



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 21, 2017 – 05:48 AM EST

PDB ID : 1RNE  
Title : THE CRYSTAL STRUCTURE OF RECOMBINANT GLYCOSYLATED  
HUMAN RENIN ALONE AND IN COMPLEX WITH A TRANSITION  
STATE ANALOG INHIBITOR  
Authors : Gruetter, M.G.; Rahuel, J.; Priestle, J.P.  
Deposited on : unknown  
Resolution : 2.40 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) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20030345

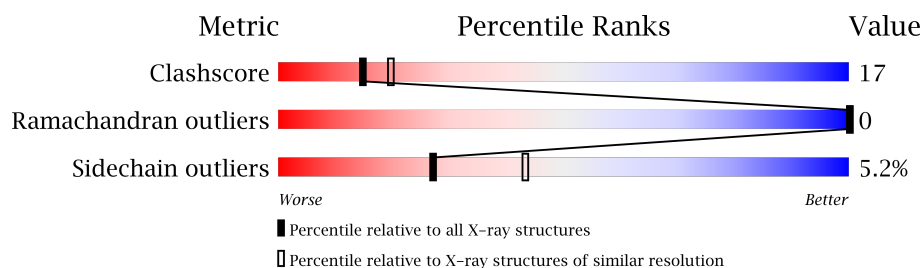
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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))
Clashscore	112137	3674 (2.40-2.40)
Ramachandran outliers	110173	3616 (2.40-2.40)
Sidechain outliers	110143	3617 (2.40-2.40)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	340	

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
2	NGA	A	345	X	-	-	-

## 2 Entry composition [i](#)

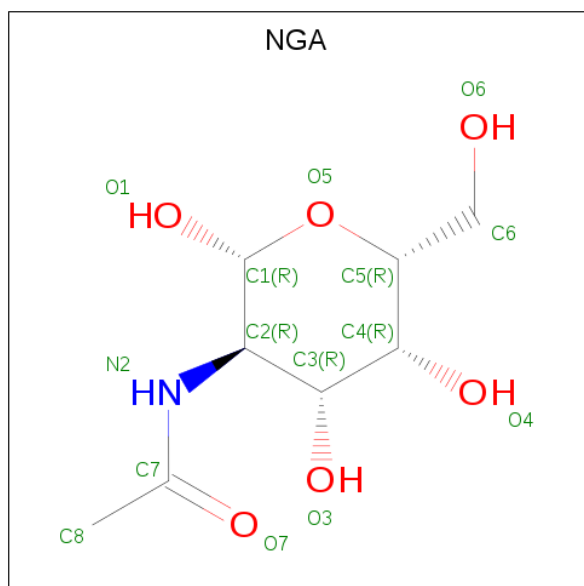
There are 4 unique types of molecules in this entry. The entry contains 2713 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 RENIN.

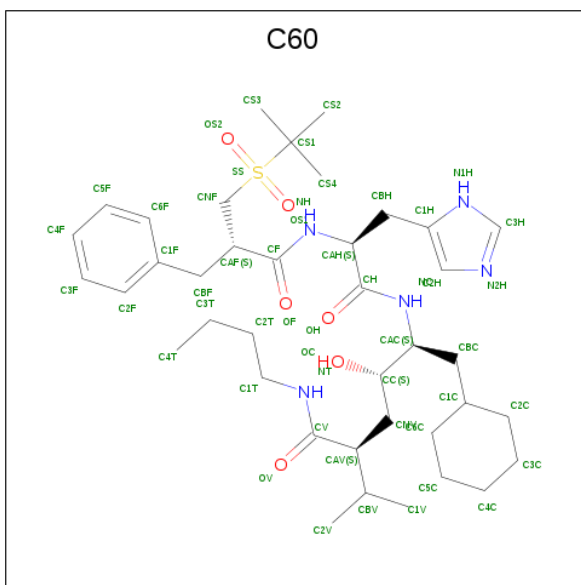
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	330	2510	1611	398	487	14	0	0	1

- Molecule 2 is N-ACETYL-D-GALACTOSAMINE (three-letter code: NGA) (formula:  $C_8H_{15}NO_6$ ).



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

- Molecule 3 is [[[3-(2-METHYL-PROPANE-2-SULFONYL)-1-BENZENYL]-2-PROPYL]-CARBONYL-HISTIDYL]-AMINO-[CYCLOHEXYLMETHYL]-[2-HYDROXY-4-ISOPROPYL]-PENTAN-5-OIC ACID BUTYLAMIDE (three-letter code: C60) (formula:  $C_{39}H_{63}N_5O_6S$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			51	39	5	6	1		

- Molecule 4 is water.

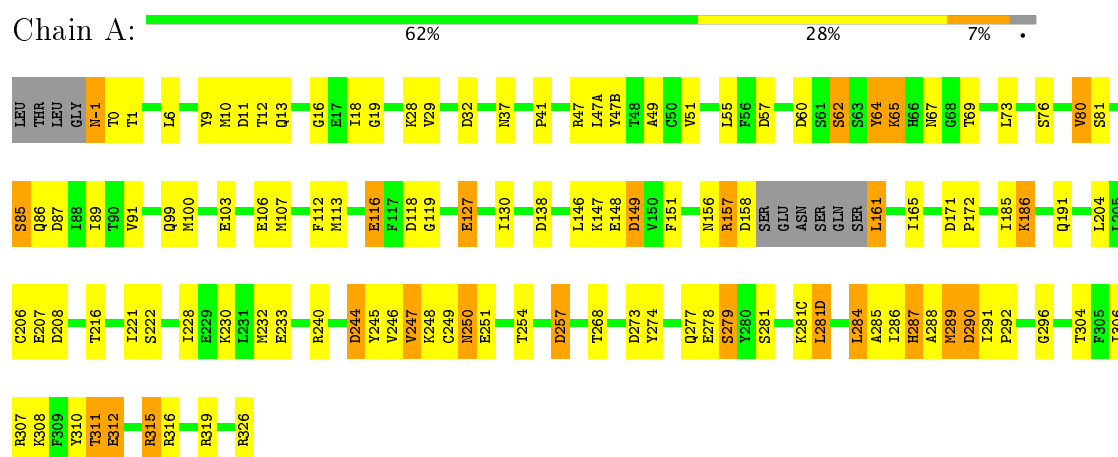
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	138	Total O 138 138	0	0

### 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 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.

Note EDS was not executed.

#### • Molecule 1: RENIN



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	90.90 Å 90.90 Å 109.40 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	(Not available) – 2.40	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-2.40)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	CORELS, TNT, X-PLOR	Depositor
R, $R_{free}$	0.176 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2713	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	37.0	wwPDB-VP

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NGA, C60

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	1.11	10/2569 (0.4%)	1.67	41/3491 (1.2%)

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	278	GLU	CD-OE1	7.61	1.34	1.25
1	A	106	GLU	CD-OE1	7.14	1.33	1.25
1	A	207	GLU	CD-OE1	6.35	1.32	1.25
1	A	116	GLU	CD-OE2	6.20	1.32	1.25
1	A	127	GLU	CD-OE2	5.89	1.32	1.25
1	A	103	GLU	CD-OE2	5.51	1.31	1.25
1	A	233	GLU	CD-OE1	5.25	1.31	1.25
1	A	312	GLU	CD-OE2	-5.24	1.19	1.25
1	A	251	GLU	CD-OE2	5.07	1.31	1.25
1	A	148	GLU	CD-OE1	5.02	1.31	1.25

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	274	TYR	CB-CG-CD2	-11.54	114.08	121.00
1	A	289	MET	CG-SD-CE	-10.59	83.26	100.20
1	A	307	ARG	NE-CZ-NH2	-10.13	115.24	120.30
1	A	118	ASP	CB-CG-OD2	-9.93	109.36	118.30
1	A	257	ASP	CB-CG-OD2	-9.36	109.88	118.30
1	A	149	ASP	CB-CG-OD1	-8.18	110.94	118.30
1	A	208	ASP	CB-CG-OD2	-8.09	111.02	118.30
1	A	157	ARG	NE-CZ-NH1	8.06	124.33	120.30
1	A	274	TYR	CB-CG-CD1	7.59	125.55	121.00
1	A	319	ARG	NE-CZ-NH2	-7.42	116.59	120.30
1	A	208	ASP	CB-CG-OD1	7.28	124.85	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	32	ASP	CB-CG-OD2	-7.15	111.86	118.30
1	A	118	ASP	CB-CG-OD1	7.14	124.72	118.30
1	A	310	TYR	CB-CG-CD1	-6.86	116.89	121.00
1	A	112	PHE	CB-CG-CD2	6.80	125.56	120.80
1	A	290	ASP	CB-CG-OD2	-6.80	112.18	118.30
1	A	85	SER	N-CA-CB	6.42	120.12	110.50
1	A	138	ASP	CB-CG-OD1	6.37	124.03	118.30
1	A	273	ASP	CB-CG-OD2	6.35	124.02	118.30
1	A	257	ASP	CB-CG-OD1	6.28	123.95	118.30
1	A	57	ASP	CB-CG-OD2	-6.26	112.66	118.30
1	A	149	ASP	CB-CG-OD2	6.25	123.93	118.30
1	A	60	ASP	CB-CG-OD2	-6.16	112.76	118.30
1	A	138	ASP	CB-CG-OD2	-6.14	112.77	118.30
1	A	273	ASP	CB-CG-OD1	-6.13	112.78	118.30
1	A	-1	ASN	CB-CA-C	5.89	122.18	110.40
1	A	315	ARG	NE-CZ-NH1	5.84	123.22	120.30
1	A	248	LYS	N-CA-CB	5.79	121.02	110.60
1	A	310	TYR	CB-CG-CD2	5.77	124.46	121.00
1	A	81	SER	CB-CA-C	-5.67	99.32	110.10
1	A	311	THR	CA-CB-CG2	-5.65	104.50	112.40
1	A	62	SER	CB-CA-C	-5.64	99.38	110.10
1	A	204	LEU	CB-CA-C	-5.62	99.52	110.20
1	A	244	ASP	CB-CG-OD1	5.53	123.27	118.30
1	A	112	PHE	CB-CG-CD1	-5.50	116.95	120.80
1	A	254	THR	CA-CB-CG2	-5.35	104.92	112.40
1	A	32	ASP	CB-CG-OD1	5.33	123.10	118.30
1	A	64	TYR	CB-CG-CD1	5.30	124.18	121.00
1	A	151	PHE	CB-CG-CD1	5.19	124.43	120.80
1	A	274	TYR	CG-CD1-CE1	-5.10	117.22	121.30
1	A	47(A)	LEU	C-N-CA	5.10	134.44	121.70

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	2510	0	2408	84	0
2	A	14	0	13	0	0
3	A	51	0	58	4	0
4	A	138	0	0	9	0
All	All	2713	0	2479	86	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 17.

All (86) 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:186:LYS:NZ	1:A:191:GLN:HE22	1.54	1.06
1:A:13:GLN:HE22	1:A:116:GLU:H	1.13	0.96
1:A:250:ASN:H	1:A:250:ASN:HD22	1.13	0.91
1:A:186:LYS:HZ3	1:A:191:GLN:NE2	1.69	0.90
1:A:49:ALA:HA	1:A:113:MET:CE	2.11	0.81
1:A:186:LYS:HZ3	1:A:191:GLN:HE22	0.84	0.81
1:A:246:VAL:HG11	1:A:281(D):LEU:HG	1.66	0.78
1:A:13:GLN:HE22	1:A:116:GLU:N	1.82	0.77
1:A:13:GLN:NE2	1:A:116:GLU:H	1.83	0.77
1:A:186:LYS:NZ	1:A:191:GLN:NE2	2.29	0.76
1:A:49:ALA:HA	1:A:113:MET:HE2	1.66	0.75
1:A:250:ASN:H	1:A:250:ASN:ND2	1.85	0.73
1:A:291:ILE:O	1:A:296:GLY:HA3	1.90	0.71
1:A:67:ASN:OD1	1:A:69:THR:HG23	1.89	0.71
1:A:257:ASP:HB3	4:A:413:HOH:O	1.91	0.70
1:A:76:SER:HB2	3:A:350:C60:OH	1.91	0.70
1:A:268:THR:O	1:A:308:LYS:HE2	1.91	0.69
1:A:284:LEU:N	1:A:284:LEU:HD23	2.07	0.69
1:A:29:VAL:HA	1:A:119:GLY:O	1.93	0.69
1:A:249:CYS:HB3	1:A:279:SER:O	1.94	0.67
1:A:158:ASP:N	4:A:414:HOH:O	2.29	0.66
1:A:246:VAL:CG1	1:A:281(D):LEU:HG	2.28	0.64
1:A:240:ARG:HB2	1:A:244:ASP:OD1	2.00	0.62
1:A:41:PRO:HB2	1:A:55:LEU:HD23	1.80	0.62
1:A:85:SER:HB3	4:A:408:HOH:O	1.98	0.61
1:A:49:ALA:HA	1:A:113:MET:HE3	1.83	0.60
1:A:41:PRO:HB3	1:A:107:MET:CE	2.32	0.59
1:A:47:ARG:C	1:A:47(B):TYR:H	2.07	0.58
1:A:86:GLN:HG2	1:A:87:ASP:N	2.18	0.58
1:A:10:MET:O	1:A:12:THR:HG23	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:41:PRO:HB3	1:A:107:MET:HE1	1.85	0.58
1:A:37:ASN:HB3	4:A:386:HOH:O	2.04	0.57
1:A:41:PRO:HB2	1:A:55:LEU:CD2	2.35	0.57
1:A:232:MET:HG3	1:A:245:TYR:CE2	2.40	0.56
1:A:288:ALA:O	1:A:289:MET:HG2	2.04	0.56
1:A:64:TYR:HE1	1:A:85:SER:HB2	1.68	0.56
1:A:64:TYR:CE1	1:A:85:SER:HB2	2.42	0.55
1:A:91:VAL:O	1:A:91:VAL:HG13	2.07	0.54
1:A:232:MET:HG3	1:A:245:TYR:CD2	2.41	0.54
1:A:216:THR:HG22	1:A:306:ILE:HD13	1.89	0.54
1:A:250:ASN:N	1:A:250:ASN:HD22	1.92	0.53
1:A:316:ARG:HB3	4:A:452:HOH:O	2.08	0.52
1:A:9:TYR:HB3	1:A:13:GLN:HB2	1.92	0.52
1:A:185:ILE:HG22	1:A:191:GLN:O	2.09	0.52
1:A:221:ILE:HG13	1:A:304:THR:HB	1.92	0.52
1:A:13:GLN:NE2	1:A:116:GLU:HB2	2.26	0.50
1:A:249:CYS:HB2	1:A:281:SER:O	2.11	0.50
1:A:-1:ASN:HB3	1:A:147:LYS:HA	1.94	0.50
3:A:350:C60:C1V	3:A:350:C60:NT	2.75	0.49
1:A:206:CYS:SG	1:A:206:CYS:O	2.70	0.49
1:A:250:ASN:ND2	1:A:250:ASN:N	2.51	0.49
1:A:149:ASP:O	1:A:315:ARG:HB2	2.13	0.49
1:A:10:MET:HE2	4:A:478:HOH:O	2.12	0.49
1:A:6:LEU:HD11	1:A:165:ILE:HG13	1.96	0.48
1:A:285:ALA:C	1:A:286:ILE:HG13	2.35	0.47
1:A:291:ILE:HD11	3:A:350:C60:H3H	1.96	0.47
1:A:51:VAL:HG12	1:A:51:VAL:O	2.15	0.47
1:A:277:GLN:NE2	1:A:279:SER:O	2.46	0.47
1:A:16:GLY:O	1:A:28:LYS:HA	2.15	0.47
1:A:246:VAL:CG1	1:A:247:VAL:N	2.78	0.47
1:A:86:GLN:HE21	1:A:100:MET:HE1	1.81	0.46
1:A:127:GLU:OE1	1:A:127:GLU:N	2.45	0.46
1:A:89:ILE:HD12	1:A:99:GLN:HB3	1.97	0.46
1:A:287:HIS:HB2	4:A:419:HOH:O	2.16	0.46
1:A:161:LEU:HA	1:A:161:LEU:HD22	1.80	0.46
1:A:65:LYS:HD3	1:A:65:LYS:C	2.36	0.46
1:A:156:ASN:OD1	1:A:157:ARG:N	2.42	0.45
1:A:0:THR:HG22	1:A:1:THR:N	2.32	0.45
1:A:0:THR:O	1:A:146:LEU:HA	2.17	0.44
1:A:171:ASP:HA	1:A:172:PRO:HD2	1.77	0.44
1:A:64:TYR:HE1	1:A:85:SER:CB	2.30	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:18:ILE:HG22	1:A:91:VAL:HB	2.01	0.43
1:A:149:ASP:N	1:A:149:ASP:OD1	2.52	0.43
1:A:244:ASP:N	1:A:244:ASP:OD1	2.51	0.43
1:A:311:THR:HG22	1:A:312:GLU:N	2.34	0.43
3:A:350:C60:H6C1	3:A:350:C60:HAC	1.69	0.42
1:A:228:ILE:O	1:A:232:MET:HG2	2.20	0.42
1:A:73:LEU:HB2	1:A:80:VAL:HG23	2.01	0.42
1:A:291:ILE:HA	1:A:292:PRO:HD3	1.87	0.42
1:A:62:SER:CB	4:A:444:HOH:O	2.67	0.41
1:A:249:CYS:SG	1:A:281(C):LYS:O	2.79	0.41
1:A:10:MET:O	1:A:11:ASP:HB2	2.18	0.41
1:A:107:MET:HB2	1:A:107:MET:HE3	1.62	0.41
1:A:290:ASP:HA	4:A:406:HOH:O	2.21	0.41
1:A:19:GLY:O	1:A:89:ILE:HA	2.22	0.40
1:A:47:ARG:CG	1:A:51:VAL:HG22	2.51	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	326/340 (96%)	317 (97%)	9 (3%)	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	270/290 (93%)	256 (95%)	14 (5%)	27	43

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	65	LYS
1	A	80	VAL
1	A	130	ILE
1	A	161	LEU
1	A	186	LYS
1	A	222	SER
1	A	230	LYS
1	A	247	VAL
1	A	250	ASN
1	A	279	SER
1	A	281(D)	LEU
1	A	284	LEU
1	A	287	HIS
1	A	326	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	13	GLN
1	A	86	GLN
1	A	173	GLN
1	A	174	HIS
1	A	191	GLN
1	A	250	ASN
1	A	317	ASN

### 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 [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

2 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
2	NGA	A	345	1	14,14,15	2.48	1 (7%)	15,19,21	2.11	2 (13%)
3	C60	A	350	-	48,53,53	1.94	5 (10%)	59,73,73	2.13	16 (27%)

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	NGA	A	345	1	1/1/5/7	0/6/23/26	0/1/1/1
3	C60	A	350	-	-	0/63/71/71	0/3/3/3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	350	C60	CS1-SS	-10.33	1.71	1.82
3	A	350	C60	OS2-SS	-3.90	1.38	1.44
3	A	350	C60	CAF-CF	-2.80	1.46	1.51
3	A	350	C60	CAH-NH	-2.36	1.40	1.45
3	A	350	C60	CBF-CAF	-2.24	1.47	1.53
2	A	345	NGA	C4-C5	8.84	1.71	1.53

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	350	C60	OS2-SS-OS1	-7.13	107.80	118.06

*Continued on next page...*

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	345	NGA	C1-O5-C5	-6.50	103.20	112.17
2	A	345	NGA	C2-N2-C7	-4.16	116.87	122.94
3	A	350	C60	CBC-CAC-CC	-3.91	106.63	112.55
3	A	350	C60	OF-CF-CAF	-3.24	117.96	122.11
3	A	350	C60	CBC-C1C-C6C	-3.15	104.41	111.70
3	A	350	C60	C3C-C2C-C1C	-3.06	106.58	112.19
3	A	350	C60	CBF-C1F-C2F	-2.50	115.85	120.91
3	A	350	C60	C2V-CBV-C1V	-2.49	103.53	110.64
3	A	350	C60	CH-CAH-NH	-2.41	104.56	111.20
3	A	350	C60	CAV-CNV-CC	-2.41	109.73	113.92
3	A	350	C60	C1V-CBV-CAV	-2.22	107.83	111.94
3	A	350	C60	CAH-CH-NC	-2.14	111.95	116.78
3	A	350	C60	OF-CF-NH	2.43	127.48	122.90
3	A	350	C60	OS2-SS-CS1	3.89	111.73	107.73
3	A	350	C60	C1T-NT-CV	3.92	129.90	122.59
3	A	350	C60	CC-CAC-NC	4.16	116.92	109.52
3	A	350	C60	OS1-SS-CNF	5.13	113.01	108.26

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	345	NGA	C4

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
3	A	350	C60	4	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

### 6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

### 6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

### 6.4 Ligands

EDS was not executed - this section is therefore empty.

### 6.5 Other polymers

EDS was not executed - this section is therefore empty.