



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 17, 2018 – 08:16 pm GMT

PDB ID : 1SDX
Title : Crystal structure of the zinc saturated C-terminal half of bovine lactoferrin at 2.0 Å resolution reveals two additional zinc binding sites
Authors : Jabeen, T.; Sharma, S.; Singhal, G.; Singh, N.; Singh, T.P.
Deposited on : 2004-02-15
Resolution : 2.06 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.3 (157068), CSD as539be (2018)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk30686

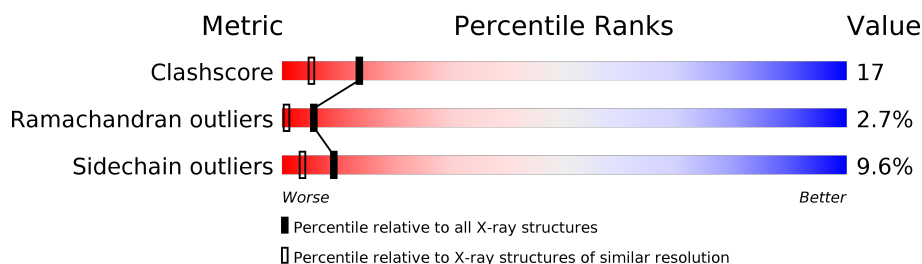
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.06 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))
Clashscore	122078	2357 (2.08-2.04)
Ramachandran outliers	120005	2338 (2.08-2.04)
Sidechain outliers	119972	2338 (2.08-2.04)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	335	
2	E	5	

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
3	NAG	A	6	X	-	-	-
6	MAN	A	10	X	-	X	-
6	MAN	A	11	X	-	-	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	MAN	A	8	-	-	X	-

2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 2973 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 Lactotransferrin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	335	Total	C	N	O	S	0	0	0
			2560	1593	448	499	20			

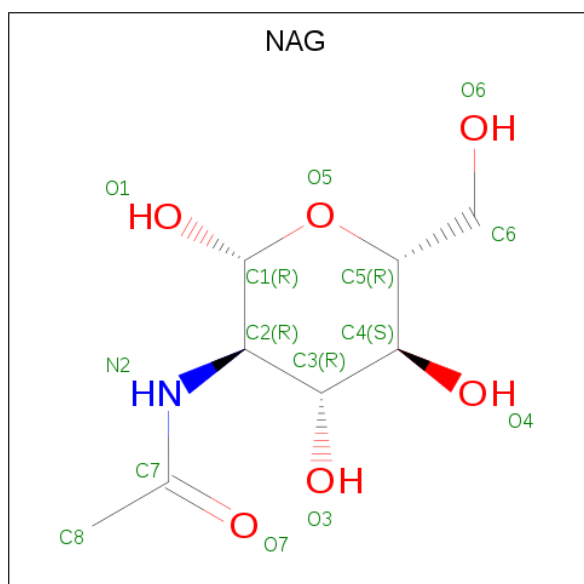
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	565	LYS	ASN	SEE REMARK 999	UNP P24627
A	608	GLU	LYS	SEE REMARK 999	UNP P24627

- Molecule 2 is a protein called Lactotransferrin.

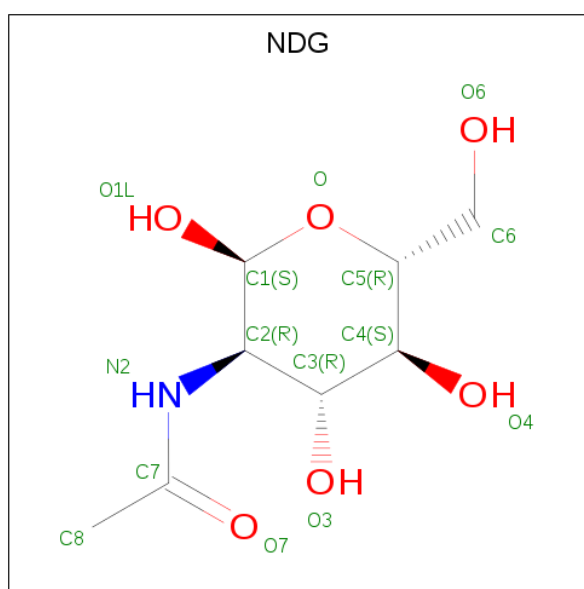
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	5	Total	C	N	O	S	0	0	0
			34	20	5	8	1			

- Molecule 3 is N-ACETYL-D-GLUCOSAMINE (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



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

- Molecule 4 is 2-(ACETYLAMINO)-2-DEOXY-A-D-GLUCOPYRANOSE (three-letter code: NDG) (formula: $C_8H_{15}NO_6$).



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

- Molecule 5 is BETA-D-MANNOSE (three-letter code: BMA) (formula: $C_6H_{12}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			11	6	5		

- Molecule 6 is ALPHA-D-MANNOSE (three-letter code: MAN) (formula: C₆H₁₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			11	6	5		
6	A	1	Total	C	O	0	0
			11	6	5		
6	A	1	Total	C	O	0	0
			11	6	5		

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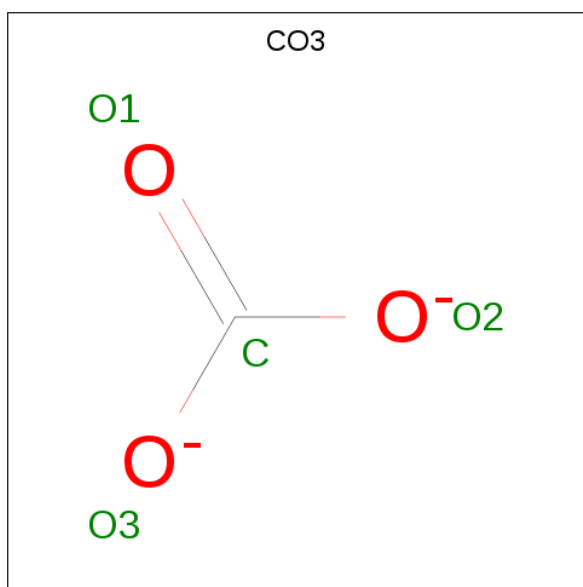
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			11	6	5		

- Molecule 7 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 8 is CARBONATE ION (three-letter code: CO3) (formula: CO₃).



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

- Molecule 9 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

- Molecule 10 is water.

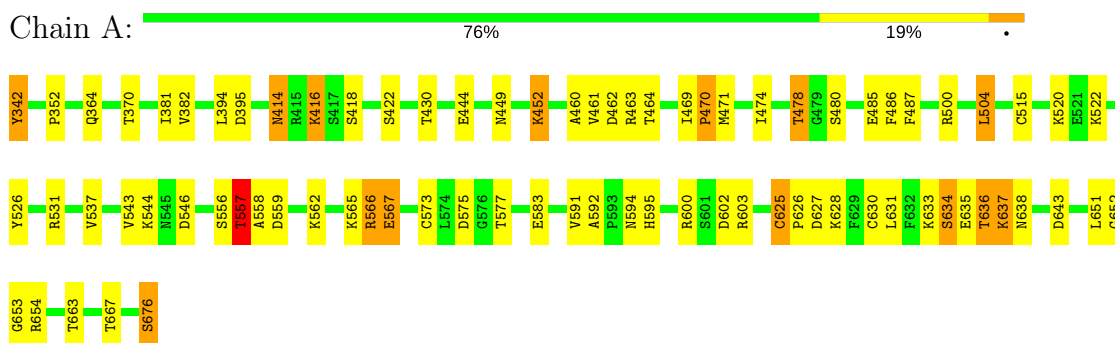
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	223	Total	O	0	0
			223	223		
10	E	5	Total	O	0	0
			5	5		

3 Residue-property plots

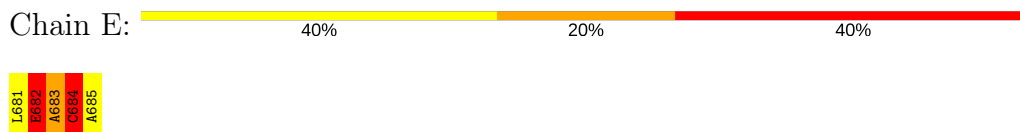
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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.

Note EDS was not executed.

• Molecule 1: Lactotransferrin



• Molecule 2: Lactotransferrin



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	63.21Å 50.57Å 66.17Å 90.00° 107.64° 90.00°	Depositor
Resolution (Å)	19.90 – 2.06	Depositor
% Data completeness (in resolution range)	89.9 (19.90-2.06)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
Refinement program	CNS 0.9	Depositor
R, R_{free}	0.192 , 0.210	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2973	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, BMA, NAG, CO3, NDG, SO4, MAN

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	0.40	0/2608	0.72	0/3533
2	E	0.76	0/33	1.96	2/42 (4.8%)
All	All	0.41	0/2641	0.75	2/3575 (0.1%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	682	GLU	N-CA-C	-6.45	93.60	111.00
2	E	683	ALA	N-CA-C	-6.24	94.14	111.00

There are no chirality outliers.

There are no planarity outliers.

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	2560	0	2480	75	0
2	E	34	0	30	5	0
3	A	56	0	48	0	0
4	A	28	0	25	0	0
5	A	11	0	9	1	0
6	A	44	0	38	13	0
7	A	3	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	A	4	0	0	0	0
9	A	5	0	0	0	0
10	A	223	0	0	7	0
10	E	5	0	0	0	0
All	All	2973	0	2630	92	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 17.

All (92) 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)
6:A:8:MAN:H61	6:A:10:MAN:C2	1.50	1.42
6:A:8:MAN:C6	6:A:10:MAN:H2	1.68	1.20
1:A:625:CYS:SG	1:A:626:PRO:HD3	1.96	1.05
6:A:8:MAN:H61	6:A:10:MAN:C3	1.88	1.04
1:A:485:GLU:HG2	1:A:500:ARG:HH21	1.32	0.95
1:A:485:GLU:HG2	1:A:500:ARG:NH2	1.87	0.90
1:A:638:ASN:HD22	1:A:643:ASP:H	1.20	0.84
6:A:10:MAN:H61	10:A:3089:HOH:O	1.75	0.84
1:A:416:LYS:HB3	1:A:416:LYS:NZ	1.94	0.81
1:A:565:LYS:CD	1:A:567:GLU:H	1.93	0.81
6:A:8:MAN:H61	6:A:10:MAN:H2	0.81	0.81
1:A:416:LYS:HD3	5:A:7:BMA:O6	1.82	0.78
1:A:565:LYS:HD3	1:A:567:GLU:H	1.48	0.78
6:A:8:MAN:C6	6:A:10:MAN:C3	2.61	0.76
1:A:430:THR:HB	1:A:594:ASN:ND2	2.02	0.75
6:A:8:MAN:C6	6:A:10:MAN:H3	2.17	0.75
1:A:625:CYS:SG	1:A:626:PRO:CD	2.74	0.73
6:A:8:MAN:C6	6:A:10:MAN:C2	2.46	0.73
1:A:638:ASN:ND2	1:A:643:ASP:H	1.86	0.73
6:A:8:MAN:O6	6:A:10:MAN:H3	1.89	0.72
1:A:478:THR:CG2	1:A:480:SER:HB3	2.19	0.72
1:A:485:GLU:CG	1:A:500:ARG:HH21	2.03	0.72
1:A:565:LYS:HD3	1:A:566:ARG:N	2.06	0.70
1:A:416:LYS:HB3	1:A:416:LYS:HZ3	1.55	0.68
1:A:676:SER:HB2	10:A:3215:HOH:O	1.94	0.68
6:A:8:MAN:H5	6:A:9:MAN:H2	1.77	0.66
1:A:342:TYR:N	1:A:342:TYR:CD1	2.63	0.66
1:A:575:ASP:OD1	1:A:577:THR:HG22	1.96	0.66
1:A:565:LYS:HD2	1:A:567:GLU:HB2	1.79	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:635:GLU:HG3	1:A:637:LYS:HE2	1.78	0.65
1:A:625:CYS:HB3	1:A:626:PRO:HD2	1.78	0.64
1:A:342:TYR:N	1:A:342:TYR:HD1	1.96	0.64
1:A:625:CYS:CB	1:A:626:PRO:CD	2.76	0.63
1:A:469:ILE:HB	1:A:470:PRO:CD	2.29	0.63
1:A:474:ILE:O	1:A:478:THR:HB	2.00	0.62
1:A:430:THR:HB	1:A:594:ASN:HD22	1.61	0.62
1:A:625:CYS:HB3	1:A:626:PRO:CD	2.30	0.60
1:A:382:VAL:HG22	2:E:682:GLU:HG2	1.84	0.59
1:A:469:ILE:HG22	1:A:470:PRO:HD3	1.86	0.58
1:A:634:SER:O	1:A:637:LYS:HD3	2.03	0.58
6:A:10:MAN:C6	10:A:3089:HOH:O	2.43	0.57
1:A:460:ALA:HB3	1:A:463:ARG:HD3	1.86	0.56
2:E:684:CYS:O	2:E:685:ALA:HB3	2.06	0.56
1:A:469:ILE:CB	1:A:470:PRO:CD	2.83	0.56
1:A:469:ILE:HB	1:A:470:PRO:HD2	1.87	0.56
6:A:8:MAN:C5	6:A:10:MAN:H2	2.35	0.55
2:E:682:GLU:CD	2:E:682:GLU:O	2.47	0.54
1:A:625:CYS:O	1:A:626:PRO:C	2.46	0.53
1:A:449:ASN:HB3	10:A:3191:HOH:O	2.09	0.53
6:A:8:MAN:H61	6:A:10:MAN:H3	1.75	0.52
1:A:635:GLU:O	1:A:636:THR:HG22	2.08	0.52
1:A:364:GLN:OE1	1:A:628:LYS:HE2	2.10	0.52
2:E:684:CYS:O	2:E:685:ALA:CB	2.59	0.51
1:A:461:VAL:O	1:A:462:ASP:HB2	2.10	0.51
1:A:452:LYS:HB2	10:A:3064:HOH:O	2.11	0.51
1:A:635:GLU:HG3	1:A:637:LYS:CE	2.40	0.51
1:A:478:THR:HG23	1:A:480:SER:HB3	1.92	0.51
1:A:565:LYS:HD2	1:A:567:GLU:H	1.74	0.51
1:A:562:LYS:NZ	1:A:562:LYS:HB2	2.26	0.51
1:A:565:LYS:C	1:A:565:LYS:HD3	2.31	0.51
1:A:471:MET:HE3	1:A:487:PHE:HE2	1.76	0.50
1:A:636:THR:HG23	1:A:636:THR:O	2.11	0.50
1:A:381:ILE:HD11	1:A:394:LEU:HD11	1.94	0.49
1:A:478:THR:HG22	1:A:480:SER:H	1.78	0.49
1:A:395:ASP:HA	1:A:595:HIS:CD2	2.48	0.49
1:A:352:PRO:HG2	1:A:520:LYS:HD2	1.97	0.47
1:A:653:GLY:O	1:A:654:ARG:C	2.53	0.47
1:A:573:CYS:HB2	1:A:577:THR:HG22	1.97	0.47
1:A:416:LYS:HB3	1:A:416:LYS:HZ2	1.78	0.46
1:A:469:ILE:CG2	1:A:470:PRO:HD3	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:452:LYS:HG2	1:A:486:PHE:CE1	2.50	0.46
1:A:636:THR:HG21	10:A:3107:HOH:O	2.15	0.46
2:E:682:GLU:OE1	2:E:682:GLU:O	2.34	0.46
1:A:565:LYS:HD2	1:A:567:GLU:CB	2.47	0.45
1:A:478:THR:HG22	1:A:480:SER:N	2.32	0.45
1:A:600:ARG:HB3	1:A:602:ASP:OD1	2.17	0.45
1:A:565:LYS:HZ2	1:A:567:GLU:HB2	1.82	0.45
1:A:504:LEU:HG	1:A:537:VAL:HG12	1.99	0.45
1:A:520:LYS:HE2	1:A:520:LYS:HB3	1.85	0.44
1:A:556:SER:C	1:A:557:THR:HG23	2.37	0.44
1:A:603:ARG:HD2	10:A:3111:HOH:O	2.17	0.44
1:A:663:THR:O	1:A:667:THR:HG23	2.17	0.44
1:A:565:LYS:NZ	1:A:567:GLU:HB2	2.34	0.43
1:A:464:THR:HG21	1:A:592:ALA:HB1	1.99	0.43
1:A:414:ASN:HD22	1:A:414:ASN:HA	1.58	0.43
1:A:522:LYS:HE3	1:A:531:ARG:CZ	2.49	0.43
1:A:625:CYS:C	1:A:630:CYS:SG	2.98	0.42
1:A:556:SER:C	1:A:558:ALA:H	2.23	0.42
1:A:562:LYS:HB2	1:A:562:LYS:HZ3	1.85	0.41
1:A:526:TYR:CZ	1:A:544:LYS:HD3	2.55	0.41
1:A:651:LEU:HD23	1:A:651:LEU:HA	1.85	0.41
1:A:573:CYS:SG	1:A:577:THR:HG23	2.62	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	333/335 (99%)	313 (94%)	14 (4%)	6 (2%)	9	2
2	E	3/5 (60%)	0	0	3 (100%)	0	0
All	All	336/340 (99%)	313 (93%)	14 (4%)	9 (3%)	5	1

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	625	CYS
2	E	682	GLU
2	E	683	ALA
2	E	684	CYS
1	A	634	SER
1	A	627	ASP
1	A	557	THR
1	A	652	GLY
1	A	543	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	278/278 (100%)	254 (91%)	24 (9%)	11	5
2	E	3/3 (100%)	0	3 (100%)	0	0
All	All	281/281 (100%)	254 (90%)	27 (10%)	9	3

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	342	TYR
1	A	370	THR
1	A	414	ASN
1	A	416	LYS
1	A	418	SER
1	A	422	SER
1	A	444	GLU
1	A	452	LYS
1	A	470	PRO
1	A	478	THR
1	A	504	LEU
1	A	515	CYS
1	A	546	ASP
1	A	557	THR

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Mol	Chain	Res	Type
1	A	559	ASP
1	A	566	ARG
1	A	567	GLU
1	A	583	GLU
1	A	591	VAL
1	A	631	LEU
1	A	633	LYS
1	A	636	THR
1	A	637	LYS
1	A	676	SER
2	E	681	LEU
2	E	682	GLU
2	E	684	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	414	ASN
1	A	594	ASN
1	A	621	ASN
1	A	638	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 16 ligands modelled in this entry, 3 are monoatomic - leaving 13 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	NAG	A	1	1,4	14,14,15	0.87	1 (7%)	17,19,21	2.51	3 (17%)
6	MAN	A	10	6	11,11,12	0.64	0	15,15,17	1.62	3 (20%)
6	MAN	A	11	3	11,11,12	0.83	0	15,15,17	1.23	2 (13%)
4	NDG	A	2	3	14,14,15	0.95	1 (7%)	17,19,21	2.00	4 (23%)
8	CO3	A	2001	7	0,3,3	0.00	-	0,3,3	0.00	-
3	NAG	A	3	1,4	14,14,15	0.98	1 (7%)	17,19,21	1.60	6 (35%)
9	SO4	A	3001	-	4,4,4	0.26	0	6,6,6	0.15	0
4	NDG	A	4	3,5	14,14,15	0.91	1 (7%)	17,19,21	1.98	4 (23%)
3	NAG	A	5	1,3	14,14,15	0.50	0	17,19,21	1.25	2 (11%)
3	NAG	A	6	3,6	14,14,15	0.88	0	17,19,21	1.78	6 (35%)
5	BMA	A	7	4,6	11,11,12	1.09	1 (9%)	15,15,17	2.26	4 (26%)
6	MAN	A	8	5,6	11,11,12	1.02	1 (9%)	15,15,17	2.91	7 (46%)
6	MAN	A	9	6	11,11,12	0.82	0	15,15,17	1.01	1 (6%)

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	NAG	A	1	1,4	-	0/6/23/26	0/1/1/1
6	MAN	A	10	6	1/1/4/5	0/2/19/22	0/1/1/1
6	MAN	A	11	3	1/1/4/5	0/2/19/22	1/1/1/1
4	NDG	A	2	3	-	0/6/23/26	0/1/1/1
8	CO3	A	2001	7	-	0/0/0/0	0/0/0/0
3	NAG	A	3	1,4	-	0/6/23/26	0/1/1/1
9	SO4	A	3001	-	-	0/0/0/0	0/0/0/0
4	NDG	A	4	3,5	-	0/6/23/26	0/1/1/1
3	NAG	A	5	1,3	-	0/6/23/26	0/1/1/1
3	NAG	A	6	3,6	1/1/5/7	0/6/23/26	0/1/1/1
5	BMA	A	7	4,6	-	0/2/19/22	0/1/1/1
6	MAN	A	8	5,6	-	0/2/19/22	0/1/1/1
6	MAN	A	9	6	-	0/2/19/22	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	8	MAN	C2-C3	-2.57	1.48	1.52
3	A	1	NAG	O5-C1	-2.54	1.39	1.43
5	A	7	BMA	O5-C1	-2.13	1.40	1.43
4	A	4	NDG	C1-C2	2.03	1.55	1.52
4	A	2	NDG	O-C5	2.31	1.48	1.43
3	A	3	NAG	C1-C2	2.45	1.55	1.52

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1	NAG	O5-C1-C2	-9.03	99.05	111.52
4	A	4	NDG	C4-C3-C2	-5.61	102.79	111.02
5	A	7	BMA	O4-C4-C3	-3.41	102.37	110.34
6	A	8	MAN	O4-C4-C5	-2.99	101.83	109.31
3	A	6	NAG	C2-N2-C7	-2.94	118.65	122.94
3	A	3	NAG	O4-C4-C3	-2.75	103.92	110.34
3	A	6	NAG	O4-C4-C5	-2.59	102.83	109.31
3	A	3	NAG	O4-C4-C5	-2.29	103.57	109.31
4	A	4	NDG	C2-N2-C7	-2.23	119.69	122.94
6	A	10	MAN	O3-C3-C2	-2.11	106.11	110.04
3	A	3	NAG	O5-C5-C6	-2.08	103.85	107.15
3	A	3	NAG	O7-C7-C8	-2.05	118.36	122.07
4	A	2	NDG	C4-C3-C2	-2.03	108.04	111.02
4	A	2	NDG	O-C5-C4	2.01	115.72	110.83
6	A	9	MAN	C2-C3-C4	2.07	114.46	110.87
3	A	1	NAG	C4-C3-C2	2.11	114.12	111.02
3	A	1	NAG	O4-C4-C3	2.15	115.36	110.34
3	A	5	NAG	O5-C5-C6	2.20	110.63	107.15
3	A	6	NAG	C3-C4-C5	2.26	114.28	110.24
3	A	6	NAG	C1-C2-N2	2.28	114.39	110.49
3	A	6	NAG	O5-C5-C6	2.37	110.90	107.15
6	A	10	MAN	C1-O5-C5	2.37	115.45	112.19
3	A	3	NAG	C3-C4-C5	2.41	114.55	110.24
3	A	6	NAG	O5-C1-C2	2.41	114.85	111.52
6	A	11	MAN	C1-O5-C5	2.45	115.55	112.19
6	A	11	MAN	O5-C1-C2	2.67	114.94	110.78
6	A	8	MAN	O3-C3-C4	2.69	116.63	110.34
3	A	5	NAG	C3-C4-C5	2.70	115.07	110.24
4	A	4	NDG	C1-C2-N2	2.78	115.23	110.49
3	A	3	NAG	C1-O5-C5	2.91	116.19	112.19
6	A	8	MAN	C1-O5-C5	3.01	116.33	112.19
4	A	4	NDG	O4-C4-C5	3.42	117.86	109.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	7	BMA	C1-C2-C3	3.96	114.67	109.66
6	A	8	MAN	C3-C4-C5	3.97	117.34	110.24
6	A	10	MAN	C1-C2-C3	4.37	115.19	109.66
5	A	7	BMA	C2-C3-C4	4.43	118.56	110.87
5	A	7	BMA	C6-C5-C4	4.48	123.57	112.99
6	A	8	MAN	O5-C1-C2	4.63	118.00	110.78
4	A	2	NDG	O-C1-C2	4.88	118.27	111.52
6	A	8	MAN	C1-C2-C3	5.06	116.06	109.66
6	A	8	MAN	C6-C5-C4	5.51	126.00	112.99
4	A	2	NDG	C1-O-C5	5.52	119.79	112.19

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
6	A	10	MAN	C1
3	A	6	NAG	C1
6	A	11	MAN	C1

There are no torsion outliers.

All (1) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	11	MAN	C1-C2-C3-C4-C5-O5

4 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	10	MAN	12	0
5	A	7	BMA	1	0
6	A	8	MAN	11	0
6	A	9	MAN	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.