



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2014 – 03:24 AM GMT

PDB ID : 2O1L
Title : Structure of a complex of C-terminal lobe of bovine lactoferrin with disaccharide at 1.97 Å resolution
Authors : Singh, N.; Sharma, S.; Perbandt, M.; Kaur, P.; Betzel, C.; Singh, T.P.
Deposited on : 2006-11-29
Resolution : 1.97 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

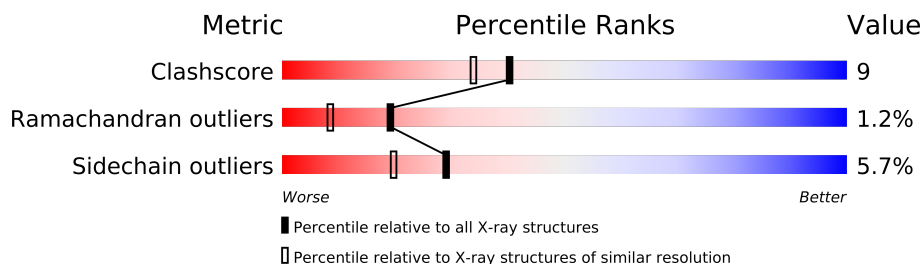
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 21963
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.97 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	79885	8091 (2.00-1.96)
Ramachandran outliers	78287	7989 (2.00-1.96)
Sidechain outliers	78261	7987 (2.00-1.96)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	345	

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 3086 atoms, of which 0 are hydrogen and 0 are deuterium.

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	341	Total	C	N	O	S	0	0	0
			2605	1622	454	508	21			

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 polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	2	Total	C	N	O	0	0
			28	16	2	10		
2	A	2	Total	C	N	O	0	0
			29	16	2	11		

There are 4 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
A	565	LYS	ASN	SEE REMARK 999	UNP P24627
A	608	GLU	LYS	SEE REMARK 999	UNP P24627

- Molecule 3 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	3	Total	C	N	O	0	0
			39	22	2	15		

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 4 is a polymer of unknown type called SUGAR (4-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	4	Total	C	N	O	0	0
			50	28	2	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

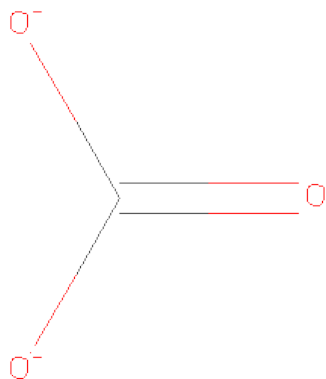
- Molecule 5 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	2	Total	Zn	0	0
			2	2		

- Molecule 6 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Fe	0	0
			1	1		

- Molecule 7 is CARBONATE ION (three-letter code: CO3) (formula: CO₃).



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

- Molecule 8 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

- Molecule 9 is water.

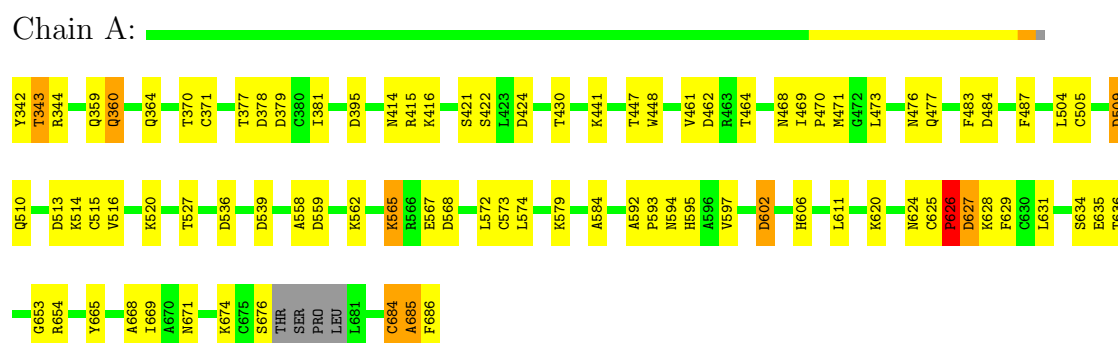
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	318	Total 318	O 318	0	0

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.

Note EDS was not executed.

- Molecule 1: Lactotransferrin



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	61.61Å 50.02Å 65.39Å 90.00° 107.11° 90.00°	Depositor
Resolution (Å)	62.02 – 1.97	Depositor
% Data completeness (in resolution range)	88.8 (62.02-1.97)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC 5	Depositor
R, R_{free}	0.214 , 0.231	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3086	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, NAG, SO4, CO3, FE, 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	0.49	0/2653	0.99	14/3591 (0.4%)

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
3	A	1	0
4	A	1	0
All	All	2	0

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	625	CYS	C-N-CD	-27.15	60.86	120.60
1	A	625	CYS	C-N-CA	8.80	158.95	122.00
1	A	626	PRO	CA-N-CD	-8.49	99.61	111.50
1	A	536	ASP	CB-CG-OD2	6.96	124.56	118.30
1	A	509	ASP	CB-CG-OD2	6.28	123.95	118.30
1	A	627	ASP	CB-CG-OD2	6.27	123.94	118.30
1	A	539	ASP	CB-CG-OD2	6.22	123.90	118.30
1	A	379	ASP	CB-CG-OD2	5.61	123.34	118.30
1	A	602	ASP	CB-CG-OD2	5.59	123.34	118.30
1	A	568	ASP	CB-CG-OD2	5.54	123.28	118.30
1	A	484	ASP	CB-CG-OD2	5.46	123.21	118.30
1	A	378	ASP	CB-CG-OD2	5.32	123.09	118.30
1	A	421	SER	N-CA-C	5.30	125.32	111.00
1	A	424	ASP	CB-CG-OD2	5.05	122.85	118.30

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	2	NAG	C1
4	A	8	MAN	C1

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2605	0	2519	49	0
2	A	57	0	52	1	0
3	A	39	0	34	1	0
4	A	50	0	43	2	0
5	A	2	0	0	0	0
6	A	1	0	0	0	0
7	A	4	0	0	0	0
8	A	10	0	0	0	0
9	A	318	0	0	8	0
All	All	3086	0	2648	51	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 9.

All (51) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:514:LYS:HB2	9:A:2098:HOH:O	1.77	0.83
1:A:685:ALA:HB3	9:A:2127:HOH:O	1.81	0.80
1:A:371:CYS:HB3	9:A:2045:HOH:O	1.82	0.80
1:A:359:GLN:HG2	9:A:2256:HOH:O	1.81	0.80
4:A:7:MAN:H5	4:A:8:MAN:H2	1.70	0.73
1:A:629:PHE:HA	9:A:2055:HOH:O	1.93	0.68
1:A:565:LYS:HD3	1:A:567:GLU:H	1.63	0.64
1:A:448:TRP:HE1	1:A:477:GLN:HE22	1.45	0.63
1:A:370:THR:HG22	1:A:371:CYS:N	2.17	0.59

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:527:THR:HG21	1:A:636:THR:O	2.02	0.59
1:A:665:TYR:CE1	1:A:669:ILE:HD11	2.37	0.59
1:A:461:VAL:O	1:A:462:ASP:HB2	2.02	0.58
1:A:370:THR:CG2	1:A:371:CYS:N	2.68	0.57
1:A:416:LYS:HE3	1:A:620:LYS:HE3	1.88	0.54
1:A:447:THR:HA	1:A:572:LEU:HD22	1.90	0.54
1:A:584:ALA:HB3	4:A:5:NAG:H82	1.90	0.54
1:A:624:ASN:HB3	1:A:628:LYS:HB2	1.90	0.53
1:A:513:ASP:OD2	1:A:520:LYS:NZ	2.39	0.53
1:A:430:THR:HB	1:A:594:ASN:ND2	2.26	0.51
1:A:395:ASP:HA	1:A:595:HIS:CD2	2.47	0.49
1:A:370:THR:CG2	1:A:371:CYS:H	2.25	0.49
1:A:441:LYS:HE2	1:A:567:GLU:O	2.12	0.48
1:A:505:CYS:O	1:A:514:LYS:HE2	2.15	0.47
1:A:415:ARG:HG2	1:A:416:LYS:N	2.30	0.47
1:A:471:MET:HE1	1:A:487:PHE:HE2	1.79	0.47
1:A:343:THR:O	1:A:343:THR:HG23	2.14	0.47
1:A:416:LYS:CE	1:A:620:LYS:HE3	2.45	0.47
1:A:513:ASP:OD2	1:A:520:LYS:HD3	2.15	0.46
1:A:469:ILE:HB	1:A:470:PRO:HD3	1.98	0.46
1:A:668:ALA:HB1	3:A:2:NAG:H83	1.98	0.46
1:A:573:CYS:SG	1:A:579:LYS:HG3	2.56	0.45
1:A:509:ASP:OD1	1:A:510:GLN:HG3	2.15	0.45
1:A:653:GLY:O	1:A:654:ARG:C	2.55	0.45
2:A:687:NAG:H5	2:A:688:NAG:O5	2.14	0.45
1:A:360:GLN:NE2	9:A:2198:HOH:O	2.49	0.45
1:A:635:GLU:N	9:A:2089:HOH:O	2.49	0.45
1:A:468:ASN:HB3	1:A:669:ILE:CD1	2.46	0.44
1:A:364:GLN:HG3	1:A:629:PHE:HB2	1.99	0.44
1:A:343:THR:OG1	1:A:606:HIS:CE1	2.72	0.43
1:A:684:CYS:C	1:A:686:PHE:H	2.21	0.43
1:A:448:TRP:HE1	1:A:477:GLN:NE2	2.16	0.42
1:A:377:THR:O	1:A:381:ILE:HG12	2.19	0.42
1:A:671:ASN:HA	1:A:674:LYS:HD2	2.01	0.42
1:A:558:ALA:O	1:A:562:LYS:HG3	2.20	0.42
1:A:516:VAL:O	1:A:516:VAL:HG23	2.20	0.42
1:A:483:PHE:C	1:A:483:PHE:CD2	2.93	0.42
1:A:473:LEU:HD13	1:A:574:LEU:HD21	2.02	0.42
1:A:597:VAL:HG11	1:A:611:LEU:HD21	2.02	0.41
1:A:464:THR:HG21	1:A:592:ALA:HB1	2.02	0.41
1:A:469:ILE:HB	1:A:470:PRO:CD	2.51	0.41
1:A:559:ASP:HB2	9:A:2234:HOH:O	2.20	0.41

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	337/345 (98%)	315 (94%)	18 (5%)	4 (1%)	19 8

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	626	PRO
1	A	634	SER
1	A	684	CYS
1	A	685	ALA

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	282/286 (99%)	266 (94%)	16 (6%)	29 20

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	342	TYR
1	A	343	THR
1	A	344	ARG
1	A	360	GLN
1	A	414	ASN
1	A	422	SER
1	A	476	ASN

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Mol	Chain	Res	Type
1	A	504	LEU
1	A	515	CYS
1	A	565	LYS
1	A	593	PRO
1	A	602	ASP
1	A	626	PRO
1	A	627	ASP
1	A	631	LEU
1	A	676	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	359	GLN
1	A	414	ASN
1	A	477	GLN
1	A	606	HIS
1	A	624	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

11 carbohydrates 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
2	NAG	A	1	1,2	12,14,15	0.69	0	15,19,21	1.20	2 (13%)
3	NAG	A	2	1,3	12,14,15	0.53	0	15,19,21	1.73	1 (6%)
3	NAG	A	3	3	12,14,15	0.64	0	15,19,21	1.12	1 (6%)
3	MAN	A	4	3	10,11,12	0.69	0	11,15,17	1.87	3 (27%)
4	NAG	A	5	1,4	12,14,15	0.83	1 (8%)	15,19,21	0.94	0
4	NAG	A	6	4	12,14,15	0.68	1 (8%)	15,19,21	0.75	0
2	NAG	A	687	2	15,15,15	0.48	0	21,21,21	1.64	5 (23%)
2	NAG	A	688	2	12,14,15	0.68	0	15,19,21	1.72	4 (26%)
4	MAN	A	7	4	10,11,12	0.69	0	11,15,17	1.87	3 (27%)
4	MAN	A	8	4	10,11,12	0.75	0	11,15,17	1.74	1 (9%)
2	NAG	A	9	2	12,14,15	0.59	0	15,19,21	0.99	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
2	NAG	A	1	1,2	-	0/6/23/26	0/1/1/1
3	NAG	A	2	1,3	1/1/5/7	0/6/23/26	0/1/1/1
3	NAG	A	3	3	-	0/6/23/26	0/1/1/1
3	MAN	A	4	3	-	0/2/19/22	0/1/1/1
4	NAG	A	5	1,4	-	0/6/23/26	0/1/1/1
4	NAG	A	6	4	-	0/6/23/26	0/1/1/1
2	NAG	A	687	2	-	0/6/26/26	0/1/1/1
2	NAG	A	688	2	-	0/6/23/26	0/1/1/1
4	MAN	A	7	4	-	0/2/19/22	0/1/1/1
4	MAN	A	8	4	1/1/4/5	0/2/19/22	0/1/1/1
2	NAG	A	9	2	-	0/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	5	NAG	O5-C5	-2.24	1.41	1.45
4	A	6	NAG	O5-C5	-2.12	1.41	1.45

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	2	NAG	O5-C5-C6	6.08	113.36	106.98
3	A	4	MAN	O5-C5-C6	4.60	111.81	106.98
4	A	7	MAN	O5-C5-C6	4.59	111.80	106.98
4	A	8	MAN	O5-C5-C6	4.46	111.67	106.98
2	A	687	NAG	C3-C4-C5	3.78	116.96	110.20
2	A	687	NAG	C4-C3-C2	3.64	115.61	110.44
2	A	688	NAG	O5-C5-C6	3.14	110.27	106.98
2	A	1	NAG	O5-C5-C6	2.89	110.02	106.98
3	A	4	MAN	C4-C3-C2	2.64	114.05	110.50
4	A	7	MAN	C4-C3-C2	2.63	114.03	110.50
2	A	687	NAG	C1-O5-C5	-2.59	108.75	113.40
2	A	687	NAG	C1-C2-C3	2.56	114.27	110.59
2	A	688	NAG	C2-N2-C7	2.54	127.36	123.09
2	A	688	NAG	C3-C4-C5	2.48	114.62	110.20
3	A	3	NAG	O5-C5-C6	2.43	109.53	106.98
2	A	687	NAG	C1-C2-N2	-2.37	108.09	110.85
2	A	688	NAG	C6-C5-C4	-2.35	107.32	113.00
4	A	7	MAN	C3-C4-C5	2.21	114.15	110.20
3	A	4	MAN	C3-C4-C5	2.21	114.15	110.20
2	A	9	NAG	C3-C4-C5	2.12	113.98	110.20
2	A	1	NAG	C3-C2-N2	-2.02	108.68	111.76

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	2	NAG	C1
4	A	8	MAN	C1

There are no torsion outliers.

There are no ring outliers.

5.6 Ligand geometry ⓘ

Of 6 ligands modelled in this entry, 3 are monoatomic - leaving 3 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
7	CO3	A	1999	6	0,3,3	0.00	-	0,3,3	0.00	-
8	SO4	A	2000	-	4,4,4	0.26	0	6,6,6	0.12	0
8	SO4	A	2001	-	4,4,4	0.21	0	6,6,6	0.10	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
7	CO3	A	1999	6	-	0/0/0/0	0/0/0/0
8	SO4	A	2000	-	-	0/0/0/0	0/0/0/0
8	SO4	A	2001	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.