



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 01:45 PM GMT

PDB ID : 2C3A  
Title : STRUCTURE OF UNLIGANDED HSV GD REVEALS A MECHANISM  
FOR RECEPTOR-MEDIATED ACTIVATION OF VIRUS ENTRY  
Authors : Krummenacher, C.; Supekár, V.M.; Whitbeck, J.C.; Lazear, E.; Connolly,  
S.A.; Eisenberg, R.J.; Cohen, G.H.; Wiley, D.C.; Carfi, A.  
Deposited on : 2005-10-05  
Resolution : 2.50 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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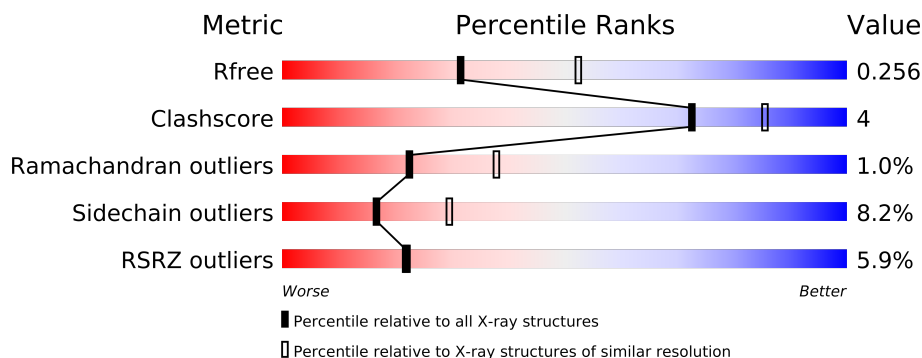
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.50 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))
$R_{free}$	66092	2784 (2.50-2.50)
Clashscore	79885	3562 (2.50-2.50)
Ramachandran outliers	78287	3480 (2.50-2.50)
Sidechain outliers	78261	3482 (2.50-2.50)
RSRZ outliers	66119	2785 (2.50-2.50)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	286	
1	B	286	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
4	ZN	B	1313	-	X

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 4185 atoms, of which 0 are hydrogen and 0 are deuterium.

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 GLYCOPROTEIN D.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	253	Total	C	N	O	S	0	0	0
			1959	1261	331	355	12			
1	B	274	Total	C	N	O	S	0	0	0
			2108	1358	350	388	12			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	37	CYS	VAL	ENGINEERED MUTATION	UNP P57083
A	302	CYS	ALA	ENGINEERED MUTATION	UNP P57083
B	37	CYS	VAL	ENGINEERED MUTATION	UNP P57083
B	302	CYS	ALA	ENGINEERED MUTATION	UNP P57083

- Molecule 2 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	2	Total	C	N	O	0	0
			28	16	2	10		
2	B	2	Total	C	N	O	0	0
			28	16	2	10		

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	37	CYS	VAL	ENGINEERED MUTATION	UNP P57083
A	302	CYS	ALA	ENGINEERED MUTATION	UNP P57083
B	37	CYS	VAL	ENGINEERED MUTATION	UNP P57083
B	302	CYS	ALA	ENGINEERED MUTATION	UNP P57083

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Cl 1 1	0	0
3	A	1	Total Cl 1 1	0	0

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	4	Total Zn 4 4	0	0
4	A	1	Total Zn 1 1	0	0

- Molecule 5 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	2	Total Na 2 2	0	0

- Molecule 6 is water.

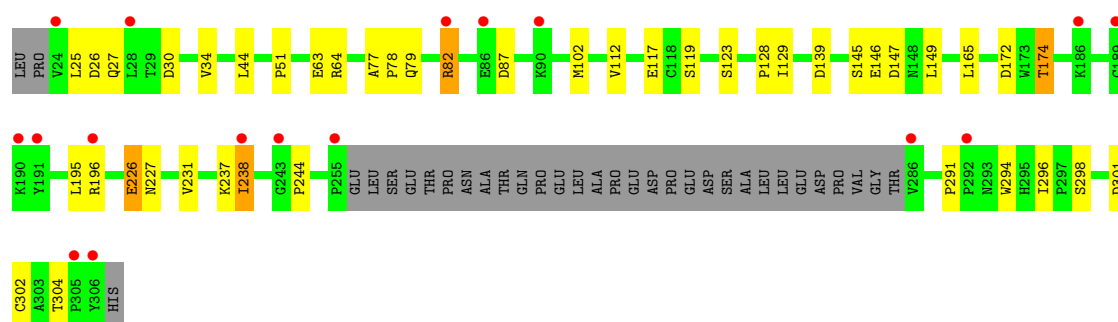
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	18	Total O 18 18	0	0
6	B	35	Total O 35 35	0	0

### 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 ( $RSRZ > 2$ ). 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.

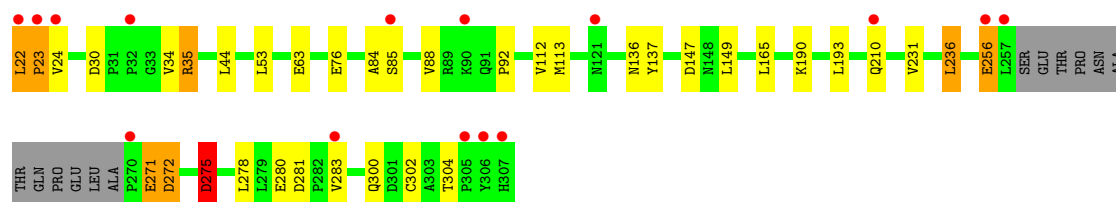
#### • Molecule 1: GLYCOPROTEIN D

Chain A: 



#### • Molecule 1: GLYCOPROTEIN D

Chain B: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	67.63Å 79.11Å 123.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 2.50 19.95 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.7 (25.00-2.50) 99.8 (19.95-2.50)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.92 (at 2.50Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, $R_{free}$	0.226 , 0.275 0.223 , 0.256	Depositor DCC
$R_{free}$ test set	1212 reflections (5.44%)	DCC
Wilson B-factor (Å <sup>2</sup> )	48.5	Xtriage
Anisotropy	0.284	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 18.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 23483 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	4185	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.17% of the height of the origin peak. No significant pseudotranslation is detected.*

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

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NA, ZN, NAG, CL

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	0.85	1/2015 (0.0%)	0.86	7/2763 (0.3%)
1	B	0.81	0/2174	0.89	5/2987 (0.2%)
All	All	0.83	1/4189 (0.0%)	0.88	12/5750 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	226	GLU	CA-CB	-17.04	1.16	1.53

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	301	ASP	CB-CG-OD2	10.58	127.82	118.30
1	A	226	GLU	N-CA-CB	8.61	126.09	110.60
1	B	281	ASP	CB-CG-OD2	6.54	124.19	118.30
1	B	275	ASP	CB-CG-OD1	6.50	124.15	118.30
1	B	30	ASP	CB-CG-OD2	6.14	123.83	118.30
1	A	87	ASP	CB-CG-OD2	5.92	123.62	118.30
1	A	172	ASP	CB-CG-OD2	5.49	123.25	118.30
1	A	26	ASP	CB-CG-OD2	5.48	123.23	118.30
1	B	147	ASP	CB-CG-OD2	5.38	123.15	118.30
1	A	139	ASP	CB-CG-OD2	5.23	123.00	118.30
1	A	30	ASP	CB-CG-OD2	5.14	122.92	118.30
1	B	272	ASP	CB-CG-OD2	5.14	122.92	118.30

There are no chirality outliers.

There are no planarity outliers.

## 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1959	0	1894	22	0
1	B	2108	0	2043	13	0
2	A	28	0	25	0	0
2	B	28	0	25	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	1	0	0	0	0
4	B	4	0	0	0	0
5	B	2	0	0	0	0
6	A	18	0	0	1	0
6	B	35	0	0	3	0
All	All	4185	0	3987	35	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 4.

All (35) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:102:MET:HE2	6:B:2019:HOH:O	1.51	1.07
1:B:35:ARG:NH2	1:B:304:THR:O	2.13	0.81
1:A:226:GLU:CA	1:A:227:ASN:N	2.45	0.78
1:A:102:MET:CE	6:B:2019:HOH:O	2.19	0.77
1:A:146:GLU:N	1:A:146:GLU:CB	2.48	0.76
1:A:146:GLU:CD	1:A:146:GLU:CB	2.56	0.74
1:A:226:GLU:CB	1:A:226:GLU:CD	2.58	0.71
1:A:226:GLU:O	1:A:227:ASN:N	2.26	0.69
1:A:226:GLU:CA	1:A:226:GLU:O	2.43	0.66
1:B:44:LEU:HD21	1:B:112:VAL:HG21	1.77	0.66
1:B:35:ARG:HD2	1:B:300:GLN:HE21	1.62	0.65
1:B:85:SER:OG	6:B:2011:HOH:O	2.14	0.65
1:A:237:LYS:HE2	1:A:244:PRO:HD3	1.82	0.60
1:A:44:LEU:HD21	1:A:112:VAL:HG21	1.84	0.59
1:A:79:GLN:OE1	1:A:82:ARG:HD3	2.04	0.58
1:B:193:LEU:CD2	1:B:236:LEU:HD13	2.35	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:27:GLN:HG2	1:A:294:TRP:CE2	2.43	0.53
1:A:174:THR:HG21	6:A:2004:HOH:O	2.09	0.52
1:A:77:ALA:HB3	1:A:78:PRO:HD3	1.90	0.52
1:B:84:ALA:HB1	1:B:88:VAL:HG23	1.94	0.48
1:B:272:ASP:O	1:B:275:ASP:HB2	2.14	0.48
1:A:44:LEU:CD2	1:A:112:VAL:HG21	2.44	0.48
1:B:136:ASN:HB3	1:B:137:TYR:CD1	2.48	0.48
1:A:27:GLN:HG2	1:A:294:TRP:CD2	2.48	0.47
1:A:117:GLU:HG3	1:A:129:ILE:HD12	1.96	0.46
1:A:34:VAL:HA	1:A:128:PRO:O	2.17	0.44
1:B:44:LEU:CD2	1:B:112:VAL:HG21	2.47	0.43
1:A:145:SER:OG	1:A:147:ASP:OD1	2.37	0.43
1:B:271:GLU:HG3	1:B:271:GLU:O	2.18	0.43
1:B:22:LEU:HB2	1:B:23:PRO:CD	2.50	0.41
1:B:84:ALA:HB1	1:B:88:VAL:CG2	2.50	0.41
1:A:238:ILE:HA	1:A:238:ILE:HD12	1.92	0.41
1:A:291:PRO:HB2	1:A:294:TRP:CD1	2.55	0.41
1:B:35:ARG:HD2	1:B:300:GLN:NE2	2.33	0.41
1:A:51:PRO:CB	1:A:174:THR:HG22	2.51	0.40

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	245/286 (86%)	231 (94%)	14 (6%)	0	100	100
1	B	270/286 (94%)	253 (94%)	12 (4%)	5 (2%)	12	19
All	All	515/572 (90%)	484 (94%)	26 (5%)	5 (1%)	22	38

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	256	GLU

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Mol	Chain	Res	Type
1	B	24	VAL
1	B	275	ASP
1	B	23	PRO
1	B	92	PRO

### 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	210/247 (85%)	193 (92%)	17 (8%)	17	30
1	B	227/247 (92%)	208 (92%)	19 (8%)	16	28
All	All	437/494 (88%)	401 (92%)	36 (8%)	17	29

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	LEU
1	A	63	GLU
1	A	64	ARG
1	A	82	ARG
1	A	119	SER
1	A	123	SER
1	A	149	LEU
1	A	165	LEU
1	A	174	THR
1	A	195	LEU
1	A	196	ARG
1	A	231	VAL
1	A	238	ILE
1	A	296	ILE
1	A	298	SER
1	A	302	CYS
1	A	304	THR
1	B	22	LEU
1	B	34	VAL
1	B	35	ARG
1	B	53	LEU

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Mol	Chain	Res	Type
1	B	63	GLU
1	B	76	GLU
1	B	113	MET
1	B	149	LEU
1	B	165	LEU
1	B	190	LYS
1	B	210	GLN
1	B	231	VAL
1	B	236	LEU
1	B	256	GLU
1	B	271	GLU
1	B	278	LEU
1	B	280	GLU
1	B	283	VAL
1	B	302	CYS

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	39	HIS
1	A	105	ASN
1	A	132	GLN
1	B	209	GLN
1	B	295	HIS
1	B	300	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains 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
2	NAG	A	1307	1,2	12,14,15	0.63	0	15,19,21	1.33	1 (6%)
2	NAG	A	1308	2	12,14,15	0.90	0	15,19,21	1.50	2 (13%)
2	NAG	B	1308	1,2	12,14,15	0.70	1 (8%)	15,19,21	1.40	3 (20%)
2	NAG	B	1309	2	12,14,15	0.71	0	15,19,21	1.47	2 (13%)

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	NAG	A	1307	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	1308	2	-	0/6/23/26	0/1/1/1
2	NAG	B	1308	1,2	-	0/6/23/26	0/1/1/1
2	NAG	B	1309	2	-	0/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1308	NAG	O5-C5	-2.02	1.41	1.45

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1308	NAG	O5-C5-C4	4.23	116.02	110.65
2	A	1307	NAG	O5-C5-C6	4.12	111.31	106.98
2	B	1308	NAG	O4-C4-C3	-3.09	103.42	110.35
2	B	1309	NAG	O3-C3-C2	2.99	115.37	109.09
2	A	1308	NAG	C3-C4-C5	2.66	114.96	110.20
2	B	1308	NAG	C3-C2-N2	-2.64	107.75	111.76
2	B	1309	NAG	C4-C3-C2	-2.29	105.72	111.32
2	B	1308	NAG	O5-C5-C6	2.08	109.16	106.98

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.6 Ligand geometry ⓘ

Of 9 ligands modelled in this entry, 9 are monoatomic - 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.

## 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 ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

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	253/286 (88%)	0.34	17 (6%) 17 17	24, 44, 69, 75	0
1	B	274/286 (95%)	0.15	15 (5%) 24 24	21, 34, 67, 77	0
All	All	527/572 (92%)	0.24	32 (6%) 22 20	21, 38, 68, 77	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	22	LEU	7.4
1	A	191	TYR	6.4
1	A	255	PRO	5.3
1	B	24	VAL	5.2
1	B	270	PRO	5.1
1	A	306	TYR	5.1
1	A	24	VAL	4.9
1	B	257	LEU	4.7
1	B	305	PRO	4.4
1	B	85	SER	4.1
1	B	306	TYR	4.0
1	A	190	LYS	3.6
1	B	307	HIS	3.5
1	B	90	LYS	3.4
1	B	256	GLU	3.3
1	B	23	PRO	3.1
1	A	196	ARG	2.9
1	A	292	PRO	2.9
1	A	82	ARG	2.8
1	A	86	GLU	2.8
1	A	286	VAL	2.6
1	B	32	PRO	2.4
1	A	90	LYS	2.4
1	A	238	ILE	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	243	GLY	2.4
1	B	210	GLN	2.3
1	A	186	LYS	2.2
1	A	189	CYS	2.2
1	B	283	VAL	2.1
1	B	121	ASN	2.1
1	A	305	PRO	2.0
1	A	28	LEU	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates ⓘ

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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	NAG	A	1307	14/15	0.26	2.26	43,49,54,61	0
2	NAG	B	1308	14/15	0.22	1.27	44,52,57,64	0
2	NAG	B	1309	14/15	0.41	-	69,71,73,74	0
2	NAG	A	1308	14/15	0.43	-	65,67,70,70	0

## 6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
4	ZN	B	1313	1/1	0.15	2.44	53,53,53,53	0
5	NA	B	1312	1/1	0.13	-0.01	26,26,26,26	0
3	CL	A	1309	1/1	0.12	-0.83	42,42,42,42	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
4	ZN	B	1316	1/1	0.08	-1.52	60,60,60,60	0
3	CL	B	1310	1/1	0.06	-1.86	29,29,29,29	0
4	ZN	B	1314	1/1	0.06	-2.16	40,40,40,40	0
4	ZN	B	1315	1/1	0.06	-2.48	36,36,36,36	0
4	ZN	A	1310	1/1	0.12	-2.74	69,69,69,69	1
5	NA	B	1311	1/1	0.10	-5.50	44,44,44,44	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.