



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 12:00 pm GMT

PDB ID : 3MBW  
Title : Crystal structure of the human ephrin A2 LBD and CRD domains in complex with ephrin A1  
Authors : Walker, J.R.; Yermekbayeva, L.; Seitova, A.; Butler-Cole, C.; Bountra, C.; Weigelt, J.; Arrowsmith, C.H.; Edwards, A.M.; Bochkarev, A.; Dhe-Paganon, S.; Structural Genomics Consortium (SGC)  
Deposited on : 2010-03-26  
Resolution : 2.81 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](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	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	recalc28949

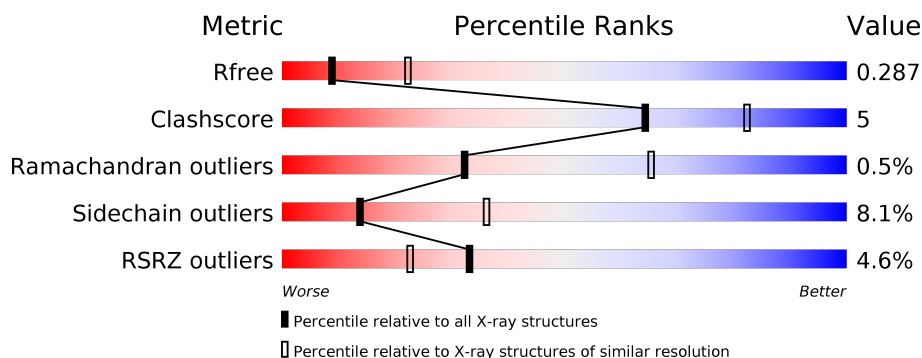
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.81 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))
$R_{free}$	100719	2917 (2.84-2.80)
Clashscore	112137	3382 (2.84-2.80)
Ramachandran outliers	110173	3324 (2.84-2.80)
Sidechain outliers	110143	3326 (2.84-2.80)
RSRZ outliers	101464	2948 (2.84-2.80)

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	330	<div> <div>5%</div> <div> <div></div> <div>74%</div> <div>12%</div> <div>•</div> <div>13%</div> </div> </div>
2	B	181	<div> <div>%</div> <div> <div></div> <div>54%</div> <div>15%</div> <div>•</div> <div>30%</div> </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	UNX	A	327	-	-	-	X

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3292 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 Ephrin type-A receptor 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	286	Total	C	N	O	S	0	1	0
			2160	1365	360	413	22			

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	ALA	-	EXPRESSION TAG	UNP P29317
A	-2	PRO	-	EXPRESSION TAG	UNP P29317
A	-1	GLU	-	EXPRESSION TAG	UNP P29317
A	0	HIS	-	EXPRESSION TAG	UNP P29317
A	1	HIS	-	EXPRESSION TAG	UNP P29317
A	2	HIS	-	EXPRESSION TAG	UNP P29317
A	3	HIS	-	EXPRESSION TAG	UNP P29317
A	4	HIS	-	EXPRESSION TAG	UNP P29317
A	5	HIS	-	EXPRESSION TAG	UNP P29317
A	6	ASP	-	EXPRESSION TAG	UNP P29317
A	7	TYR	-	EXPRESSION TAG	UNP P29317
A	8	ASP	-	EXPRESSION TAG	UNP P29317
A	9	ILE	-	EXPRESSION TAG	UNP P29317
A	10	PRO	-	EXPRESSION TAG	UNP P29317
A	11	THR	-	EXPRESSION TAG	UNP P29317
A	12	THR	-	EXPRESSION TAG	UNP P29317
A	13	GLU	-	EXPRESSION TAG	UNP P29317
A	14	ASN	-	EXPRESSION TAG	UNP P29317
A	15	LEU	-	EXPRESSION TAG	UNP P29317
A	16	TYR	-	EXPRESSION TAG	UNP P29317
A	17	PHE	-	EXPRESSION TAG	UNP P29317
A	18	GLN	-	EXPRESSION TAG	UNP P29317
A	19	GLY	-	EXPRESSION TAG	UNP P29317
A	20	ALA	-	EXPRESSION TAG	UNP P29317
A	21	MET	-	EXPRESSION TAG	UNP P29317
A	22	ASP	-	EXPRESSION TAG	UNP P29317

- Molecule 2 is a protein called Ephrin-A1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	127	Total	C	N	O	S	0	0	0
			1055	674	182	194	5			

There are 27 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-9	ALA	-	EXPRESSION TAG	UNP P20827
B	-8	PRO	-	EXPRESSION TAG	UNP P20827
B	-7	GLU	-	EXPRESSION TAG	UNP P20827
B	-6	HIS	-	EXPRESSION TAG	UNP P20827
B	-5	HIS	-	EXPRESSION TAG	UNP P20827
B	-4	HIS	-	EXPRESSION TAG	UNP P20827
B	-3	HIS	-	EXPRESSION TAG	UNP P20827
B	-2	HIS	-	EXPRESSION TAG	UNP P20827
B	-1	HIS	-	EXPRESSION TAG	UNP P20827
B	0	ASP	-	EXPRESSION TAG	UNP P20827
B	1	TYR	-	EXPRESSION TAG	UNP P20827
B	2	ASP	-	EXPRESSION TAG	UNP P20827
B	3	ILE	-	EXPRESSION TAG	UNP P20827
B	4	PRO	-	EXPRESSION TAG	UNP P20827
B	5	THR	-	EXPRESSION TAG	UNP P20827
B	6	THR	-	EXPRESSION TAG	UNP P20827
B	7	GLU	-	EXPRESSION TAG	UNP P20827
B	8	ASN	-	EXPRESSION TAG	UNP P20827
B	9	LEU	-	EXPRESSION TAG	UNP P20827
B	10	TYR	-	EXPRESSION TAG	UNP P20827
B	11	PHE	-	EXPRESSION TAG	UNP P20827
B	12	GLN	-	EXPRESSION TAG	UNP P20827
B	13	GLY	-	EXPRESSION TAG	UNP P20827
B	14	ALA	-	EXPRESSION TAG	UNP P20827
B	15	MET	-	EXPRESSION TAG	UNP P20827
B	16	ASP	-	EXPRESSION TAG	UNP P20827
B	159	VAL	ASP	SEE REMARK 999	UNP P20827

- Molecule 3 is UNKNOWN ATOM OR ION (three-letter code: UNX) (formula: X).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	X	0	0
			1	1		

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	4	Total	C	N	O	0	0
			49	28	2	19		

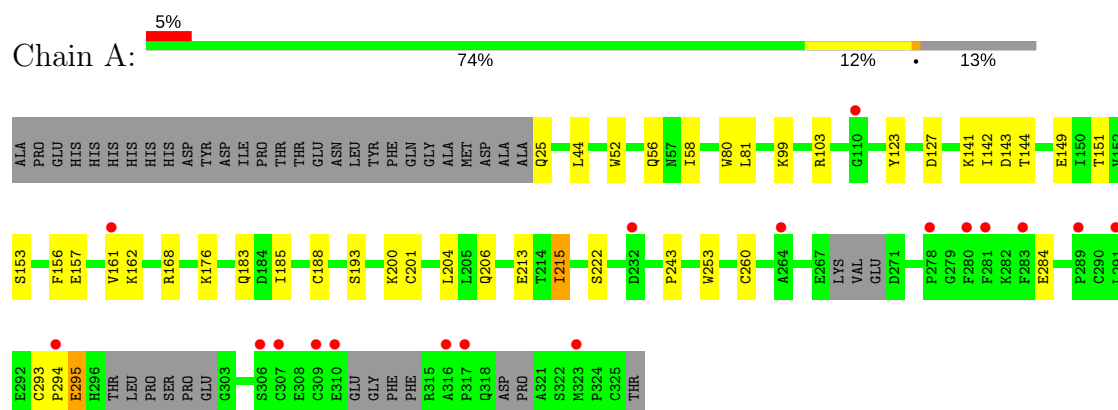
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	22	Total	O	0	0
			22	22		
5	B	5	Total	O	0	0
			5	5		

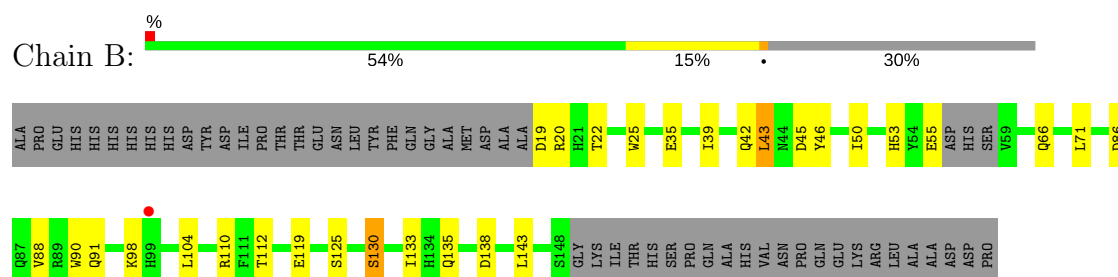
### 3 Residue-property plots [i](#)

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: Ephrin type-A receptor 2



#### • Molecule 2: Ephrin-A1



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	58.27Å 215.79Å 107.26Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.10 – 2.81 34.10 – 2.81	Depositor EDS
% Data completeness (in resolution range)	(Not available) (34.10-2.81) 99.7 (34.10-2.81)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.33 (at 2.81Å)	Xtriage
Refinement program	BUSTER 2.8.0	Depositor
R, $R_{free}$	0.226 , 0.272 0.238 , 0.287	Depositor DCC
$R_{free}$ test set	874 reflections (5.46%)	DCC
Wilson B-factor (Å <sup>2</sup> )	64.4	Xtriage
Anisotropy	0.327	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 68.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	3292	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.21% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



## 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: UNX, BMA, 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	0.46	0/2211	0.71	0/3004
2	B	0.44	0/1086	0.69	0/1471
All	All	0.45	0/3297	0.70	0/4475

There are no bond length outliers.

There are no bond angle outliers.

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	2160	0	1999	17	1
2	B	1055	0	973	16	0
3	A	1	0	0	0	1
4	B	49	0	40	4	0
5	A	22	0	0	0	0
5	B	5	0	0	0	0
All	All	3292	0	3012	30	1

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.

All (30) 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:53:HIS:H	4:B:501:NAG:H83	1.36	0.91
2:B:53:HIS:N	4:B:501:NAG:H83	2.07	0.69
1:A:123:TYR:HB3	1:A:142:ILE:HD11	1.78	0.65
2:B:53:HIS:O	4:B:501:NAG:H82	1.97	0.64
1:A:103:ARG:HA	1:A:151:THR:HB	1.84	0.59
2:B:22:THR:HG23	2:B:50:ILE:HD13	1.86	0.58
2:B:112:THR:HG22	2:B:119:GLU:HG3	1.86	0.57
2:B:66:GLN:HB3	2:B:133:ILE:HD12	1.90	0.54
2:B:20:ARG:HB3	2:B:46:TYR:HB2	1.90	0.53
1:A:52:TRP:CD1	1:A:81:LEU:HB2	2.44	0.52
1:A:99:LYS:HB2	1:A:193:SER:HB3	1.92	0.51
2:B:53:HIS:H	4:B:501:NAG:C8	2.15	0.51
2:B:39:ILE:HG12	2:B:143:LEU:HD11	1.94	0.50
1:A:149:GLU:HB3	1:A:162:LYS:HD3	1.93	0.50
1:A:58:ILE:HD12	2:B:88:VAL:HG12	1.94	0.49
2:B:25:TRP:NE1	2:B:130:SER:OG	2.45	0.48
1:A:200:LYS:HG2	1:A:215:ILE:HA	1.95	0.48
2:B:22:THR:HG23	2:B:50:ILE:CD1	2.43	0.48
1:A:243:PRO:HG3	1:A:260:CYS:SG	2.54	0.47
1:A:103:ARG:HB3	1:A:188:CYS:HB3	2.00	0.44
1:A:80:TRP:CE2	1:A:183:GLN:HG3	2.53	0.44
1:A:294:PRO:HA	1:A:295:GLU:HA	1.77	0.43
2:B:43:LEU:HD11	2:B:110:ARG:HB2	2.01	0.42
1:A:127:ASP:HA	1:A:176:LYS:HB3	2.02	0.42
1:A:56:GLN:HE22	2:B:104:LEU:HA	1.84	0.42
1:A:201:CYS:HB3	1:A:253:TRP:CE2	2.55	0.41
1:A:141:LYS:NZ	1:A:144:THR:OG1	2.41	0.41
1:A:56:GLN:HB3	2:B:90:TRP:CZ3	2.55	0.41
1:A:103:ARG:HD2	1:A:156:PHE:HZ	1.86	0.41
2:B:112:THR:HG22	2:B:119:GLU:CG	2.51	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:143:ASP:OD1	3:A:327:UNX:UNK[3_555]	2.05	0.15

## 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	277/330 (84%)	260 (94%)	16 (6%)	1 (0%)	38	70
2	B	123/181 (68%)	118 (96%)	4 (3%)	1 (1%)	22	53
All	All	400/511 (78%)	378 (94%)	20 (5%)	2 (0%)	32	65

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	86	ASP
1	A	222	SER

### 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	224/275 (82%)	210 (94%)	14 (6%)	21	50
2	B	112/165 (68%)	99 (88%)	13 (12%)	6	19
All	All	336/440 (76%)	309 (92%)	27 (8%)	14	37

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

Mol	Chain	Res	Type
1	A	25	GLN
1	A	44	LEU
1	A	153	SER
1	A	157	GLU

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Mol	Chain	Res	Type
1	A	161	VAL
1	A	168	ARG
1	A	185	ILE
1	A	204	LEU
1	A	206	GLN
1	A	213	GLU
1	A	215	ILE
1	A	284	GLU
1	A	293	CYS
1	A	295	GLU
2	B	19	ASP
2	B	35	GLU
2	B	42	GLN
2	B	43	LEU
2	B	45	ASP
2	B	55	GLU
2	B	71	LEU
2	B	91	GLN
2	B	98	LYS
2	B	125	SER
2	B	130	SER
2	B	135	GLN
2	B	138	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	57	ASN
1	A	259	GLN
2	B	74	HIS
2	B	81	GLN
2	B	91	GLN
2	B	135	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules 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 ⓘ

4 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$
4	NAG	B	501	2,4	14,14,15	0.96	0	15,19,21	2.03	4 (26%)
4	NAG	B	502	4	14,14,15	1.29	1 (7%)	15,19,21	1.97	5 (33%)
4	BMA	B	503	4	11,11,12	2.09	4 (36%)	13,15,17	2.53	7 (53%)
4	MAN	B	505	4	9,10,12	1.90	4 (44%)	13,14,17	1.93	4 (30%)

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	B	501	2,4	-	0/6/23/26	0/1/1/1
4	NAG	B	502	4	-	0/6/23/26	0/1/1/1
4	BMA	B	503	4	-	0/2/19/22	0/1/1/1
4	MAN	B	505	4	-	0/0/17/22	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	505	MAN	O2-C2	2.11	1.48	1.43
4	B	503	BMA	O2-C2	2.15	1.48	1.43
4	B	505	MAN	C4-C5	2.40	1.57	1.53
4	B	505	MAN	C1-C2	2.41	1.57	1.52
4	B	505	MAN	C4-C3	2.50	1.58	1.52
4	B	502	NAG	C4-C3	2.88	1.59	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	503	BMA	C4-C3	2.88	1.59	1.52
4	B	503	BMA	C4-C5	3.15	1.59	1.53
4	B	503	BMA	C2-C3	3.82	1.57	1.52

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	501	NAG	C4-C3-C2	-3.73	105.55	111.02
4	B	502	NAG	C2-N2-C7	-3.42	117.95	122.94
4	B	502	NAG	O3-C3-C2	-2.91	103.14	109.39
4	B	502	NAG	C1-O5-C5	-2.86	108.22	112.17
4	B	503	BMA	C6-C5-C4	2.00	117.69	113.00
4	B	501	NAG	C2-N2-C7	2.09	126.00	122.94
4	B	501	NAG	O3-C3-C4	2.11	114.96	110.36
4	B	502	NAG	C6-C5-C4	2.32	118.44	113.00
4	B	505	MAN	O3-C3-C4	2.37	115.52	110.36
4	B	503	BMA	C1-C2-C3	2.53	112.86	109.65
4	B	503	BMA	O2-C2-C3	2.57	115.23	110.17
4	B	505	MAN	C2-C3-C4	2.77	115.72	110.88
4	B	503	BMA	O3-C3-C4	2.86	116.58	110.36
4	B	505	MAN	C3-C4-C5	3.09	114.53	109.68
4	B	503	BMA	O2-C2-C1	3.12	115.52	109.18
4	B	503	BMA	C2-C3-C4	3.56	117.08	110.88
4	B	505	MAN	O2-C2-C1	3.73	116.76	109.18
4	B	502	NAG	C1-C2-N2	3.79	116.96	110.49
4	B	501	NAG	C1-C2-N2	5.43	119.77	110.49
4	B	503	BMA	O5-C1-C2	5.45	119.34	110.79

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	501	NAG	4	0

## 5.6 Ligand geometry

Of 1 ligands modelled in this entry, 1 is unknown - leaving 0 for Mogul analysis.

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.

No monomer is involved in short contacts.

## 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	286/330 (86%)	0.17	18 (6%) 21 12	29, 72, 137, 174	0
2	B	127/181 (70%)	-0.06	1 (0%) 86 81	53, 81, 109, 126	0
All	All	413/511 (80%)	0.10	19 (4%) 33 23	29, 78, 132, 174	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	310	GLU	4.5
1	A	317	PRO	3.7
1	A	283	PHE	3.3
1	A	291	LEU	3.3
1	A	306	SER	2.9
1	A	316	ALA	2.9
1	A	294	PRO	2.8
1	A	281	PHE	2.8
1	A	307	CYS	2.8
1	A	309	CYS	2.5
1	A	289	PRO	2.4
1	A	161	VAL	2.4
1	A	278	PRO	2.4
1	A	264	ALA	2.3
1	A	110	GLY	2.3
2	B	99	HIS	2.0
1	A	232	ASP	2.0
1	A	323	MET	2.0
1	A	280	PHE	2.0

### 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](#)

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
4	NAG	B	501	14/15	0.94	0.20	0.13	63,68,70,71	0
4	MAN	B	505	10/12	0.83	0.27	-	104,107,108,109	0
4	BMA	B	503	11/12	0.77	0.31	-	95,99,101,101	0
4	NAG	B	502	14/15	0.90	0.29	-	80,85,87,88	0

### 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(Å <sup>2</sup> )	Q<0.9
3	UNX	A	327	1/1	0.38	0.79	31.71	49,49,49,49	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.