



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:29 AM GMT

PDB ID : 2X97  
Title : Crystal structure of AnCE-RXP407 complex  
Authors : Akif, M.; Georgiadis, D.; Mahajan, A.; Dive, V.; Sturrock, E.D.; Isaac, R.E.; Acharya, K.R.  
Deposited on : 2010-03-14  
Resolution : 1.85 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 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
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) : trunk26865

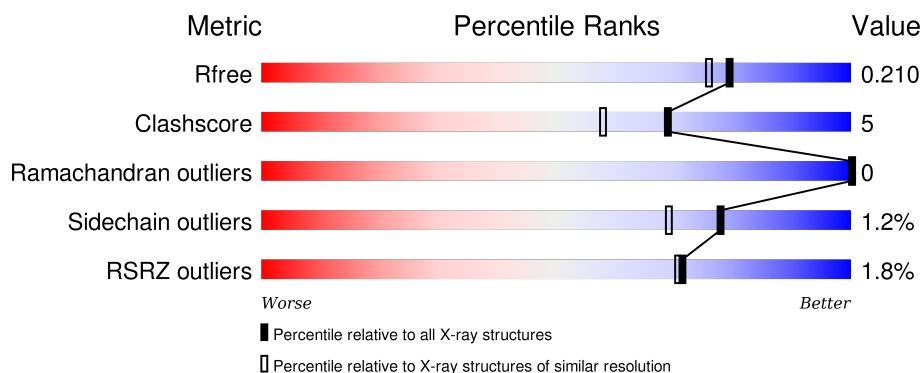
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.85 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	1745 (1.86-1.86)
Clashscore	102246	1898 (1.86-1.86)
Ramachandran outliers	100387	1875 (1.86-1.86)
Sidechain outliers	100360	1875 (1.86-1.86)
RSRZ outliers	91569	1747 (1.86-1.86)

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	598	

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
5	NAG	A	1623	-	-	-	X

## 2 Entry composition [i](#)

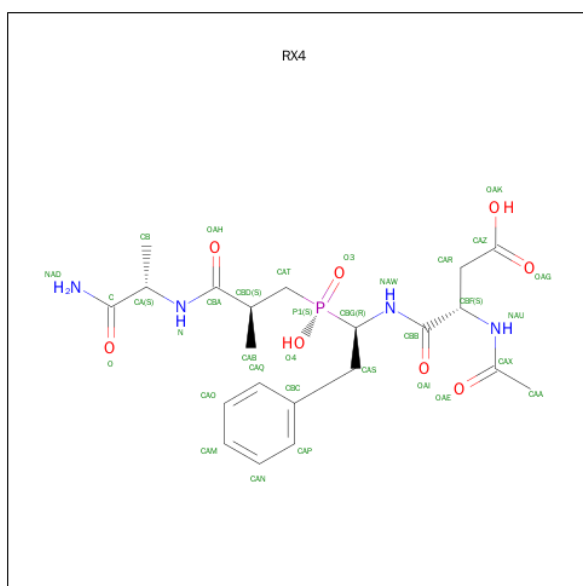
There are 6 unique types of molecules in this entry. The entry contains 5756 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 ANGIOTENSIN CONVERTING ENZYME.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	598	Total	C	N	O	S	0	2	0
			4883	3124	806	933	20			

- Molecule 2 is N 2 -ACETYL-N-{(1R)-1-[(S)-[(2S)-3-{[(2S)-1-AMINO-1-OXOPROPAN-2-YL]AMINO}-2-METHYL-3-OXOPROPYL](HYDROXY)PHOSPHORYL]-2-PHENYLETHYL}-L-ALPHA-ASPARAGINE (three-letter code: RX4) (formula: C<sub>21</sub>H<sub>31</sub>N<sub>4</sub>O<sub>8</sub>P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			34	21	4	8	1		

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 4 is a polymer of unknown type called SUGAR (6-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	6	Total	C	N	O	0	0
			72	40	2	30		

- Molecule 5 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula:  $C_8H_{15}NO_6$ ).



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

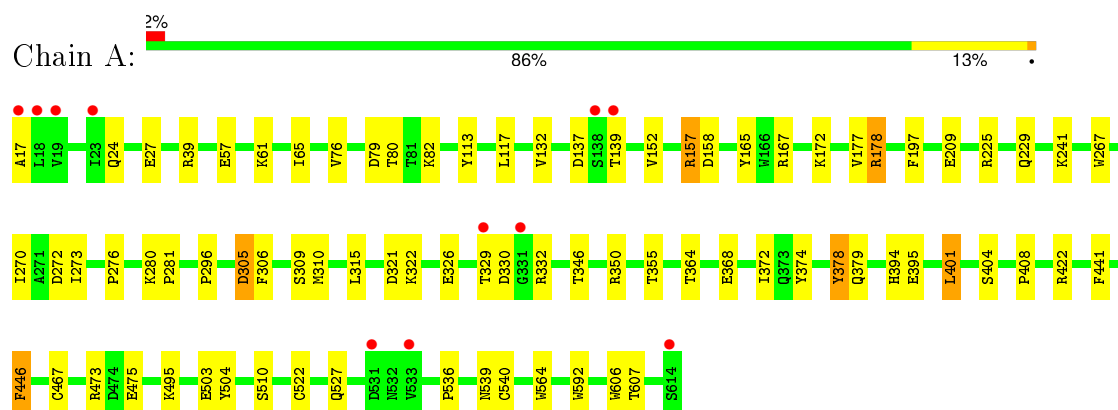
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	738	Total	O	0	0
			738	738		

### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $\text{RSRZ} > 2$ ). 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: ANGIOTENSIN CONVERTING ENZYME



## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	173.09 Å   173.09 Å   101.90 Å 90.00°   90.00°   120.00°	Depositor
Resolution (Å)	31.62 – 1.85 30.19 – 1.85	Depositor EDS
% Data completeness (in resolution range)	93.3 (31.62-1.85) 73.8 (30.19-1.85)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.57 (at 1.85 Å)	Xtriage
Refinement program	REFMAC 5.5.0072	Depositor
R, $R_{free}$	0.172 , 0.201 0.182 , 0.210	Depositor DCC
$R_{free}$ test set	3564 reflections (5.25%)	DCC
Wilson B-factor (Å <sup>2</sup> )	26.2	Xtriage
Anisotropy	0.051	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 46.9	EDS
Estimated twinning fraction	0.017 for -2/3*h-1/3*k-4/3*l,-1/3*h-2/3*k+4/3*l,-1/3*h+1/3*k+1/3*l 0.018 for -h,1/3*h-1/3*k-4/3*l,-1/3*h-2/3*k+1/3*l 0.011 for -1/3*h+1/3*k+4/3*l,-k,2/3*h+1/3*k+1/3*l 0.014 for -h,2/3*h+1/3*k+4/3*l,1/3*h+2/3*k-1/3*l 0.028 for -1/3*h-2/3*k+4/3*l,-2/3*h-1/3*k-4/3*l,1/3*h-1/3*k-1/3*l 0.026 for 1/3*h+2/3*k-4/3*l,-k,-2/3*h-1/3*k-1/3*l 0.116 for h,-h-k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 71479 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	5756	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.86% 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: MAN, ZN, BMA, NAG, RX4

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	1.34	21/5015 (0.4%)	1.02	13/6796 (0.2%)

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	24	GLN	CB-CG	-9.89	1.25	1.52
1	A	467	CYS	CB-SG	9.24	1.98	1.82
1	A	157	ARG	CZ-NH1	7.68	1.43	1.33
1	A	346	THR	CB-OG1	6.54	1.56	1.43
1	A	27	GLU	CG-CD	6.05	1.61	1.51
1	A	522	CYS	CB-SG	-5.91	1.72	1.81
1	A	378	TYR	CE2-CZ	-5.87	1.30	1.38
1	A	527	GLN	CD-OE1	5.75	1.36	1.24
1	A	57	GLU	CG-CD	5.69	1.60	1.51
1	A	378	TYR	CE1-CZ	5.60	1.45	1.38
1	A	379	GLN	CD-OE1	5.45	1.35	1.24
1	A	441	PHE	CD1-CE1	5.42	1.50	1.39
1	A	209	GLU	CG-CD	5.31	1.59	1.51
1	A	504	TYR	CE1-CZ	5.21	1.45	1.38
1	A	446	PHE	CE1-CZ	5.14	1.47	1.37
1	A	475	GLU	CG-CD	5.14	1.59	1.51
1	A	607	THR	CB-OG1	5.11	1.53	1.43
1	A	197	PHE	CD2-CE2	5.11	1.49	1.39
1	A	564	TRP	CZ3-CH2	5.09	1.48	1.40
1	A	267	TRP	CE3-CZ3	5.07	1.47	1.38
1	A	606	TRP	CE3-CZ3	5.00	1.47	1.38

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	157	ARG	NE-CZ-NH2	-14.65	112.98	120.30
1	A	157	ARG	NE-CZ-NH1	8.24	124.42	120.30
1	A	177	VAL	CG1-CB-CG2	7.83	123.42	110.90
1	A	167	ARG	NE-CZ-NH1	7.50	124.05	120.30
1	A	39	ARG	NE-CZ-NH1	7.29	123.95	120.30
1	A	178	ARG	NE-CZ-NH2	-7.21	116.69	120.30
1	A	473	ARG	NE-CZ-NH1	-6.86	116.87	120.30
1	A	401	LEU	CB-CG-CD2	5.89	121.02	111.00
1	A	241	LYS	CD-CE-NZ	-5.80	98.35	111.70
1	A	305	ASP	CB-CG-OD1	5.67	123.40	118.30
1	A	178	ARG	NE-CZ-NH1	5.59	123.10	120.30
1	A	158	ASP	CB-CG-OD1	5.47	123.23	118.30
1	A	310	MET	CG-SD-CE	-5.10	92.03	100.20

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	4883	0	4677	45	0
2	A	34	0	29	1	0
3	A	1	0	0	0	0
4	A	72	0	61	0	0
5	A	28	0	26	0	0
6	A	738	0	0	6	0
All	All	5756	0	4793	45	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.

All (45) 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:322:LYS:HB2	1:A:350:ARG:HD3	1.34	1.08

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:322:LYS:CB	1:A:350:ARG:HD3	1.86	1.06
1:A:270:ILE:HB	1:A:273:ILE:HD11	1.40	1.01
1:A:178:ARG:NH1	6:A:2267:HOH:O	2.03	0.87
1:A:322:LYS:HB3	1:A:350:ARG:HD3	1.67	0.77
1:A:61:LYS:O	1:A:65:ILE:HG12	1.89	0.72
1:A:270:ILE:HB	1:A:273:ILE:CD1	2.19	0.70
1:A:296:PRO:HA	1:A:355:THR:HG21	1.77	0.65
1:A:305:ASP:OD1	6:A:2448:HOH:O	2.13	0.65
1:A:270:ILE:CB	1:A:273:ILE:HD11	2.21	0.62
1:A:270:ILE:O	1:A:273:ILE:HG12	1.99	0.61
1:A:270:ILE:O	1:A:273:ILE:CG1	2.48	0.61
1:A:306:PHE:CD2	1:A:401:LEU:HD13	2.37	0.59
1:A:276:PRO:HB3	1:A:592:TRP:CH2	2.40	0.57
1:A:157:ARG:NH2	1:A:272:ASP:OD1	2.38	0.57
1:A:79:ASP:HA	1:A:82:LYS:HD3	1.87	0.57
1:A:321:ASP:OD1	6:A:2461:HOH:O	2.18	0.57
1:A:296:PRO:HG2	1:A:326:GLU:HG2	1.89	0.55
1:A:330:ASP:OD1	1:A:332:ARG:HD3	2.08	0.54
1:A:422:ARG:NH2	6:A:2550:HOH:O	2.31	0.53
1:A:17:ALA:HB3	6:A:2002:HOH:O	2.10	0.52
1:A:270:ILE:O	1:A:273:ILE:HG13	2.12	0.49
1:A:296:PRO:HG3	1:A:355:THR:HG22	1.96	0.48
1:A:296:PRO:CA	1:A:355:THR:HG21	2.44	0.47
1:A:395:GLU:HB2	1:A:510:SER:HB2	1.97	0.47
1:A:422:ARG:O	1:A:422:ARG:HG2	2.15	0.47
1:A:137:ASP:OD1	1:A:139:THR:HB	2.15	0.47
1:A:422:ARG:NH2	1:A:536:PRO:HG3	2.30	0.46
1:A:280:LYS:HB3	1:A:281:PRO:HD2	1.97	0.46
1:A:306:PHE:HZ	1:A:404:SER:HG	1.59	0.46
1:A:306:PHE:CZ	1:A:404:SER:OG	2.70	0.45
1:A:76:VAL:O	1:A:80:THR:HG23	2.16	0.45
1:A:132:VAL:HG12	1:A:172:LYS:HE3	1.98	0.45
1:A:296:PRO:HG3	1:A:355:THR:CG2	2.47	0.44
1:A:225:ARG:O	1:A:229:GLN:HG2	2.18	0.44
1:A:408:PRO:HD2	1:A:539:ASN:OD1	2.18	0.44
1:A:495:LYS:NZ	2:A:1615:RX4:HAD2	2.15	0.43
1:A:329:THR:HB	6:A:2473:HOH:O	2.18	0.43
1:A:364:THR:O	1:A:368:GLU:HG2	2.18	0.43
1:A:305:ASP:O	1:A:309:SER:HB3	2.19	0.43
1:A:270:ILE:HB	1:A:273:ILE:CG1	2.49	0.42
1:A:113:TYR:CE2	1:A:117:LEU:HD11	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:137:ASP:C	1:A:139:THR:H	2.21	0.42
1:A:315:LEU:CD2	1:A:372:ILE:HG21	2.49	0.42
1:A:152:VAL:HG11	1:A:165:TYR:CD2	2.56	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	598/598 (100%)	586 (98%)	12 (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	521/520 (100%)	515 (99%)	6 (1%)	78	69

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	374	TYR
1	A	378	TYR
1	A	394	HIS

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Mol	Chain	Res	Type
1	A	446	PHE
1	A	503	GLU
1	A	540	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 ⓘ

6 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	A	1617	1,4	14,14,15	0.88	0	15,19,21	1.32	2 (13%)
4	NAG	A	1618	4	14,14,15	0.77	0	15,19,21	2.27	6 (40%)
4	BMA	A	1619	4	11,11,12	0.58	0	14,15,17	1.20	1 (7%)
4	BMA	A	1620	4	11,11,12	0.90	0	14,15,17	2.82	5 (35%)
4	MAN	A	1621	4	11,11,12	1.57	2 (18%)	14,15,17	2.84	9 (64%)
4	MAN	A	1624	4	11,11,12	0.78	0	14,15,17	1.11	1 (7%)

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	1617	1,4	-	0/6/23/26	0/1/1/1
4	NAG	A	1618	4	-	0/6/23/26	0/1/1/1
4	BMA	A	1619	4	-	0/2/19/22	0/1/1/1
4	BMA	A	1620	4	-	0/2/19/22	1/1/1/1
4	MAN	A	1621	4	-	0/2/19/22	0/1/1/1
4	MAN	A	1624	4	-	0/2/19/22	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1621	MAN	C1-C2	2.60	1.58	1.52
4	A	1621	MAN	C2-C3	3.80	1.57	1.52

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1620	BMA	O2-C2-C3	-4.57	100.92	110.12
4	A	1620	BMA	C2-C3-C4	-4.51	103.38	111.04
4	A	1618	NAG	O6-C6-C5	-4.05	97.96	111.33
4	A	1621	MAN	C1-O5-C5	-4.00	107.17	112.25
4	A	1620	BMA	C3-C4-C5	-3.76	103.64	110.20
4	A	1618	NAG	C6-C5-C4	-3.33	104.79	113.02
4	A	1618	NAG	O4-C4-C5	-2.87	101.63	109.24
4	A	1621	MAN	O4-C4-C3	-2.63	104.41	110.34
4	A	1617	NAG	O6-C6-C5	-2.07	104.51	111.33
4	A	1621	MAN	C6-C5-C4	2.17	118.37	113.02
4	A	1621	MAN	O5-C1-C2	2.22	114.46	110.86
4	A	1624	MAN	C1-O5-C5	2.41	115.31	112.25
4	A	1621	MAN	O4-C4-C5	2.58	116.09	109.24
4	A	1621	MAN	O2-C2-C1	2.65	114.51	109.21
4	A	1618	NAG	C2-N2-C7	2.82	126.66	123.04
4	A	1619	BMA	O5-C5-C6	2.88	113.58	107.35
4	A	1617	NAG	C1-O5-C5	3.23	116.35	112.25
4	A	1618	NAG	C3-C2-N2	3.30	118.47	110.56
4	A	1621	MAN	O5-C5-C6	3.59	115.11	107.35
4	A	1620	BMA	C1-C2-C3	3.63	113.84	109.54
4	A	1618	NAG	C1-O5-C5	3.81	117.08	112.25
4	A	1621	MAN	O3-C3-C2	3.88	117.02	110.00
4	A	1620	BMA	O3-C3-C4	5.33	122.34	110.34
4	A	1621	MAN	C1-C2-C3	5.85	116.46	109.54

There are no chirality outliers.

There are no torsion outliers.

All (1) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1620	BMA	C1-C2-C3-C4-C5-O5

No monomer is involved in short contacts.

## 5.6 Ligand geometry

Of 4 ligands modelled in this entry, 1 is monoatomic - leaving 3 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	RX4	A	1615	3	27,34,34	1.53	6 (22%)	30,47,47	2.15	8 (26%)
5	NAG	A	1622	1	14,14,15	0.64	0	15,19,21	1.79	3 (20%)
5	NAG	A	1623	1	14,14,15	0.55	0	15,19,21	0.69	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
2	RX4	A	1615	3	-	0/36/43/43	0/1/1/1
5	NAG	A	1622	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1623	1	-	0/6/23/26	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1615	RX4	CBF-NAU	2.27	1.51	1.45
2	A	1615	RX4	O-C	2.38	1.28	1.23
2	A	1615	RX4	CAB-CBD	2.45	1.61	1.52
2	A	1615	RX4	CA-N	2.69	1.51	1.46
2	A	1615	RX4	CAA-CAX	2.78	1.56	1.50
2	A	1615	RX4	CBD-CBA	3.57	1.58	1.51

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1615	RX4	O4-P1-O3	-8.16	101.18	113.72
2	A	1615	RX4	CAM-CAN-CAP	-3.99	114.34	120.19
2	A	1615	RX4	CAB-CBD-CBA	-3.98	103.71	109.34
5	A	1622	NAG	C4-C3-C2	-3.21	106.25	111.23
2	A	1615	RX4	CBB-CBF-NAU	-2.48	104.26	111.26
2	A	1615	RX4	CAS-CBC-CAP	-2.37	115.94	120.90
2	A	1615	RX4	CAO-CAQ-CBC	-2.21	117.13	120.65
2	A	1615	RX4	CAQ-CBC-CAP	2.02	121.36	118.13
2	A	1615	RX4	CAO-CAM-CAN	2.16	123.72	119.93
5	A	1622	NAG	C1-O5-C5	3.39	116.55	112.25
5	A	1622	NAG	C3-C2-N2	3.68	119.38	110.56

There are no chirality outliers.

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
2	A	1615	RX4	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	598/598 (100%)	-0.50	11 (1%) 71 71	16, 25, 40, 52	7 (1%)

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	533	VAL	3.9
1	A	139	THR	3.8
1	A	329	THR	3.4
1	A	19	VAL	3.2
1	A	18	LEU	2.9
1	A	331	GLY	2.8
1	A	614	SER	2.6
1	A	531	ASP	2.3
1	A	23	ILE	2.2
1	A	17	ALA	2.1
1	A	138	SER	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( $\text{\AA}^2$ )	Q<0.9
4	NAG	A	1617	14/15	0.98	0.06	-0.78	22,26,29,30	0
4	MAN	A	1621	11/12	0.84	0.31	-	64,68,70,71	0
4	NAG	A	1618	14/15	0.99	0.10	-	24,29,37,40	0
4	MAN	A	1624	11/12	0.94	0.15	-	43,45,49,51	0
4	BMA	A	1620	11/12	0.87	0.14	-	52,57,66,68	0
4	BMA	A	1619	11/12	0.97	0.12	-	38,43,53,58	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( $\text{\AA}^2$ )	Q<0.9
5	NAG	A	1623	14/15	0.81	0.43	7.86	44,45,45,45	14
2	RX4	A	1615	34/34	0.99	0.08	-0.39	17,23,36,38	0
3	ZN	A	1616	1/1	1.00	0.07	-	22,22,22,22	0
5	NAG	A	1622	14/15	0.73	0.55	-	39,42,46,46	14

## 6.5 Other polymers [i](#)

There are no such residues in this entry.