



# wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 19, 2016 – 09:11 PM GMT

PDB ID : 4ZNE  
Title : IgG1 Fc-FcγRI ecd complex  
Authors : Oganessian, V.Y.; Dall'Acqua, W.F.  
Deposited on : 2015-05-04  
Resolution : 2.42 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](mailto: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.

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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-20026982  
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-20026982

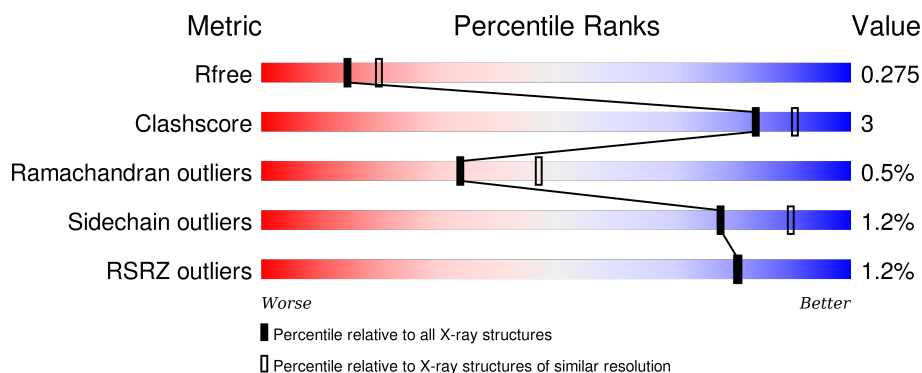
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.42 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	3386 (2.44-2.40)
Clashscore	102246	3897 (2.44-2.40)
Ramachandran outliers	100387	3837 (2.44-2.40)
Sidechain outliers	100360	3838 (2.44-2.40)
RSRZ outliers	91569	3396 (2.44-2.40)

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  $\geq 3$ , 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	267	<div> <div>3%</div> <div>78%</div> <div>11%</div> <div>9%</div> </div>
2	E	227	<div> <div>93%</div> <div>5%</div> </div>
2	J	227	<div> <div>83%</div> <div>9%</div> <div>8%</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
3	NAG	A	305	X	-	-	-
3	NAG	A	309	X	-	-	-
3	NAG	E	501	-	-	-	X
3	NAG	J	502	-	-	-	X
4	MAN	E	503	X	-	-	-
4	MAN	J	503	X	-	-	-
5	ZN	A	310	-	-	-	X
5	ZN	J	509	-	-	-	X

## 2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 5767 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 High affinity immunoglobulin gamma Fc receptor I.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	242	Total	C	N	O	S	0	0	0
			1932	1228	333	363	8			

- Molecule 2 is a protein called Ig gamma-1 chain C region.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	215	Total	C	N	O	S	0	0	0
			1711	1090	287	328	6			
2	J	209	Total	C	N	O	S	0	0	0
			1668	1061	281	320	6			

- Molecule 3 is N-ACETYL-D-GLUCOSAMINE (three-letter code: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).



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

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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	E	1	Total	C	N	O	0	0
			14	8	1	5		
3	E	1	Total	C	N	O	0	0
			14	8	1	5		
3	E	1	Total	C	N	O	0	0
			14	8	1	5		
3	E	1	Total	C	N	O	0	0
			14	8	1	5		
3	J	1	Total	C	N	O	0	0
			14	8	1	5		
3	J	1	Total	C	N	O	0	0
			14	8	1	5		
3	J	1	Total	C	N	O	0	0
			14	8	1	5		
3	J	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 4 is ALPHA-D-MANNOSE (three-letter code: MAN) (formula: C<sub>6</sub>H<sub>12</sub>O<sub>6</sub>).

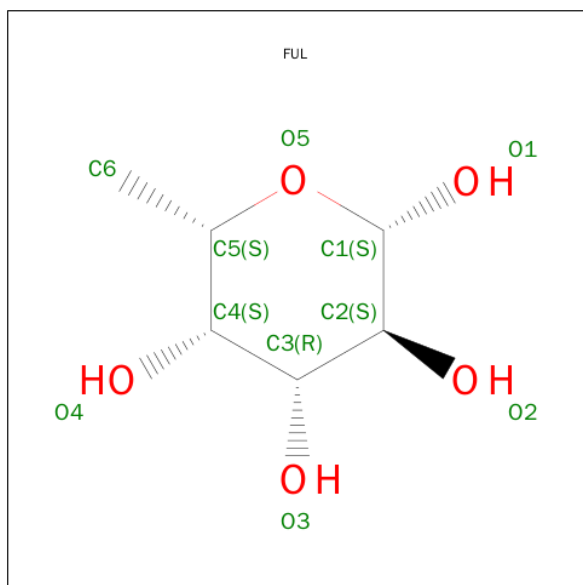


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			11	6	5		
4	A	1	Total	C	O	0	0
			11	6	5		
4	E	1	Total	C	O	0	0
			11	6	5		
4	E	1	Total	C	O	0	0
			11	6	5		
4	E	1	Total	C	O	0	0
			11	6	5		
4	J	1	Total	C	O	0	0
			11	6	5		
4	J	1	Total	C	O	0	0
			11	6	5		
4	J	1	Total	C	O	0	0
			11	6	5		

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	J	1	Total	Zn	0	0
			1	1		
5	A	1	Total	Zn	0	0
			1	1		
5	E	1	Total	Zn	0	0
			1	1		

- Molecule 6 is BETA-L-FUCOSE (three-letter code: FUL) (formula: C<sub>6</sub>H<sub>12</sub>O<sub>5</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	E	1	Total	C	O	0	0
			10	6	4		
6	J	1	Total	C	O	0	0
			10	6	4		

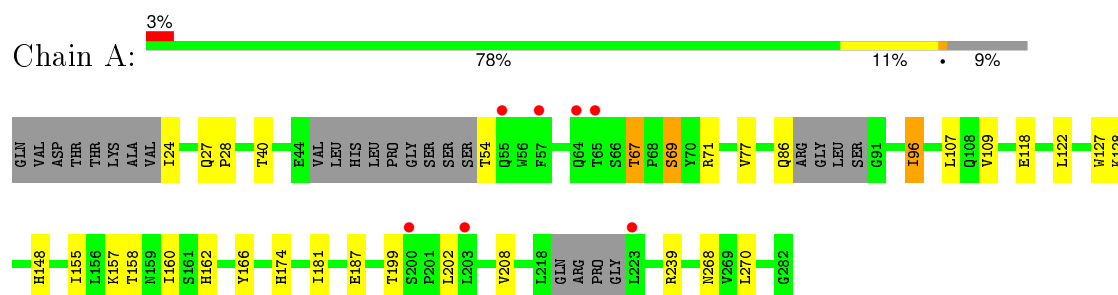
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	16	Total	O	0	0
			16	16		
7	E	64	Total	O	0	0
			64	64		
7	J	55	Total	O	0	0
			55	55		

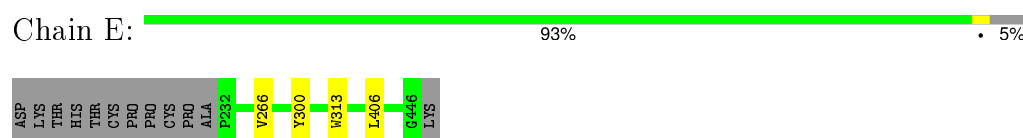
### 3 Residue-property plots [i](#)

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

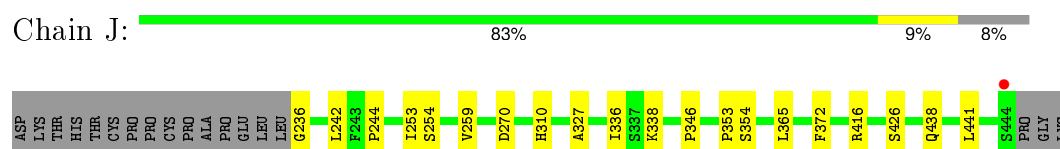
- Molecule 1: High affinity immunoglobulin gamma Fc receptor I



- Molecule 2: Ig gamma-1 chain C region



- Molecule 2: Ig gamma-1 chain C region





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	134.74Å 126.75Å 71.84Å 90.00° 118.44° 90.00°	Depositor
Resolution (Å)	86.55 – 2.42 86.55 – 2.42	Depositor EDS
% Data completeness (in resolution range)	98.6 (86.55-2.42) 98.7 (86.55-2.42)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.18 (at 2.42Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, $R_{free}$	0.201 , 0.254 0.228 , 0.275	Depositor DCC
$R_{free}$ test set	1946 reflections (5.12%)	DCC
Wilson B-factor (Å <sup>2</sup> )	48.0	Xtriage
Anisotropy	0.188	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 31.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 39992 reflections (0.003%)	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	5767	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.56% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: ZN, MAN, FUL, NAG

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	0.67	0/1976	0.72	0/2680
2	E	0.67	1/1759 (0.1%)	0.70	0/2397
2	J	0.66	0/1714	0.68	1/2335 (0.0%)
All	All	0.66	1/5449 (0.0%)	0.70	1/7412 (0.0%)

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
1	A	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	313	TRP	CD2-CE2	5.70	1.48	1.41

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	J	416	ARG	NE-CZ-NH2	-5.31	117.64	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	199	THR	Peptide
1	A	27	GLN	Peptide

## 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	1932	0	1887	18	0
2	E	1711	0	1679	2	0
2	J	1668	0	1633	14	0
3	A	98	0	87	0	0
3	E	56	0	49	0	0
3	J	56	0	49	1	0
4	A	22	0	20	0	0
4	E	33	0	26	0	0
4	J	33	0	26	0	0
5	A	1	0	0	0	0
5	E	1	0	0	0	0
5	J	1	0	0	0	0
6	E	10	0	10	0	0
6	J	10	0	10	1	0
7	A	16	0	0	1	0
7	E	64	0	0	0	0
7	J	55	0	0	1	0
All	All	5767	0	5476	33	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 3.

The worst 5 of 33 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)
2:J:244:PRO:HB3	2:J:336:ILE:HD11	1.78	0.66
1:A:67:THR:HG22	1:A:69:SER:H	1.60	0.66
1:A:109:VAL:HG12	1:A:122:LEU:HD23	1.77	0.65
1:A:107:LEU:HD21	1:A:166:TYR:HB3	1.80	0.64
1:A:174:HIS:ND1	7:A:401:HOH:O	2.32	0.59

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	234/267 (88%)	224 (96%)	7 (3%)	3 (1%)	15	20
2	E	213/227 (94%)	210 (99%)	3 (1%)	0	100	100
2	J	207/227 (91%)	201 (97%)	6 (3%)	0	100	100
All	All	654/721 (91%)	635 (97%)	16 (2%)	3 (0%)	34	47

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	28	PRO
1	A	239	ARG
1	A	67	THR

### 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	215/236 (91%)	210 (98%)	5 (2%)	58	77
2	E	199/210 (95%)	199 (100%)	0	100	100
2	J	194/210 (92%)	192 (99%)	2 (1%)	82	92
All	All	608/656 (93%)	601 (99%)	7 (1%)	78	90

5 of 7 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	202	LEU
2	J	354	SER
1	A	208	VAL
1	A	77	VAL
2	J	254	SER

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	206	ASN
2	E	361	ASN
2	E	389	ASN
2	E	438	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

15 non-standard protein/DNA/RNA residues 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	301	1,3	14,14,15	0.63	0	15,19,21	1.27	1 (6%)
3	NAG	A	302	3,4	14,14,15	0.82	0	15,19,21	2.43	6 (40%)
3	NAG	A	304	1	14,14,15	0.56	0	15,19,21	1.39	4 (26%)
3	NAG	A	305	1	14,14,15	1.24	1 (7%)	15,19,21	1.52	3 (20%)
3	NAG	A	306	1,4	14,14,15	0.63	0	15,19,21	1.32	2 (13%)
3	NAG	A	308	1,3	14,14,15	0.74	0	15,19,21	1.56	3 (20%)
3	NAG	A	309	3	14,14,15	1.39	1 (7%)	15,19,21	2.33	5 (33%)
3	NAG	E	501	3,2,6	14,14,15	0.50	0	15,19,21	1.87	3 (20%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	NAG	E	502	3,4	14,14,15	0.49	0	15,19,21	1.51	3 (20%)
3	NAG	E	505	4	14,14,15	0.52	0	15,19,21	0.86	0
3	NAG	E	507	4	14,14,15	0.77	1 (7%)	15,19,21	0.95	1 (6%)
3	NAG	J	501	3,2,6	14,14,15	0.54	0	15,19,21	1.89	4 (26%)
3	NAG	J	502	3,4	14,14,15	0.62	0	15,19,21	1.07	1 (6%)
3	NAG	J	505	4	14,14,15	0.53	0	15,19,21	1.10	1 (6%)
3	NAG	J	507	4	14,14,15	0.49	0	15,19,21	1.21	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	301	1,3	-	0/6/23/26	0/1/1/1
3	NAG	A	302	3,4	-	0/6/23/26	0/1/1/1
3	NAG	A	304	1	-	0/6/23/26	0/1/1/1
3	NAG	A	305	1	1/1/5/7	0/6/23/26	0/1/1/1
3	NAG	A	306	1,4	-	0/6/23/26	0/1/1/1
3	NAG	A	308	1,3	-	0/6/23/26	0/1/1/1
3	NAG	A	309	3	1/1/5/7	0/6/23/26	0/1/1/1
3	NAG	E	501	3,2,6	-	0/6/23/26	0/1/1/1
3	NAG	E	502	3,4	-	0/6/23/26	0/1/1/1
3	NAG	E	505	4	-	0/6/23/26	0/1/1/1
3	NAG	E	507	4	-	0/6/23/26	0/1/1/1
3	NAG	J	501	3,2,6	-	0/6/23/26	0/1/1/1
3	NAG	J	502	3,4	-	0/6/23/26	0/1/1/1
3	NAG	J	505	4	-	0/6/23/26	0/1/1/1
3	NAG	J	507	4	-	0/6/23/26	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	507	NAG	O5-C1	-2.01	1.40	1.43
3	A	309	NAG	C1-C2	3.14	1.56	1.52
3	A	305	NAG	C1-C2	3.49	1.57	1.52

The worst 5 of 38 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	306	NAG	O5-C5-C4	-3.31	104.65	110.13

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	J	501	NAG	C6-C5-C4	-3.30	104.72	112.99
3	E	507	NAG	C4-C3-C2	-2.58	107.34	111.34
3	A	302	NAG	C4-C3-C2	-2.39	107.63	111.34
3	J	502	NAG	O7-C7-C8	-2.38	117.68	122.07

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	309	NAG	C1
3	A	305	NAG	C1

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	J	501	NAG	1	0

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 28 ligands modelled in this entry, 3 are monoatomic - leaving 25 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	301	1,3	14,14,15	0.63	0	15,19,21	1.27	1 (6%)
3	NAG	A	302	3,4	14,14,15	0.82	0	15,19,21	2.43	6 (40%)
4	MAN	A	303	3	11,11,12	0.73	0	15,15,17	2.65	3 (20%)
3	NAG	A	304	1	14,14,15	0.56	0	15,19,21	1.39	4 (26%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	NAG	A	305	1	14,14,15	1.24	1 (7%)	15,19,21	1.52	3 (20%)
3	NAG	A	306	1,4	14,14,15	0.63	0	15,19,21	1.32	2 (13%)
4	MAN	A	307	3	11,11,12	0.60	0	15,15,17	0.80	0
3	NAG	A	308	1,3	14,14,15	0.74	0	15,19,21	1.56	3 (20%)
3	NAG	A	309	3	14,14,15	1.39	1 (7%)	15,19,21	2.33	5 (33%)
3	NAG	E	501	3,2,6	14,14,15	0.50	0	15,19,21	1.87	3 (20%)
3	NAG	E	502	3,4	14,14,15	0.49	0	15,19,21	1.51	3 (20%)
4	MAN	E	503	3,4	11,11,12	0.33	0	15,15,17	1.32	2 (13%)
4	MAN	E	504	3,4	11,11,12	0.82	0	15,15,17	0.98	1 (6%)
3	NAG	E	505	4	14,14,15	0.52	0	15,19,21	0.86	0
4	MAN	E	506	3,4	11,11,12	0.78	0	15,15,17	1.33	1 (6%)
3	NAG	E	507	4	14,14,15	0.77	1 (7%)	15,19,21	0.95	1 (6%)
6	FUL	E	508	3	10,10,11	0.87	0	13,14,16	2.69	5 (38%)
3	NAG	J	501	3,2,6	14,14,15	0.54	0	15,19,21	1.89	4 (26%)
3	NAG	J	502	3,4	14,14,15	0.62	0	15,19,21	1.07	1 (6%)
4	MAN	J	503	3,4	11,11,12	0.33	0	15,15,17	1.46	1 (6%)
4	MAN	J	504	3,4	11,11,12	0.69	0	15,15,17	1.09	1 (6%)
3	NAG	J	505	4	14,14,15	0.53	0	15,19,21	1.10	1 (6%)
4	MAN	J	506	3,4	11,11,12	0.64	0	15,15,17	0.95	0
3	NAG	J	507	4	14,14,15	0.49	0	15,19,21	1.21	1 (6%)
6	FUL	J	508	3	10,10,11	0.56	0	13,14,16	2.35	5 (38%)

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	301	1,3	-	0/6/23/26	0/1/1/1
3	NAG	A	302	3,4	-	0/6/23/26	0/1/1/1
4	MAN	A	303	3	-	0/2/19/22	0/1/1/1
3	NAG	A	304	1	-	0/6/23/26	0/1/1/1
3	NAG	A	305	1	1/1/5/7	0/6/23/26	0/1/1/1
3	NAG	A	306	1,4	-	0/6/23/26	0/1/1/1
4	MAN	A	307	3	-	0/2/19/22	0/1/1/1
3	NAG	A	308	1,3	-	0/6/23/26	0/1/1/1
3	NAG	A	309	3	1/1/5/7	0/6/23/26	0/1/1/1
3	NAG	E	501	3,2,6	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	E	502	3,4	-	0/6/23/26	0/1/1/1
4	MAN	E	503	3,4	1/1/4/5	0/2/19/22	0/1/1/1
4	MAN	E	504	3,4	-	0/2/19/22	0/1/1/1
3	NAG	E	505	4	-	0/6/23/26	0/1/1/1
4	MAN	E	506	3,4	-	0/2/19/22	0/1/1/1
3	NAG	E	507	4	-	0/6/23/26	0/1/1/1
6	FUL	E	508	3	-	0/0/17/20	0/1/1/1
3	NAG	J	501	3,2,6	-	0/6/23/26	0/1/1/1
3	NAG	J	502	3,4	-	0/6/23/26	0/1/1/1
4	MAN	J	503	3,4	1/1/4/5	0/2/19/22	0/1/1/1
4	MAN	J	504	3,4	-	0/2/19/22	0/1/1/1
3	NAG	J	505	4	-	0/6/23/26	0/1/1/1
4	MAN	J	506	3,4	-	0/2/19/22	0/1/1/1
3	NAG	J	507	4	-	0/6/23/26	0/1/1/1
6	FUL	J	508	3	-	0/0/17/20	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	507	NAG	O5-C1	-2.01	1.40	1.43
3	A	309	NAG	C1-C2	3.14	1.56	1.52
3	A	305	NAG	C1-C2	3.49	1.57	1.52

The worst 5 of 57 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	J	508	FUL	C1-C2-C3	-4.74	103.81	109.55
6	E	508	FUL	O3-C3-C2	-4.39	101.96	110.01
6	J	508	FUL	O5-C1-C2	-4.20	104.18	110.89
4	E	506	MAN	O2-C2-C3	-3.51	103.11	110.19
3	A	306	NAG	O5-C5-C4	-3.31	104.65	110.13

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	J	503	MAN	C1
4	E	503	MAN	C1
3	A	309	NAG	C1
3	A	305	NAG	C1

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	J	501	NAG	1	0
6	J	508	FUL	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 [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	242/267 (90%)	0.01	7 (2%) 55 54	20, 38, 58, 72	0
2	E	215/227 (94%)	-0.14	0 100 100	25, 34, 48, 60	0
2	J	209/227 (92%)	-0.34	1 (0%) 91 91	24, 35, 50, 73	0
All	All	666/721 (92%)	-0.15	8 (1%) 81 81	20, 35, 55, 73	0

The worst 5 of 8 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	64	GLN	3.0
1	A	57	PHE	2.7
2	J	444	SER	2.6
1	A	55	GLN	2.5
1	A	203	LEU	2.5

### 6.2 Non-standard residues in protein, DNA, RNA chains [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
3	NAG	J	502	14/15	0.94	0.17	2.44	51,53,58,59	0
3	NAG	E	501	14/15	0.94	0.16	2.37	39,49,59,62	0
3	NAG	J	501	14/15	0.96	0.18	1.64	46,53,60,70	0
3	NAG	A	301	14/15	0.94	0.18	0.43	64,76,85,91	0
3	NAG	J	507	14/15	0.93	0.14	0.19	58,64,68,69	0
3	NAG	A	308	14/15	0.90	0.15	0.18	79,99,106,120	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	NAG	A	304	14/15	0.65	0.36	-	104,121,131,131	0
3	NAG	J	505	14/15	0.80	0.23	-	100,113,124,124	0
3	NAG	A	306	14/15	0.89	0.15	-	73,89,102,118	0
3	NAG	E	502	14/15	0.96	0.18	-	41,45,49,50	0
3	NAG	A	305	14/15	0.60	0.29	-	85,108,125,126	0
3	NAG	E	505	14/15	0.86	0.17	-	95,104,118,124	0
3	NAG	A	302	14/15	0.82	0.21	-	75,105,119,122	0
3	NAG	A	309	14/15	0.81	0.36	-	126,140,145,150	0
3	NAG	E	507	14/15	0.95	0.14	-	50,59,64,65	0

### 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, 95<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
5	ZN	J	509	1/1	0.95	0.28	7.58	47,47,47,47	0
5	ZN	A	310	1/1	0.96	0.23	7.10	67,67,67,67	0
3	NAG	J	502	14/15	0.94	0.17	2.44	51,53,58,59	0
3	NAG	E	501	14/15	0.94	0.16	2.37	39,49,59,62	0
3	NAG	J	501	14/15	0.96	0.18	1.64	46,53,60,70	0
3	NAG	A	301	14/15	0.94	0.18	0.43	64,76,85,91	0
3	NAG	J	507	14/15	0.93	0.14	0.19	58,64,68,69	0
3	NAG	A	308	14/15	0.90	0.15	0.18	79,99,106,120	0
4	MAN	J	503	11/12	0.95	0.15	-	52,59,62,66	0
6	FUL	J	508	10/11	0.90	0.15	-	84,89,92,93	0
4	MAN	J	506	11/12	0.94	0.14	-	54,57,61,62	0
4	MAN	E	503	11/12	0.97	0.17	-	42,47,52,57	0
4	MAN	E	506	11/12	0.95	0.15	-	52,55,59,59	0
4	MAN	A	307	11/12	0.69	0.23	-	100,125,128,138	0
4	MAN	A	303	11/12	0.54	0.22	-	98,122,129,131	0
3	NAG	E	502	14/15	0.96	0.18	-	41,45,49,50	0
4	MAN	J	504	11/12	0.89	0.14	-	72,87,91,101	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	NAG	A	302	14/15	0.82	0.21	-	75,105,119,122	0
3	NAG	A	309	14/15	0.81	0.36	-	126,140,145,150	0
3	NAG	E	507	14/15	0.95	0.14	-	50,59,64,65	0
5	ZN	E	509	1/1	0.99	0.26	-	49,49,49,49	0
3	NAG	A	306	14/15	0.89	0.15	-	73,89,102,118	0
3	NAG	A	305	14/15	0.60	0.29	-	85,108,125,126	0
6	FUL	E	508	10/11	0.91	0.23	-	60,66,73,82	0
3	NAG	E	505	14/15	0.86	0.17	-	95,104,118,124	0
3	NAG	A	304	14/15	0.65	0.36	-	104,121,131,131	0
3	NAG	J	505	14/15	0.80	0.23	-	100,113,124,124	0
4	MAN	E	504	11/12	0.93	0.17	-	65,77,81,91	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.