



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 13, 2018 – 02:00 pm GMT

PDB ID : 2Q7Y
Title : Structure of the endogenous iNKT cell ligand iGb3 bound to mCD1d
Authors : Zajonc, D.M.; Wilson, I.A.; Teyton, L.
Deposited on : 2007-06-07
Resolution : 1.95 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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) : 1.13
EDS : trunk31020
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac : 5.8.0158
CCP4 : 7.0 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk31020

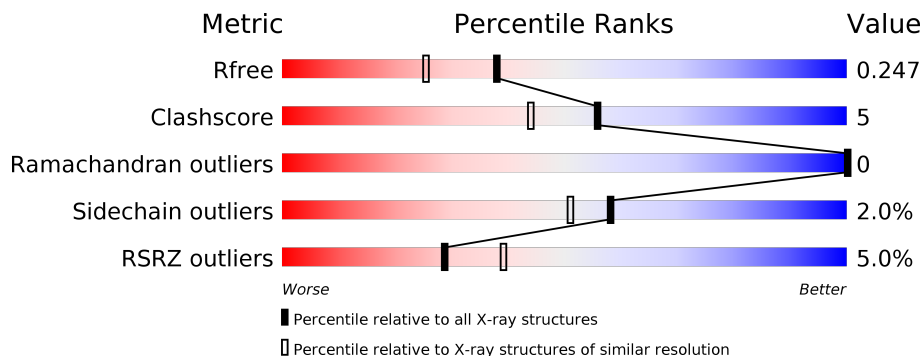
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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}	111664	2220 (1.96-1.96)
Clashscore	122126	2333 (1.96-1.96)
Ramachandran outliers	120053	2314 (1.96-1.96)
Sidechain outliers	120020	2314 (1.96-1.96)
RSRZ outliers	108989	2174 (1.96-1.96)

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. 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	285	<div> <div>6%</div> <div> <div></div> <div>78%</div> <div>16%</div> <div>• 6%</div> </div> </div>
1	C	285	<div> <div>5%</div> <div> <div></div> <div>82%</div> <div>12%</div> <div>• 6%</div> </div> </div>
2	B	99	<div> <div>2%</div> <div> <div></div> <div>83%</div> <div>14%</div> <div>• •</div> </div> </div>
2	D	99	<div> <div>2%</div> <div> <div></div> <div>85%</div> <div>13%</div> <div>• •</div> </div> </div>

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 6589 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 T-cell surface glycoprotein CD1d1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	269	Total	C	N	O	S	0	0	0
			2165	1381	372	399	13			
1	C	269	Total	C	N	O	S	0	0	0
			2165	1381	372	399	13			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	201	HIS	ASP	CONFLICT	UNP P11609
A	280	HIS	-	EXPRESSION TAG	UNP P11609
A	281	HIS	-	EXPRESSION TAG	UNP P11609
A	282	HIS	-	EXPRESSION TAG	UNP P11609
A	283	HIS	-	EXPRESSION TAG	UNP P11609
A	284	HIS	-	EXPRESSION TAG	UNP P11609
A	285	HIS	-	EXPRESSION TAG	UNP P11609
C	201	HIS	ASP	CONFLICT	UNP P11609
C	280	HIS	-	EXPRESSION TAG	UNP P11609
C	281	HIS	-	EXPRESSION TAG	UNP P11609
C	282	HIS	-	EXPRESSION TAG	UNP P11609
C	283	HIS	-	EXPRESSION TAG	UNP P11609
C	284	HIS	-	EXPRESSION TAG	UNP P11609
C	285	HIS	-	EXPRESSION TAG	UNP P11609

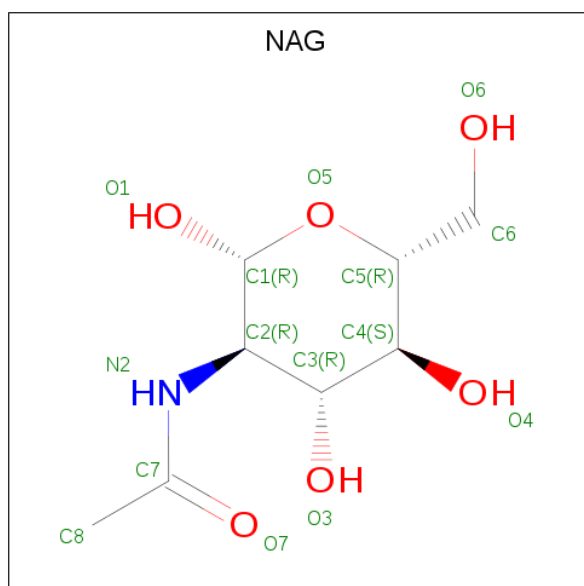
- Molecule 2 is a protein called Beta-2-microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	98	Total	C	N	O	S	0	0	0
			810	517	137	149	7			
2	D	98	Total	C	N	O	S	0	0	0
			810	517	137	149	7			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	85	ALA	ASP	VARIANT	UNP P01887
D	85	ALA	ASP	VARIANT	UNP P01887

- 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		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	C	1	Total	C	N	O	0	0
			14	8	1	5		
3	C	1	Total	C	N	O	0	0
			14	8	1	5		
3	C	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 4 is BETA-D-MANNOSE (three-letter code: BMA) (formula: $C_6H_{12}O_6$).



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

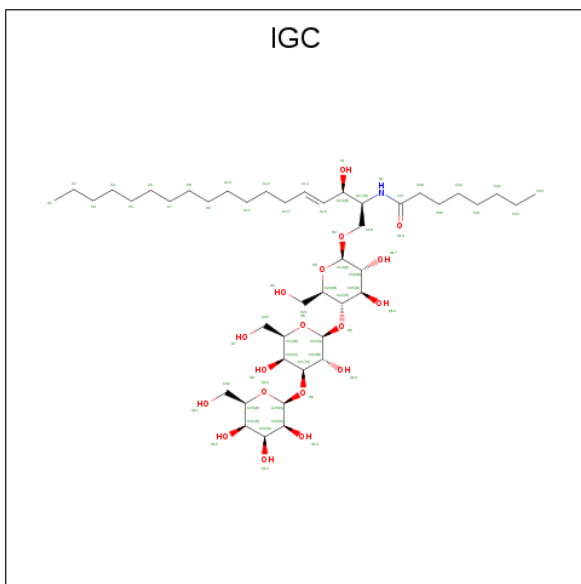
- Molecule 5 is ALPHA-D-MANNOSE (three-letter code: MAN) (formula: $C_6H_{12}O_6$).



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

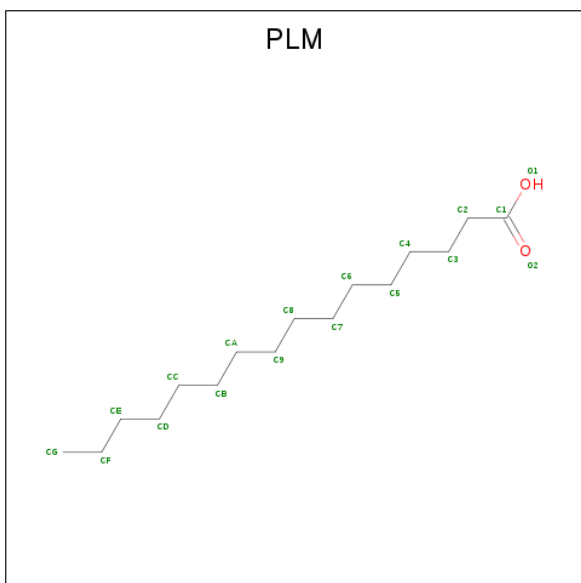
- Molecule 6 is N-[(1S,2R,3E)-1-({[ALPHA-D-GALACTOPYRANOSYL-(1->3)-BETA-D

-GALACTOPYRANOSYL-(1->4)-BETA-D-GLUCOPYRANOSYL]OXY}METHYL)-2-HYDROXYHEPTADEC-3-EN-1-YL]OCTANAMIDE (three-letter code: IGC) (formula: $C_{44}H_{81}NO_{18}$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	N	O	0	0
			41	32	1	8		
6	C	1	Total	C	N	O	0	0
			52	38	1	13		

- Molecule 7 is PALMITIC ACID (three-letter code: PLM) (formula: $C_{16}H_{32}O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			18	16	2		
7	C	1	Total	C	O	0	0
			18	16	2		

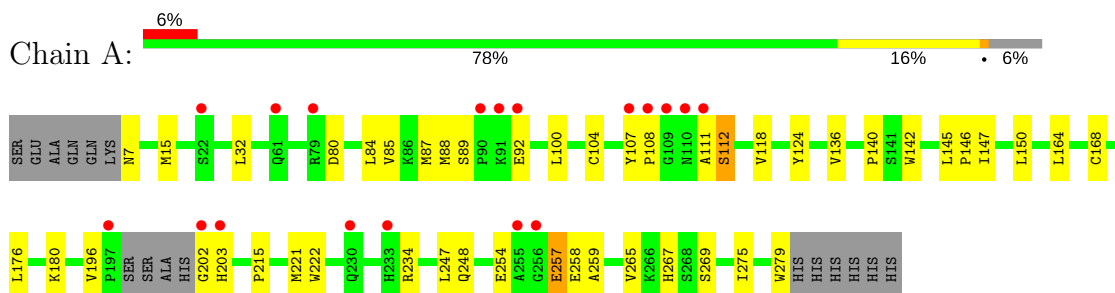
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	96	Total	O	0	0
			96	96		
8	B	52	Total	O	0	0
			52	52		
8	C	145	Total	O	0	0
			145	145		
8	D	58	Total	O	0	0
			58	58		

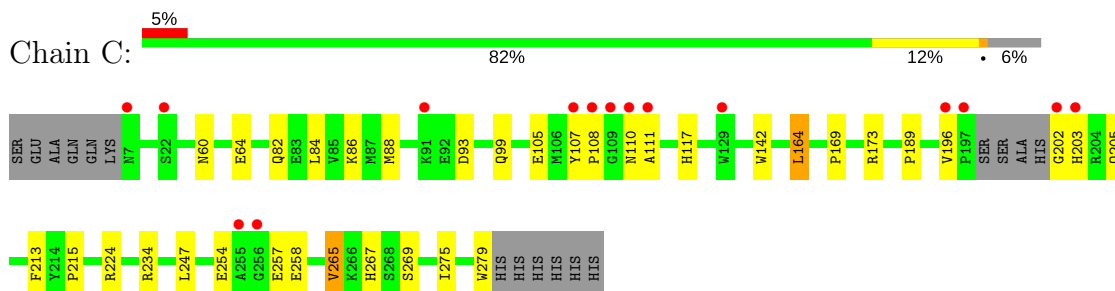
3 Residue-property plots [i](#)

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.

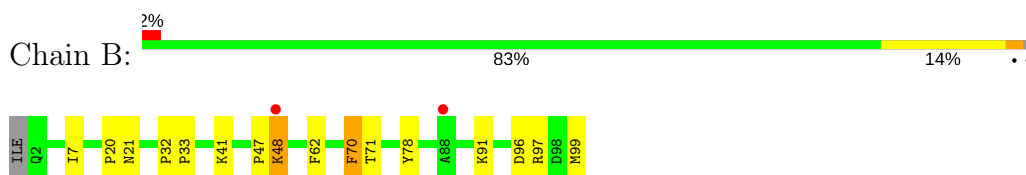
- Molecule 1: T-cell surface glycoprotein CD1d1



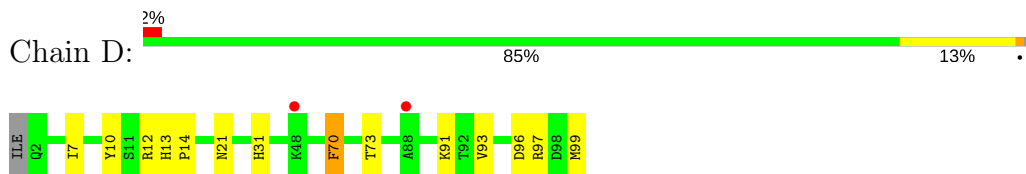
- Molecule 1: T-cell surface glycoprotein CD1d1



- Molecule 2: Beta-2-microglobulin



- Molecule 2: Beta-2-microglobulin



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	58.94Å 96.22Å 77.93Å 90.00° 106.23° 90.00°	Depositor
Resolution (Å)	27.91 – 1.95 27.90 – 1.95	Depositor EDS
% Data completeness (in resolution range)	97.5 (27.91-1.95) 97.5 (27.90-1.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.12 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.199 , 0.245 0.198 , 0.247	Depositor DCC
R_{free} test set	1219 reflections (2.06%)	wwPDB-VP
Wilson B-factor (Å ²)	27.0	Xtriage
Anisotropy	0.009	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 49.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6589	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 34.36 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 6.9034e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: PLM, BMA, IGC, NAG, 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	0.72	4/2228 (0.2%)	0.75	1/3028 (0.0%)
1	C	0.72	1/2228 (0.0%)	0.75	1/3028 (0.0%)
2	B	0.67	0/836	0.75	0/1133
2	D	0.76	0/836	0.89	2/1133 (0.2%)
All	All	0.72	5/6128 (0.1%)	0.77	4/8322 (0.0%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	202	GLY	N-CA	9.46	1.60	1.46
1	C	202	GLY	N-CA	8.54	1.58	1.46
1	A	257	GLU	CD-OE2	7.72	1.34	1.25
1	A	257	GLU	CD-OE1	6.06	1.32	1.25
1	A	259	ALA	C-N	5.44	1.42	1.33

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	12	ARG	NE-CZ-NH1	6.01	123.30	120.30
1	A	112	SER	N-CA-C	5.74	126.49	111.00
2	D	12	ARG	NE-CZ-NH2	-5.54	117.53	120.30
1	C	164	LEU	CB-CG-CD1	5.47	120.30	111.00

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	2165	0	2073	31	0
1	C	2165	0	2071	19	0
2	B	810	0	783	10	0
2	D	810	0	783	8	0
3	A	70	0	62	0	0
3	C	56	0	51	0	0
4	A	11	0	8	0	0
5	A	22	0	20	0	0
6	A	41	0	60	0	0
6	C	52	0	70	3	0
7	A	18	0	31	0	0
7	C	18	0	31	0	0
8	A	96	0	0	0	0
8	B	52	0	0	0	0
8	C	145	0	0	0	0
8	D	58	0	0	0	0
All	All	6589	0	6043	67	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.

The worst 5 of 67 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:111:ALA:O	1:A:112:SER:HB3	1.59	1.02
1:A:107:TYR:HB3	1:A:108:PRO:HD2	1.48	0.92
2:D:7:ILE:HD12	2:D:91:LYS:HD3	1.52	0.92
6:C:602:IGC:O16	6:C:602:IGC:H23	1.77	0.85
2:D:96:ASP:OD1	2:D:96:ASP:O	1.94	0.83

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	265/285 (93%)	259 (98%)	6 (2%)	0	100	100
1	C	265/285 (93%)	261 (98%)	4 (2%)	0	100	100
2	B	96/99 (97%)	94 (98%)	2 (2%)	0	100	100
2	D	96/99 (97%)	94 (98%)	2 (2%)	0	100	100
All	All	722/768 (94%)	708 (98%)	14 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	235/249 (94%)	231 (98%)	4 (2%)	63	58
1	C	235/249 (94%)	230 (98%)	5 (2%)	56	49
2	B	92/93 (99%)	90 (98%)	2 (2%)	55	47
2	D	92/93 (99%)	90 (98%)	2 (2%)	55	47
All	All	654/684 (96%)	641 (98%)	13 (2%)	58	51

5 of 13 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	70	PHE
1	C	93	ASP
1	C	265	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	B	48	LYS
1	C	164	LEU

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	62	GLN
1	A	267	HIS
1	C	267	HIS
2	D	31	HIS

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](#)

16 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	NAG	A	500	1	14,14,15	0.63	0	17,19,21	1.36	3 (17%)
3	NAG	A	501	1,3	14,14,15	0.70	0	17,19,21	1.13	2 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	A	502	3	14,14,15	0.64	0	17,19,21	1.73	4 (23%)
3	NAG	A	511	1,3	14,14,15	0.70	0	17,19,21	1.20	3 (17%)
3	NAG	A	512	3,4	14,14,15	0.55	0	17,19,21	1.22	1 (5%)
4	BMA	A	513	3,5	11,11,12	0.62	0	15,15,17	0.86	0
5	MAN	A	514	4	11,11,12	0.51	0	15,15,17	1.17	1 (6%)
5	MAN	A	515	4	11,11,12	0.43	0	15,15,17	1.17	1 (6%)
6	IGC	A	601	-	40,41,65	1.19	4 (10%)	46,49,85	1.96	2 (4%)
7	PLM	A	701	-	14,17,17	0.38	0	13,17,17	0.54	0
3	NAG	C	500	1	14,14,15	0.57	0	17,19,21	1.06	2 (11%)
3	NAG	C	501	1	14,14,15	0.65	0	17,19,21	1.35	1 (5%)
3	NAG	C	502	1,3	14,14,15	0.82	0	17,19,21	1.03	2 (11%)
3	NAG	C	503	3	14,14,15	0.58	0	17,19,21	1.31	2 (11%)
6	IGC	C	602	-	52,53,65	1.41	8 (15%)	64,67,85	1.26	8 (12%)
7	PLM	C	702	-	14,17,17	0.37	0	13,17,17	0.52	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
3	NAG	A	500	1	-	0/6/23/26	0/1/1/1
3	NAG	A	501	1,3	-	0/6/23/26	0/1/1/1
3	NAG	A	502	3	-	0/6/23/26	0/1/1/1
3	NAG	A	511	1,3	-	0/6/23/26	0/1/1/1
3	NAG	A	512	3,4	-	0/6/23/26	0/1/1/1
4	BMA	A	513	3,5	-	0/2/19/22	0/1/1/1
5	MAN	A	514	4	-	0/2/19/22	0/1/1/1
5	MAN	A	515	4	-	0/2/19/22	0/1/1/1
6	IGC	A	601	-	-	0/37/57/109	0/1/1/3
7	PLM	A	701	-	-	0/13/15/15	0/0/0/0
3	NAG	C	500	1	-	0/6/23/26	0/1/1/1
3	NAG	C	501	1	-	0/6/23/26	0/1/1/1
3	NAG	C	502	1,3	-	0/6/23/26	0/1/1/1
3	NAG	C	503	3	-	0/6/23/26	0/1/1/1
6	IGC	C	602	-	-	0/43/83/109	0/2/2/3
7	PLM	C	702	-	-	0/13/15/15	0/0/0/0

The worst 5 of 12 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	601	IGC	C13-C14	-2.21	1.38	1.50
6	C	602	IGC	O5-C23	2.00	1.47	1.41
6	A	601	IGC	C38-C37	2.02	1.55	1.51
6	C	602	IGC	C16-C15	2.06	1.53	1.50
6	A	601	IGC	O3-C19	2.07	1.47	1.41

The worst 5 of 32 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	601	IGC	C16-C15-C14	-8.62	113.64	125.22
3	C	501	NAG	O5-C1-C2	-3.59	106.57	111.52
6	C	602	IGC	C35-C22-C20	-3.52	102.75	110.93
6	C	602	IGC	C16-C15-C14	-3.26	120.84	125.22
3	C	503	NAG	C2-N2-C7	-2.86	118.77	122.94

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	C	602	IGC	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	269/285 (94%)	0.36	18 (6%) 18 26	16, 31, 63, 91	0
1	C	269/285 (94%)	0.21	15 (5%) 24 33	15, 26, 55, 78	0
2	B	98/99 (98%)	0.07	2 (2%) 65 73	18, 31, 50, 57	0
2	D	98/99 (98%)	-0.05	2 (2%) 65 73	15, 27, 45, 57	0
All	All	734/768 (95%)	0.21	37 (5%) 29 39	15, 28, 57, 91	0

The worst 5 of 37 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	111	ALA	7.6
1	C	110	ASN	6.6
1	A	255	ALA	6.2
1	C	108	PRO	5.7
1	C	109	GLY	5.7

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. 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(\AA^2)	Q<0.9
6	IGC	C	602	52/63	0.63	0.24	34,50,81,85	0
6	IGC	A	601	41/63	0.66	0.23	41,48,70,73	0
4	BMA	A	513	11/12	0.80	0.27	48,55,61,66	0
3	NAG	A	502	14/15	0.81	0.36	50,55,59,60	0
3	NAG	C	502	14/15	0.82	0.12	32,36,41,41	0
5	MAN	A	514	11/12	0.83	0.28	68,71,76,79	0
3	NAG	A	500	14/15	0.84	0.24	47,50,58,58	0
7	PLM	A	701	18/18	0.84	0.33	40,47,58,58	0
3	NAG	C	500	14/15	0.85	0.27	54,60,67,71	0
3	NAG	A	512	14/15	0.86	0.23	43,47,52,53	0
3	NAG	C	503	14/15	0.86	0.25	48,51,59,60	0
3	NAG	A	511	14/15	0.88	0.13	34,38,43,43	0
5	MAN	A	515	11/12	0.89	0.19	44,48,51,53	0
7	PLM	C	702	18/18	0.91	0.28	37,39,53,55	0
3	NAG	A	501	14/15	0.95	0.09	27,33,40,41	0
3	NAG	C	501	14/15	0.97	0.09	29,33,40,41	0

6.5 Other polymers [i](#)

There are no such residues in this entry.