



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 17, 2016 – 11:57 PM EDT

PDB ID : 5DJA
Title : Fc Heterodimer Design 9.1 Y407M + T366I
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Deposited on : 2015-09-01
Resolution : 2.90 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
<http://wwpdb.org/validation/2016/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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027257
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-20027257

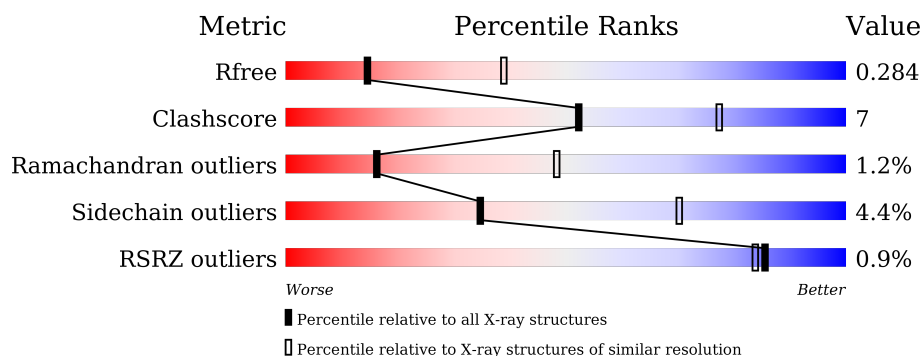
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

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	227	<div> <div>72%</div> <div>19%</div> <div>8%</div> </div>
2	B	240	<div> <div>68%</div> <div>18%</div> <div>14%</div> </div>
3	C	13	<div> <div>23%</div> <div>69%</div> <div>31%</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
7	GAL	A	506	-	-	-	X
7	GAL	B	506	-	-	-	X

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 3584 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 Ig gamma-1 chain C region.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	208	Total	C	N	O	S	1	0	0
			1634	1038	273	315	8			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	356	GLU	ASP	variant	UNP P01857
A	358	MET	LEU	variant	UNP P01857
A	407	MET	TYR	engineered mutation	UNP P01857

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	207	Total	C	N	O	S	0	0	0
			1634	1043	271	314	6			

There are 19 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	208	HIS	-	expression tag	UNP P01857
B	209	HIS	-	expression tag	UNP P01857
B	210	HIS	-	expression tag	UNP P01857
B	211	HIS	-	expression tag	UNP P01857
B	212	HIS	-	expression tag	UNP P01857
B	213	HIS	-	expression tag	UNP P01857
B	214	HIS	-	expression tag	UNP P01857
B	215	HIS	-	expression tag	UNP P01857
B	216	SER	-	expression tag	UNP P01857
B	217	GLY	-	expression tag	UNP P01857
B	218	SER	-	expression tag	UNP P01857
B	219	GLY	-	expression tag	UNP P01857
B	220	SER	-	expression tag	UNP P01857

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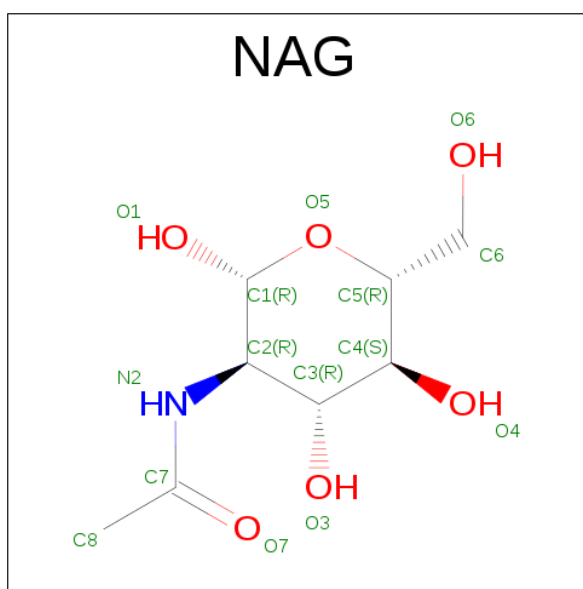
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Chain	Residue	Modelled	Actual	Comment	Reference
B	252	GLU	MET	engineered mutation	UNP P01857
B	253	ALA	ILE	engineered mutation	UNP P01857
B	356	GLU	ASP	variant	UNP P01857
B	358	MET	LEU	variant	UNP P01857
B	366	ILE	THR	engineered mutation	UNP P01857
B	435	ALA	HIS	engineered mutation	UNP P01857

- Molecule 3 is a protein called Fc-III peptide.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	13	Total	C	N	O	S	0	0	0
			103	67	17	17	2			

- Molecule 4 is N-ACETYL-D-GLUCOSAMINE (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



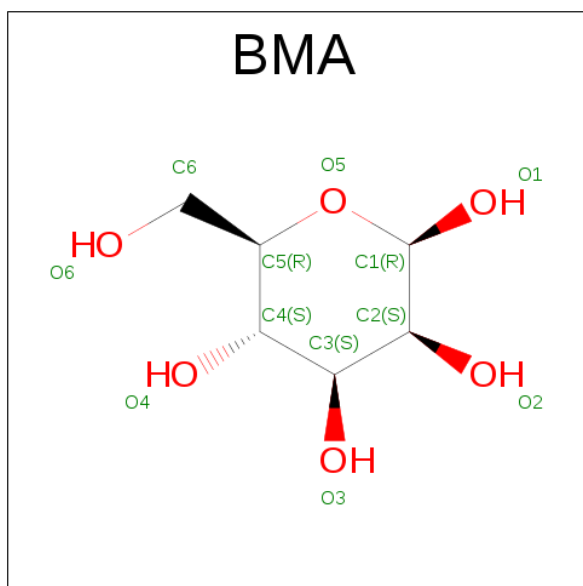
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	B	1	Total	C	N	O	0	0
			14	8	1	5		
4	B	1	Total	C	N	O	0	0
			14	8	1	5		

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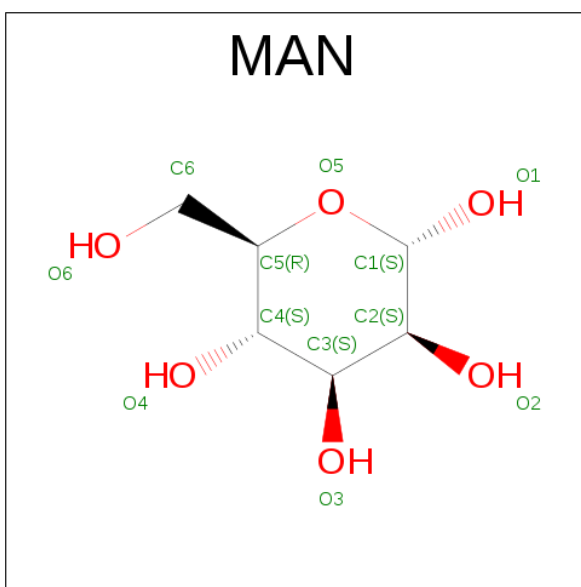
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	N	O	0	0
			14	8	1	5		

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



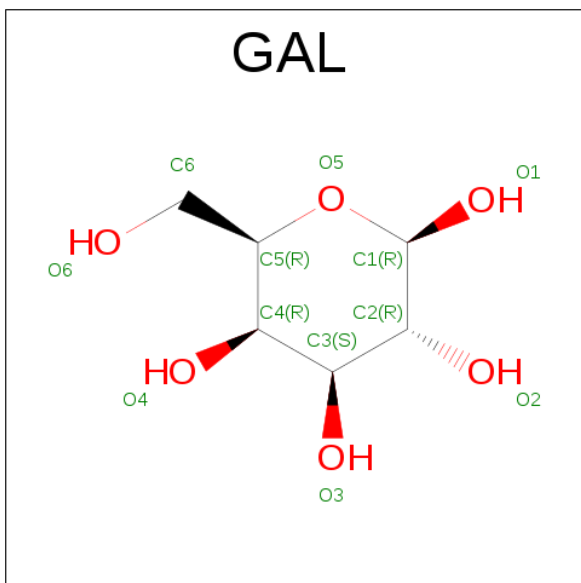
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			11	6	5		
5	B	1	Total	C	O	0	0
			11	6	5		

- Molecule 6 is ALPHA-D-MANNOSE (three-letter code: MAN) (formula: $C_6H_{12}O_6$).



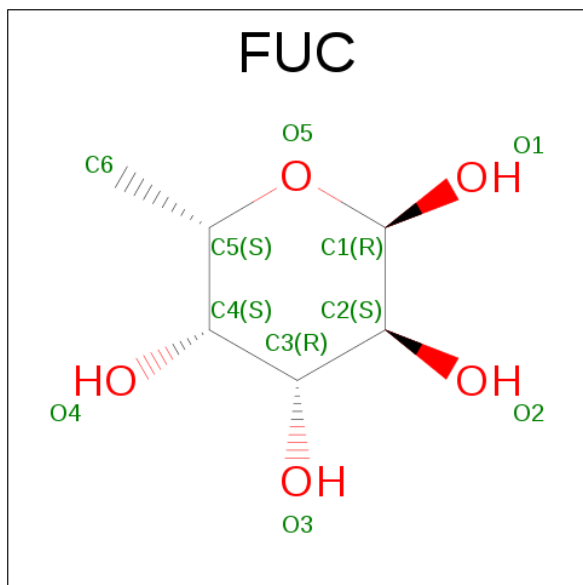
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			11	6	5		
6	A	1	Total	C	O	0	0
			11	6	5		
6	B	1	Total	C	O	0	0
			11	6	5		
6	B	1	Total	C	O	0	0
			11	6	5		

- Molecule 7 is BETA-D-GALACTOSE (three-letter code: GAL) (formula: C₆H₁₂O₆).



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

- Molecule 8 is ALPHA-L-FUCOSE (three-letter code: FUC) (formula: C₆H₁₂O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	O	0	0
			10	6	4		
8	B	1	Total	C	O	0	0
			10	6	4		

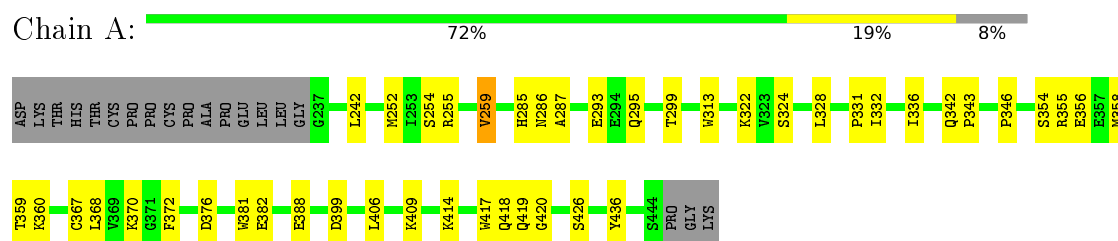
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	8	Total	O	0	0
			8	8		
9	B	13	Total	O	0	0
			13	13		

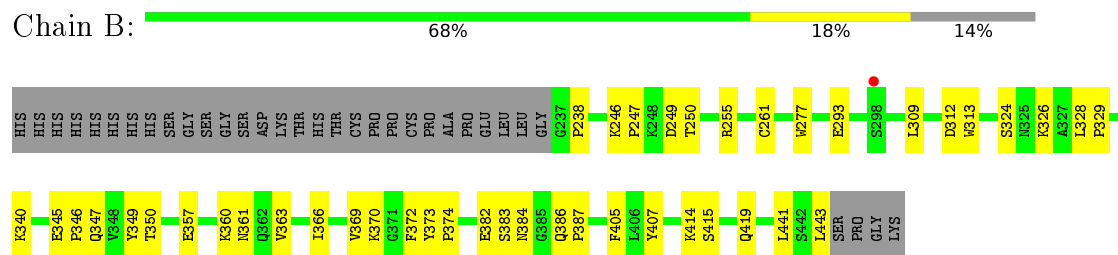
3 Residue-property plots [i](#)

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.

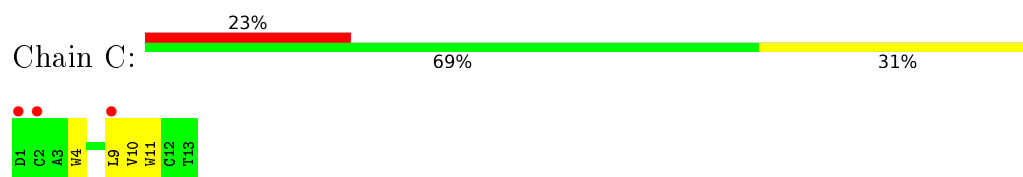
- Molecule 1: Ig gamma-1 chain C region



- Molecule 2: Ig gamma-1 chain C region



- Molecule 3: Fc-III peptide



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	60.66Å 67.20Å 78.61Å 90.00° 105.07° 90.00°	Depositor
Resolution (Å)	30.00 – 2.90 44.15 – 2.90	Depositor EDS
% Data completeness (in resolution range)	96.3 (30.00-2.90) 96.3 (44.15-2.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.19	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.80 (at 2.90Å)	Xtriage
Refinement program	REFMAC 5.7.0017	Depositor
R, R_{free}	0.207 , 0.281 0.210 , 0.284	Depositor DCC
R_{free} test set	666 reflections (5.32%)	DCC
Wilson B-factor (Å ²)	44.4	Xtriage
Anisotropy	0.137	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 36.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	3584	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.37% 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, 0.375 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: MAN, GAL, BMA, NAG, FUC

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	0.64	2/1678 (0.1%)	0.68	1/2289 (0.0%)
2	B	0.55	1/1679 (0.1%)	0.63	0/2290
3	C	1.02	2/107 (1.9%)	0.65	0/146
All	All	0.62	5/3464 (0.1%)	0.65	1/4725 (0.0%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	360	LYS	CG-CD	-13.21	1.07	1.52
1	A	417	TRP	CD2-CE2	5.88	1.48	1.41
3	C	4	TRP	CD2-CE2	5.12	1.47	1.41
3	C	11	TRP	CD2-CE2	5.04	1.47	1.41
2	B	313	TRP	CD2-CE2	5.02	1.47	1.41

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	360	LYS	CB-CG-CD	-12.94	77.96	111.60

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	1634	0	1576	28	0
2	B	1634	0	1586	21	0
3	C	103	0	89	2	0
4	A	42	0	35	0	0
4	B	42	0	35	0	0
5	A	11	0	8	0	0
5	B	11	0	8	0	0
6	A	22	0	19	0	0
6	B	22	0	19	0	0
7	A	11	0	10	0	0
7	B	11	0	10	0	0
8	A	10	0	10	0	0
8	B	10	0	10	0	0
9	A	8	0	0	0	0
9	B	13	0	0	1	0
All	All	3584	0	3415	47	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 7.

All (47) 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:B:238:PRO:HD2	2:B:328:LEU:HD13	1.68	0.76
1:A:409:LYS:HB2	2:B:407:TYR:OH	1.94	0.68
2:B:346:PRO:HB3	2:B:372:PHE:HB3	1.80	0.64
1:A:259:VAL:HG11	1:A:336:ILE:HD11	1.81	0.63
2:B:255:ARG:HD3	9:B:609:HOH:O	1.98	0.62
1:A:406:LEU:HD12	1:A:406:LEU:C	2.23	0.58
1:A:252:MET:HB2	1:A:255:ARG:HG3	1.87	0.57
2:B:357:GLU:O	2:B:360:LYS:HG2	2.06	0.56
2:B:346:PRO:CB	2:B:372:PHE:HB3	2.36	0.55
2:B:347:GLN:NE2	2:B:349:TYR:OH	2.38	0.55
2:B:246:LYS:HB2	2:B:249:ASP:OD2	2.08	0.54
1:A:367:CYS:HB2	1:A:381:TRP:CZ2	2.43	0.53
1:A:436:TYR:HB2	3:C:10:VAL:HG11	1.90	0.52
1:A:295:GLN:HB2	1:A:299:THR:O	2.11	0.50
1:A:354:SER:C	1:A:356:GLU:N	2.64	0.50
1:A:418:GLN:O	1:A:420:GLY:N	2.43	0.49
1:A:346:PRO:HB3	1:A:372:PHE:HB3	1.95	0.49
1:A:252:MET:CB	1:A:255:ARG:HG3	2.44	0.48
1:A:242:LEU:HD21	1:A:336:ILE:HG13	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:406:LEU:HD12	1:A:406:LEU:O	2.15	0.47
1:A:322:LYS:HE3	1:A:331:PRO:CB	2.44	0.47
2:B:247:PRO:O	2:B:250:THR:OG1	2.28	0.47
1:A:354:SER:C	1:A:356:GLU:H	2.18	0.46
2:B:361:ASN:N	2:B:361:ASN:HD22	2.12	0.46
1:A:313:TRP:CZ3	1:A:336:ILE:HD13	2.51	0.46
1:A:328:LEU:HD21	1:A:332:ILE:HD12	1.97	0.45
1:A:358:MET:O	1:A:414:LYS:HD2	2.16	0.45
2:B:350:THR:HB	2:B:441:LEU:HD22	1.98	0.45
1:A:285:HIS:C	1:A:287:ALA:H	2.21	0.45
1:A:418:GLN:C	1:A:420:GLY:H	2.20	0.44
2:B:369:VAL:O	2:B:405:PHE:HA	2.18	0.43
2:B:414:LYS:O	2:B:415:SER:C	2.57	0.43
2:B:383:SER:O	2:B:386:GLN:HB2	2.19	0.43
1:A:368:LEU:HD22	2:B:366:ILE:HD11	2.00	0.42
1:A:342:GLN:HA	1:A:343:PRO:HD3	1.87	0.42
1:A:254:SER:OG	3:C:9:LEU:O	2.35	0.42
2:B:386:GLN:HA	2:B:387:PRO:HD2	1.87	0.42
1:A:285:HIS:O	1:A:287:ALA:N	2.51	0.41
1:A:242:LEU:CD2	1:A:336:ILE:HG13	2.51	0.41
1:A:382:GLU:HA	1:A:388:GLU:HG2	2.02	0.41
2:B:261:CYS:HB2	2:B:277:TRP:CZ2	2.55	0.41
1:A:313:TRP:HZ3	1:A:336:ILE:HD13	1.85	0.41
2:B:345:GLU:HA	2:B:346:PRO:HD3	1.89	0.41
2:B:373:TYR:CG	2:B:374:PRO:HA	2.55	0.41
2:B:238:PRO:HG2	2:B:328:LEU:HD22	2.02	0.41
1:A:336:ILE:C	1:A:336:ILE:HD12	2.42	0.41
2:B:309:LEU:HB2	2:B:312:ASP:HB2	2.03	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	206/227 (91%)	194 (94%)	9 (4%)	3 (2%)	13	42
2	B	205/240 (85%)	186 (91%)	17 (8%)	2 (1%)	19	54
3	C	11/13 (85%)	11 (100%)	0	0	100	100
All	All	422/480 (88%)	391 (93%)	26 (6%)	5 (1%)	16	48

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	286	ASN
1	A	419	GLN
1	A	355	ARG
2	B	329	PRO
2	B	419	GLN

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	187/210 (89%)	179 (96%)	8 (4%)	35	71
2	B	186/219 (85%)	177 (95%)	9 (5%)	31	67
3	C	10/11 (91%)	10 (100%)	0	100	100
All	All	383/440 (87%)	366 (96%)	17 (4%)	35	70

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	259	VAL
1	A	293	GLU
1	A	324	SER
1	A	359	THR
1	A	370	LYS
1	A	376	ASP
1	A	399	ASP
1	A	426	SER

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Mol	Chain	Res	Type
2	B	293	GLU
2	B	324	SER
2	B	326	LYS
2	B	340	LYS
2	B	363	VAL
2	B	370	LYS
2	B	382	GLU
2	B	384	ASN
2	B	443	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	311	GLN
1	A	386	GLN
1	A	390	ASN
1	A	418	GLN
2	B	342	GLN
2	B	347	GLN
2	B	361	ASN
2	B	418	GLN

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	$\# Z > 2$
4	NAG	A	501	1,8,4	14,14,15	0.63	0	15,19,21	1.22	2 (13%)
4	NAG	A	502	5,4	14,14,15	0.55	0	15,19,21	1.13	2 (13%)
5	BMA	A	503	4,6	11,11,12	0.76	0	15,15,17	1.46	2 (13%)
6	MAN	A	504	5,4	11,11,12	0.55	0	15,15,17	1.28	1 (6%)
4	NAG	A	505	7,6	14,14,15	0.70	0	15,19,21	1.63	4 (26%)
7	GAL	A	506	4	11,11,12	0.80	0	15,15,17	1.64	1 (6%)
8	FUC	A	507	4	10,10,11	0.76	0	13,14,16	1.24	1 (7%)
6	MAN	A	508	5	11,11,12	0.66	0	15,15,17	1.58	4 (26%)
4	NAG	B	501	8,2,4	14,14,15	0.53	0	15,19,21	1.09	1 (6%)
4	NAG	B	502	5,4	14,14,15	0.48	0	15,19,21	1.03	1 (6%)
5	BMA	B	503	4,6	11,11,12	0.51	0	15,15,17	1.71	2 (13%)
6	MAN	B	504	5,4	11,11,12	0.67	0	15,15,17	0.76	0
4	NAG	B	505	7,6	14,14,15	0.65	0	15,19,21	0.81	0
7	GAL	B	506	4	11,11,12	0.87	0	15,15,17	1.40	1 (6%)
6	MAN	B	507	5	11,11,12	0.52	0	15,15,17	2.55	4 (26%)
8	FUC	B	508	4	10,10,11	0.68	0	13,14,16	0.77	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
4	NAG	A	501	1,8,4	-	0/6/23/26	0/1/1/1
4	NAG	A	502	5,4	-	0/6/23/26	0/1/1/1
5	BMA	A	503	4,6	-	0/2/19/22	0/1/1/1
6	MAN	A	504	5,4	-	0/2/19/22	0/1/1/1
4	NAG	A	505	7,6	-	0/6/23/26	0/1/1/1
7	GAL	A	506	4	-	0/2/19/22	0/1/1/1
8	FUC	A	507	4	-	0/0/17/20	0/1/1/1
6	MAN	A	508	5	-	0/2/19/22	0/1/1/1
4	NAG	B	501	8,2,4	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	B	502	5,4	-	0/6/23/26	0/1/1/1
5	BMA	B	503	4,6	-	0/2/19/22	0/1/1/1
6	MAN	B	504	5,4	-	0/2/19/22	0/1/1/1
4	NAG	B	505	7,6	-	0/6/23/26	0/1/1/1
7	GAL	B	506	4	-	0/2/19/22	0/1/1/1
6	MAN	B	507	5	-	0/2/19/22	0/1/1/1
8	FUC	B	508	4	-	0/0/17/20	0/1/1/1

There are no bond length outliers.

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	507	MAN	C1-C2-C3	-4.28	104.37	109.55
6	A	508	MAN	O5-C1-C2	-2.74	106.51	110.89
4	A	505	NAG	C3-C4-C5	-2.65	105.50	110.23
4	A	505	NAG	C1-O5-C5	-2.41	108.59	112.14
4	A	501	NAG	O7-C7-C8	-2.38	117.69	122.07
4	B	501	NAG	C3-C4-C5	-2.29	106.15	110.23
4	A	502	NAG	O7-C7-C8	-2.19	118.03	122.07
4	A	501	NAG	C3-C4-C5	-2.14	106.40	110.23
5	A	503	BMA	O5-C5-C6	2.05	111.74	107.34
4	A	505	NAG	O5-C5-C6	2.06	111.74	107.34
4	A	502	NAG	C1-O5-C5	2.13	115.28	112.14
6	A	508	MAN	O5-C5-C4	2.32	113.98	110.13
6	A	508	MAN	C2-C3-C4	2.44	115.30	111.05
4	B	502	NAG	C1-O5-C5	2.53	115.86	112.14
4	A	505	NAG	C4-C3-C2	3.13	116.19	111.34
6	A	508	MAN	C3-C4-C5	3.22	115.97	110.23
8	A	507	FUC	O5-C5-C6	3.48	112.53	106.28
6	B	507	MAN	C3-C4-C5	3.64	116.72	110.23
5	B	503	BMA	C1-C2-C3	3.69	114.02	109.55
6	A	504	MAN	C1-O5-C5	3.77	117.68	112.14
5	B	503	BMA	C1-O5-C5	4.02	118.05	112.14
6	B	507	MAN	O5-C5-C4	4.20	117.09	110.13
5	A	503	BMA	C1-C2-C3	4.21	114.65	109.55
7	B	506	GAL	C1-C2-C3	4.31	114.78	109.55
7	A	506	GAL	C1-C2-C3	5.20	115.85	109.55
6	B	507	MAN	C1-O5-C5	6.07	121.06	112.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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 [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	208/227 (91%)	-0.37	0 100 100	12, 34, 53, 73	1 (0%)
2	B	207/240 (86%)	-0.36	1 (0%) 91 90	13, 35, 57, 66	0
3	C	13/13 (100%)	1.11	3 (23%) 1 0	42, 67, 82, 86	0
All	All	428/480 (89%)	-0.32	4 (0%) 85 84	12, 35, 58, 86	1 (0%)

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	9	LEU	3.1
3	C	1	ASP	2.6
2	B	298	SER	2.4
3	C	2	CYS	2.2

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. 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, 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	LLDF	B-factors(\AA^2)	Q<0.9
7	GAL	B	506	11/12	0.79	0.37	7.76	59,77,90,93	0
7	GAL	A	506	11/12	0.87	0.32	7.62	66,83,91,103	0
4	NAG	A	505	14/15	0.92	0.17	0.31	41,46,81,82	0
4	NAG	B	505	14/15	0.95	0.17	-0.48	48,57,67,67	0
4	NAG	B	501	14/15	0.94	0.19	-0.57	41,51,58,60	0
4	NAG	A	502	14/15	0.95	0.13	-1.95	28,39,48,54	0
4	NAG	B	502	14/15	0.97	0.13	-2.05	28,38,47,51	0
6	MAN	B	504	11/12	0.94	0.12	-	39,45,54,58	0
8	FUC	B	508	10/11	0.94	0.14	-	46,49,56,62	0
4	NAG	A	501	14/15	0.92	0.19	-	37,50,61,65	0
5	BMA	A	503	11/12	0.93	0.10	-	37,40,53,56	0
6	MAN	A	508	11/12	0.89	0.12	-	45,57,67,73	0
8	FUC	A	507	10/11	0.93	0.23	-	56,65,79,81	0
6	MAN	B	507	11/12	0.88	0.13	-	44,50,56,56	0
6	MAN	A	504	11/12	0.90	0.15	-	40,55,64,68	0
5	BMA	B	503	11/12	0.98	0.09	-	36,40,42,45	0

6.5 Other polymers [i](#)

There are no such residues in this entry.