



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 29, 2018 – 02:02 PM EST

PDB ID : 5LWY  
Title : Revised crystal structure of the human adiponectin receptor 2 in complex with a C18 free fatty acid  
Authors : Leyrat, C.; Vasiliauskaite-Brooks, I.; Granier, S.  
Deposited on : 2016-09-19  
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) : 1.9-1692  
EDS : rb-20030736  
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) : rb-20030736

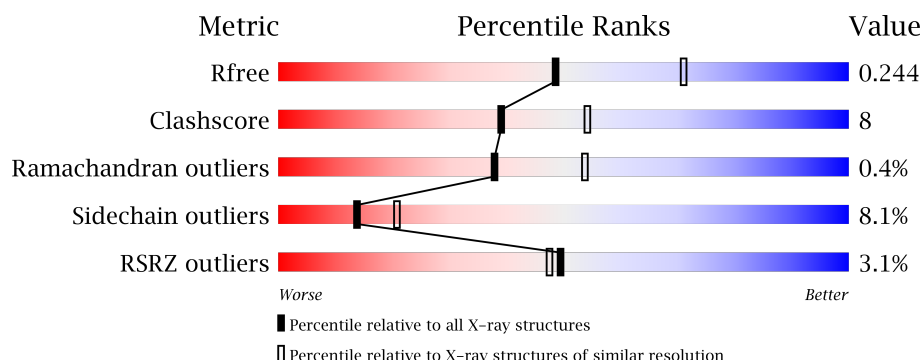
# 1 Overall quality at a glance

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))
$R_{free}$	100719	3166 (2.40-2.40)
Clashscore	112137	3674 (2.40-2.40)
Ramachandran outliers	110173	3616 (2.40-2.40)
Sidechain outliers	110143	3617 (2.40-2.40)
RSRZ outliers	101464	3195 (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.

Mol	Chain	Length	Quality of chain
1	A	292	<div> <div>4%</div> <div> <div></div> <div>76%</div> <div>18%</div> <div>• •</div> </div> </div>
2	H	119	<div> <div>3%</div> <div> <div></div> <div>82%</div> <div>16%</div> <div>•</div> </div> </div>
3	L	107	<div> <div></div> <div> <div></div> <div>84%</div> <div>16%</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
5	OLB	A	402	-	-	-	X
5	OLB	A	403	-	-	-	X
5	OLB	A	404	-	-	-	X
5	OLB	A	405	-	-	-	X
5	OLB	A	407	-	-	-	X
5	OLB	A	409	-	-	-	X
6	OLA	A	410	-	-	-	X
7	GOL	L	201	-	-	-	X

## 2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 4613 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 Adiponectin receptor protein 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	285	2325	1563	377	368	17	0	2	0

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	GLY	-	expression tag	UNP Q86V24
A	-3	GLY	-	expression tag	UNP Q86V24
A	-2	SER	-	expression tag	UNP Q86V24
A	-1	GLU	-	expression tag	UNP Q86V24
A	0	PHE	-	expression tag	UNP Q86V24

- Molecule 2 is a protein called V REGION HEAVY CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	H	119	927	592	149	182	4	0	0	0

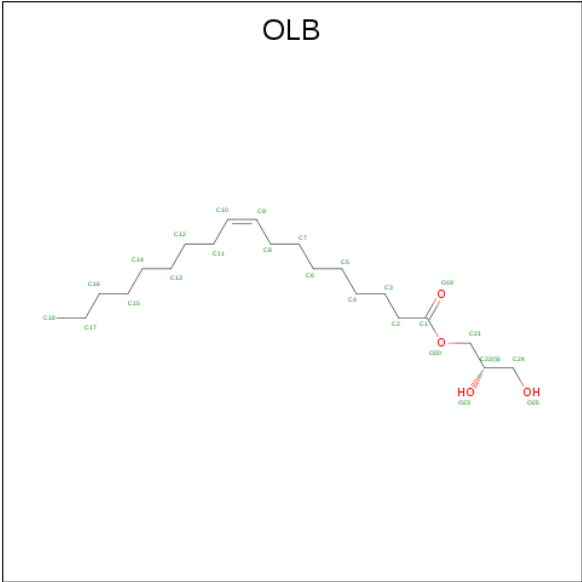
- Molecule 3 is a protein called V REGION LIGHT CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	L	107	820	517	137	163	3	0	0	0

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

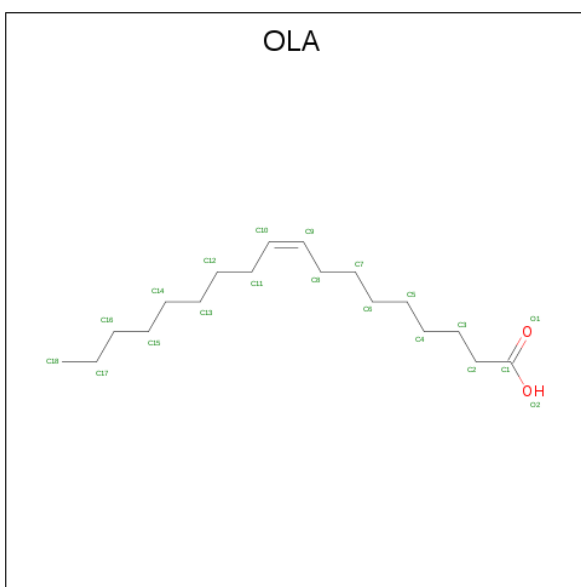
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Zn	0	0
			1	1		

- Molecule 5 is (2S)-2,3-dihydroxypropyl (9Z)-octadec-9-enoate (three-letter code: OLB) (formula: C<sub>21</sub>H<sub>40</sub>O<sub>4</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			25	21	4		
5	A	1	Total	C	O	0	0
			25	21	4		
5	A	1	Total	C	O	0	0
			25	21	4		
5	A	1	Total	C	O	6	0
			25	21	4		
5	A	1	Total	C	O	0	0
			25	21	4		
5	A	1	Total	C	O	0	0
			25	21	4		
5	A	1	Total	C	O	0	0
			25	21	4		

- Molecule 6 is OLEIC ACID (three-letter code: OLA) (formula: C<sub>18</sub>H<sub>34</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			20	18	2		

- Molecule 7 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	L	1	Total	C	O	0	0
			6	3	3		

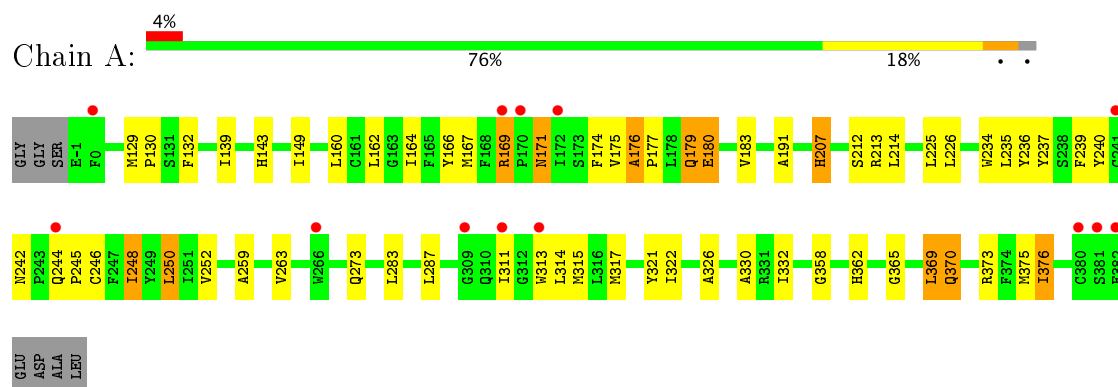
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	153	Total 153	O 153	0	0
8	H	89	Total 89	O 89	0	0
8	L	72	Total 72	O 72	0	0

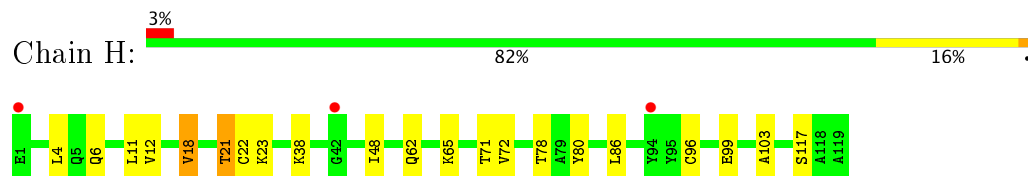
### 3 Residue-property plots [i](#)

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.

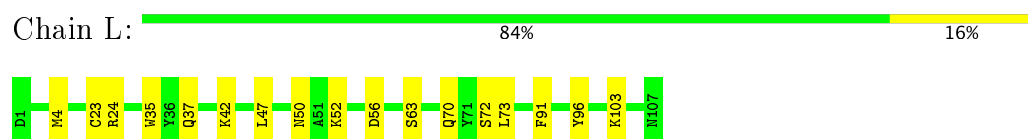
- Molecule 1: Adiponectin receptor protein 2



- Molecule 2: V REGION HEAVY CHAIN



- Molecule 3: V REGION LIGHT CHAIN





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 2 <sub>1</sub> 2 <sub>1</sub> 2 <sub>1</sub>	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	74.58Å 101.03Å 108.63Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.52 – 2.40 19.52 – 2.40	Depositor EDS
% Data completeness (in resolution range)	98.3 (19.52-2.40) 98.3 (19.52-2.40)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.57 (at 2.41Å)	Xtriage
Refinement program	BUSTER 2.10.2	Depositor
R, $R_{free}$	0.198 , 0.224 0.216 , 0.244	Depositor DCC
$R_{free}$ test set	1607 reflections (5.00%)	DCC
Wilson B-factor (Å <sup>2</sup> )	47.2	Xtriage
Anisotropy	0.653	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 75.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	4613	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 5.63% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, ZN, OLB, OLA

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	0.44	0/2413	0.64	0/3282
2	H	0.36	0/951	0.64	0/1291
3	L	0.37	0/840	0.59	0/1140
All	All	0.41	0/4204	0.63	0/5713

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

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	2325	0	2289	56	0
2	H	927	0	888	10	0
3	L	820	0	786	6	0
4	A	1	0	0	0	0
5	A	200	0	320	8	0
6	A	20	0	33	2	0
7	L	6	0	8	0	0
8	A	153	0	0	4	0
8	H	89	0	0	0	0
8	L	72	0	0	0	0
All	All	4613	0	4324	72	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 8.

All (72) 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:183:VAL:HG13	8:A:501:HOH:O	1.45	1.15
1:A:179:GLN:NE2	8:A:501:HOH:O	1.84	1.09
1:A:171:ASN:HB3	1:A:177:PRO:HG2	1.29	1.07
1:A:171:ASN:HB3	1:A:177:PRO:CG	1.90	1.00
1:A:237:TYR:CE2	1:A:370:GLN:HA	2.08	0.88
1:A:174:PHE:HB2	1:A:177:PRO:HG3	1.59	0.85
1:A:176:ALA:O	1:A:180:GLU:HB2	1.76	0.84
1:A:166:TYR:CE1	1:A:169:ARG:NH2	2.51	0.79
1:A:237:TYR:HE2	1:A:370:GLN:HA	1.47	0.77
1:A:245:PRO:HA	1:A:248:ILE:HD11	1.69	0.75
1:A:321:TYR:HA	6:A:410:OLA:H111	1.73	0.70
1:A:183:VAL:CG1	8:A:501:HOH:O	2.20	0.70
1:A:245:PRO:HA	1:A:248:ILE:CD1	2.23	0.69
1:A:322:ILE:HD12	5:A:408:OLB:H5	1.77	0.67
2:H:12:VAL:HG11	2:H:18:VAL:HG13	1.83	0.61
1:A:234[B]:TRP:CZ3	1:A:369:LEU:HD11	2.37	0.60
1:A:143:HIS:HB2	8:A:505:HOH:O	2.04	0.57
1:A:171:ASN:HD22	1:A:177:PRO:HB2	1.70	0.57
2:H:22:CYS:HG	2:H:96:CYS:HG	1.39	0.56
5:A:404:OLB:O23	5:A:409:OLB:H19	2.05	0.56
1:A:287:LEU:HD12	5:A:408:OLB:H21	1.89	0.55
1:A:330:ALA:HB3	1:A:332:ILE:HD12	1.88	0.55
1:A:175:VAL:O	1:A:176:ALA:HB2	2.06	0.54
2:H:12:VAL:HG21	2:H:86:LEU:HD13	1.89	0.54
1:A:160:LEU:O	1:A:164:ILE:HG12	2.08	0.53
1:A:373:ARG:NH2	1:A:376:ILE:HD11	2.24	0.53
1:A:166:TYR:CD1	1:A:169:ARG:NH2	2.77	0.53
1:A:248:ILE:O	1:A:252:VAL:HG23	2.09	0.53
1:A:283:LEU:HD11	1:A:322:ILE:HD13	1.89	0.52
1:A:171:ASN:ND2	1:A:177:PRO:CB	2.72	0.52
1:A:175:VAL:O	1:A:175:VAL:HG12	2.09	0.52
1:A:225:LEU:HD11	5:A:405:OLB:H241	1.91	0.52
1:A:171:ASN:ND2	1:A:177:PRO:HB2	2.24	0.51
2:H:6:GLN:HA	2:H:21:THR:O	2.11	0.51
1:A:139:ILE:HG12	1:A:149:ILE:HG21	1.93	0.49
2:H:18:VAL:HG22	2:H:86:LEU:HD11	1.93	0.49
1:A:259:ALA:O	1:A:263:VAL:HG23	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:38:LYS:HB2	2:H:48:ILE:HD11	1.95	0.48
1:A:171:ASN:OD1	1:A:171:ASN:N	2.47	0.48
1:A:365:GLY:O	1:A:369:LEU:HD12	2.14	0.47
3:L:4:MET:CE	3:L:23:CYS:SG	3.02	0.47
1:A:240:TYR:CE2	1:A:373:ARG:HD2	2.49	0.47
2:H:71:THR:OG1	2:H:80:TYR:HB2	2.14	0.46
1:A:166:TYR:HE1	1:A:169:ARG:HH22	1.62	0.46
1:A:174:PHE:CD1	1:A:174:PHE:N	2.84	0.46
1:A:236:TYR:HD2	1:A:237:TYR:CE1	2.33	0.46
1:A:166:TYR:HE1	1:A:169:ARG:NH2	2.05	0.46
3:L:35:TRP:CE2	3:L:73:LEU:HB2	2.50	0.46
2:H:12:VAL:HG11	2:H:18:VAL:CG1	2.44	0.45
3:L:37:GLN:HB2	3:L:47:LEU:HD11	2.00	0.44
1:A:130:PRO:HG3	3:L:50:ASN:HB3	1.99	0.44
1:A:237:TYR:CE2	1:A:370:GLN:CA	2.93	0.44
1:A:358:GLY:HA3	6:A:410:OLA:H162	1.98	0.44
1:A:287:LEU:CD1	5:A:408:OLB:H21	2.47	0.44
2:H:23:LYS:HA	2:H:78:THR:HG22	1.99	0.43
2:H:99:GLU:HG2	2:H:103:ALA:HA	2.00	0.43
1:A:129:MET:O	1:A:207:HIS:HA	2.18	0.43
1:A:313[B]:TRP:CZ2	1:A:362:HIS:HA	2.54	0.43
1:A:317:MET:HE2	5:A:409:OLB:C18	2.48	0.43
1:A:239:PHE:CB	1:A:246:CYS:HB2	2.49	0.43
1:A:313[B]:TRP:CE3	1:A:314:LEU:HG	2.54	0.43
1:A:326:ALA:HA	5:A:408:OLB:H20	2.01	0.43
3:L:4:MET:HE3	3:L:23:CYS:SG	2.59	0.43
1:A:171:ASN:HD22	1:A:177:PRO:CB	2.32	0.42
1:A:132:PHE:HB2	5:A:402:OLB:H4A	2.01	0.42
1:A:242:ASN:O	1:A:245:PRO:HD2	2.21	0.41
1:A:235:LEU:HD12	1:A:250:LEU:HD12	2.02	0.41
1:A:171:ASN:HB3	1:A:177:PRO:CB	2.48	0.41
1:A:175:VAL:O	1:A:175:VAL:CG1	2.69	0.41
1:A:248:ILE:HG13	1:A:248:ILE:H	1.66	0.41
3:L:91:PHE:HA	3:L:96:TYR:CD1	2.55	0.40
1:A:191:ALA:HA	1:A:226:LEU:HD13	2.03	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	285/292 (98%)	275 (96%)	8 (3%)	2 (1%)	25	37
2	H	117/119 (98%)	115 (98%)	2 (2%)	0	100	100
3	L	105/107 (98%)	102 (97%)	3 (3%)	0	100	100
All	All	507/518 (98%)	492 (97%)	13 (3%)	2 (0%)	38	54

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	176	ALA
1	A	207	HIS

### 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	244/246 (99%)	225 (92%)	19 (8%)	15	23
2	H	99/99 (100%)	91 (92%)	8 (8%)	14	21
3	L	90/90 (100%)	82 (91%)	8 (9%)	11	17
All	All	433/435 (100%)	398 (92%)	35 (8%)	14	21

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	162	LEU

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Mol	Chain	Res	Type
1	A	167	MET
1	A	169	ARG
1	A	171	ASN
1	A	179	GLN
1	A	180	GLU
1	A	212	SER
1	A	213	ARG
1	A	214	LEU
1	A	244	GLN
1	A	248	ILE
1	A	250	LEU
1	A	273	GLN
1	A	311	ILE
1	A	315	MET
1	A	369	LEU
1	A	370	GLN
1	A	375	MET
1	A	376	ILE
2	H	4	LEU
2	H	11	LEU
2	H	18	VAL
2	H	21	THR
2	H	62	GLN
2	H	65	LYS
2	H	72	VAL
2	H	117	SER
3	L	24	ARG
3	L	42	LYS
3	L	52	LYS
3	L	56	ASP
3	L	63	SER
3	L	70	GLN
3	L	72	SER
3	L	103	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	244	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](#)

Of 11 ligands modelled in this entry, 1 is monoatomic - leaving 10 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	$\# Z  > 2$
5	OLB	A	402	-	24,24,24	0.19	0	25,25,25	0.24	0
5	OLB	A	403	-	24,24,24	0.21	0	25,25,25	0.32	0
5	OLB	A	404	-	24,24,24	0.24	0	25,25,25	0.48	0
5	OLB	A	405	-	24,24,24	0.20	0	25,25,25	0.30	0
5	OLB	A	406	-	24,24,24	0.29	0	25,25,25	0.27	0
5	OLB	A	407	-	24,24,24	0.24	0	25,25,25	0.25	0
5	OLB	A	408	-	24,24,24	0.28	0	25,25,25	0.45	0
5	OLB	A	409	-	24,24,24	0.17	0	25,25,25	0.41	0
6	OLA	A	410	-	16,19,19	0.19	0	15,19,19	0.38	0
7	GOL	L	201	-	5,5,5	0.14	0	5,5,5	0.33	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
5	OLB	A	402	-	-	0/24/24/24	0/0/0/0
5	OLB	A	403	-	-	0/24/24/24	0/0/0/0
5	OLB	A	404	-	-	0/24/24/24	0/0/0/0
5	OLB	A	405	-	-	0/24/24/24	0/0/0/0
5	OLB	A	406	-	-	0/24/24/24	0/0/0/0
5	OLB	A	407	-	-	0/24/24/24	0/0/0/0
5	OLB	A	408	-	-	0/24/24/24	0/0/0/0
5	OLB	A	409	-	-	0/24/24/24	0/0/0/0
6	OLA	A	410	-	-	0/15/17/17	0/0/0/0
7	GOL	L	201	-	-	0/4/4/4	0/0/0/0

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.

6 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	402	OLB	1	0
5	A	404	OLB	1	0
5	A	405	OLB	1	0
5	A	408	OLB	4	0
5	A	409	OLB	2	0
6	A	410	OLA	2	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	285/292 (97%)	0.07	13 (4%) 33 31	53, 81, 151, 176	2 (0%)
2	H	119/119 (100%)	-0.16	3 (2%) 58 55	56, 70, 86, 100	0
3	L	107/107 (100%)	-0.29	0 100 100	50, 65, 79, 92	0
All	All	511/518 (98%)	-0.06	16 (3%) 49 47	50, 73, 138, 176	2 (0%)

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	172	ILE	6.2
1	A	170	PRO	3.9
1	A	311	ILE	3.9
1	A	382	GLU	3.2
1	A	381	SER	3.1
2	H	42	GLY	3.1
1	A	380	CYS	3.0
1	A	169	ARG	2.9
1	A	266	TRP	2.7
1	A	244	GLN	2.6
1	A	309	GLY	2.6
2	H	94	TYR	2.5
1	A	241	CYS	2.5
1	A	313[A]	TRP	2.4
1	A	0	PHE	2.2
2	H	1	GLU	2.1

### 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, 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
5	OLB	A	407	25/25	0.77	0.49	10.10	88,94,104,105	0
5	OLB	A	405	25/25	0.72	0.46	6.67	83,86,99,103	6
5	OLB	A	403	25/25	0.52	0.29	5.19	81,94,116,121	0
5	OLB	A	404	25/25	0.61	0.30	4.66	120,131,140,142	0
6	OLA	A	410	20/20	0.85	0.25	3.54	72,76,78,78	0
5	OLB	A	409	25/25	0.57	0.40	3.39	85,111,146,152	0
5	OLB	A	402	25/25	0.78	0.23	2.31	85,88,115,124	0
7	GOL	L	201	6/6	0.83	0.22	2.15	76,78,81,82	0
5	OLB	A	408	25/25	0.68	0.31	1.52	100,106,112,114	0
4	ZN	A	401	1/1	0.99	0.12	-	61,61,61,61	0
5	OLB	A	406	25/25	0.57	0.46	-	85,93,118,122	0

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

There are no such residues in this entry.