ASN
1	B	77	TYR
1	B	108	LYS
1	B	179	LEU
1	B	326	LEU
1	C	58	ASN
1	C	77	TYR
1	C	179	LEU
1	C	258	SER
1	C	326	LEU
1	D	58	ASN

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Mol	Chain	Res	Type
1	D	77	TYR
1	D	179	LEU
1	D	323	LEU
1	D	326	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	26	ASN
1	A	34	ASN
1	A	122	ASN
1	A	134	ASN
1	A	152	ASN
1	A	158	HIS
1	A	189	ASN
1	A	330	GLN
1	B	34	ASN
1	B	51	ASN
1	B	58	ASN
1	B	122	ASN
1	B	158	HIS
1	B	189	ASN
1	B	318	GLN
1	B	330	GLN
1	C	26	ASN
1	C	34	ASN
1	C	58	ASN
1	C	104	ASN
1	C	122	ASN
1	C	189	ASN
1	C	247	ASN
1	C	330	GLN
1	D	26	ASN
1	D	34	ASN
1	D	39	ASN
1	D	58	ASN
1	D	122	ASN
1	D	189	ASN
1	D	330	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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

20 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	$\# Z > 2$
3	SO4	D	702	-	4,4,4	1.85	2 (50%)	6,6,6	0.84	0
3	SO4	C	703	-	4,4,4	1.85	2 (50%)	6,6,6	0.84	0
2	NAG	C	601	1	14,14,15	0.58	0	17,19,21	0.64	0
3	SO4	A	703	-	4,4,4	1.85	2 (50%)	6,6,6	0.79	0
3	SO4	A	702	-	4,4,4	1.82	2 (50%)	6,6,6	0.83	0
5	PEG	B	704	-	6,6,6	0.84	0	5,5,5	1.53	1 (20%)
3	SO4	B	703	-	4,4,4	1.88	2 (50%)	6,6,6	0.82	0
3	SO4	B	702	-	4,4,4	1.85	2 (50%)	6,6,6	0.84	0
5	PEG	D	705	-	6,6,6	0.87	0	5,5,5	1.53	1 (20%)
5	PEG	C	704	-	6,6,6	0.90	0	5,5,5	1.56	1 (20%)
3	SO4	D	701	-	4,4,4	1.84	2 (50%)	6,6,6	0.86	0
3	SO4	D	703	-	4,4,4	1.88	2 (50%)	6,6,6	0.81	0
4	TRS	A	704	-	7,7,7	0.39	0	9,9,9	1.39	1 (11%)
5	PEG	A	705	-	6,6,6	0.86	0	5,5,5	1.54	1 (20%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SO4	C	702	-	4,4,4	1.83	2 (50%)	6,6,6	0.84	0
2	NAG	D	601	1	14,14,15	0.57	0	17,19,21	0.67	0
3	SO4	A	701	-	4,4,4	1.84	2 (50%)	6,6,6	0.83	0
4	TRS	D	704	-	7,7,7	0.64	0	9,9,9	1.62	2 (22%)
2	NAG	A	601	1	14,14,15	0.55	0	17,19,21	0.65	0
2	NAG	B	601	1	14,14,15	0.54	0	17,19,21	0.64	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
5	PEG	A	705	-	-	3/4/4/4	-
2	NAG	D	601	1	-	0/6/23/26	0/1/1/1
5	PEG	D	705	-	-	3/4/4/4	-
4	TRS	D	704	-	-	3/9/9/9	-
2	NAG	A	601	1	-	0/6/23/26	0/1/1/1
5	PEG	C	704	-	-	3/4/4/4	-
4	TRS	A	704	-	-	5/9/9/9	-
5	PEG	B	704	-	-	3/4/4/4	-
2	NAG	C	601	1	-	0/6/23/26	0/1/1/1
2	NAG	B	601	1	-	0/6/23/26	0/1/1/1

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	703	SO4	O1-S	3.12	1.63	1.44
3	B	703	SO4	O1-S	3.10	1.63	1.44
3	C	703	SO4	O1-S	3.05	1.62	1.44
3	B	702	SO4	O1-S	3.03	1.62	1.44
3	D	702	SO4	O1-S	3.02	1.62	1.44
3	A	703	SO4	O1-S	3.02	1.62	1.44
3	D	701	SO4	O1-S	3.01	1.62	1.44
3	A	701	SO4	O1-S	3.00	1.62	1.44
3	C	702	SO4	O1-S	2.99	1.62	1.44
3	A	702	SO4	O1-S	2.94	1.62	1.44
3	A	703	SO4	O3-S	-2.08	1.31	1.48
3	D	702	SO4	O3-S	-2.08	1.31	1.48
3	A	702	SO4	O3-S	-2.08	1.31	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	701	SO4	O3-S	-2.08	1.31	1.48
3	B	702	SO4	O3-S	-2.08	1.31	1.48
3	C	702	SO4	O3-S	-2.07	1.31	1.48
3	B	703	SO4	O3-S	-2.07	1.31	1.48
3	C	703	SO4	O3-S	-2.06	1.31	1.48
3	D	701	SO4	O3-S	-2.05	1.31	1.48
3	D	703	SO4	O3-S	-2.03	1.31	1.48

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	705	PEG	C3-O2-C2	2.71	125.11	113.26
5	C	704	PEG	C3-O2-C2	2.65	124.85	113.26
4	D	704	TRS	C2-C-N	-2.64	101.43	108.17
5	B	704	PEG	C3-O2-C2	2.63	124.78	113.26
5	D	705	PEG	C3-O2-C2	2.62	124.72	113.26
4	A	704	TRS	C2-C-N	-2.33	102.23	108.17
4	D	704	TRS	C3-C-N	2.19	113.74	108.17

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	704	TRS	N-C-C1-O1
4	A	704	TRS	C1-C-C3-O3
4	A	704	TRS	C2-C-C3-O3
4	A	704	TRS	N-C-C3-O3
4	D	704	TRS	C1-C-C3-O3
4	D	704	TRS	C2-C-C3-O3
4	D	704	TRS	N-C-C3-O3
5	A	705	PEG	O1-C1-C2-O2
5	B	704	PEG	O1-C1-C2-O2
5	C	704	PEG	O1-C1-C2-O2
5	D	705	PEG	O1-C1-C2-O2
5	A	705	PEG	O2-C3-C4-O4
5	D	705	PEG	O2-C3-C4-O4
5	B	704	PEG	O2-C3-C4-O4
5	C	704	PEG	O2-C3-C4-O4
5	D	705	PEG	C4-C3-O2-C2
4	A	704	TRS	C3-C-C1-O1
5	A	705	PEG	C4-C3-O2-C2
5	C	704	PEG	C4-C3-O2-C2

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Mol	Chain	Res	Type	Atoms
5	B	704	PEG	C4-C3-O2-C2

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	704	TRS	3	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	332/332 (100%)	-0.31	2 (0%) 85 87	10, 19, 27, 37	4 (1%)
1	B	332/332 (100%)	-0.31	2 (0%) 85 87	14, 20, 28, 37	1 (0%)
1	C	332/332 (100%)	-0.30	1 (0%) 90 91	11, 20, 29, 39	1 (0%)
1	D	332/332 (100%)	-0.28	3 (0%) 81 83	13, 21, 29, 38	2 (0%)
All	All	1328/1328 (100%)	-0.30	8 (0%) 85 87	10, 20, 29, 39	8 (0%)

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	254	ASN	3.1
1	A	1	ALA	2.6
1	D	107	TRP	2.6
1	B	265	ASN	2.5
1	B	1	ALA	2.4
1	A	265	ASN	2.1
1	D	265	ASN	2.1
1	D	51	ASN	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 oligosaccharides in this entry.

6.4 Ligands

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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	B-factors(Å ²)	Q<0.9
5	PEG	C	704	7/7	0.52	0.16	46,46,47,48	0
5	PEG	A	705	7/7	0.54	0.18	47,47,48,48	0
5	PEG	B	704	7/7	0.67	0.15	49,49,50,51	0
5	PEG	D	705	7/7	0.68	0.15	49,50,50,50	0
2	NAG	D	601	14/15	0.74	0.11	33,34,37,40	0
4	TRS	A	704	8/8	0.76	0.20	36,37,37,37	0
2	NAG	A	601	14/15	0.82	0.09	33,34,37,38	0
3	SO4	B	702	5/5	0.83	0.15	63,64,64,64	0
3	SO4	A	701	5/5	0.83	0.12	71,71,71,71	0
2	NAG	B	601	14/15	0.85	0.09	33,34,37,39	0
4	TRS	D	704	8/8	0.85	0.12	32,33,34,36	0
3	SO4	D	702	5/5	0.86	0.15	63,63,63,64	0
3	SO4	D	701	5/5	0.87	0.12	56,57,57,58	0
2	NAG	C	601	14/15	0.87	0.09	33,34,38,38	0
3	SO4	D	703	5/5	0.89	0.17	48,48,49,49	0
3	SO4	B	703	5/5	0.91	0.14	47,47,47,47	0
3	SO4	A	702	5/5	0.91	0.14	62,63,63,63	0
3	SO4	C	702	5/5	0.92	0.10	63,63,63,64	0
3	SO4	C	703	5/5	0.95	0.09	45,45,46,46	0
3	SO4	A	703	5/5	0.96	0.09	44,44,44,45	0

6.5 Other polymers

There are no such residues in this entry.