



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 17, 2017 – 02:42 AM EDT

PDB ID : 3E7C  
Title : Glucocorticoid Receptor LBD bound to GSK866  
Authors : Madauss, K.P.; Williams, S.P.; Mclay, I.; Stewart, E.L.; Bledsoe, R.K.  
Deposited on : unknown  
Resolution : 2.15 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-20030345  
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-20030345

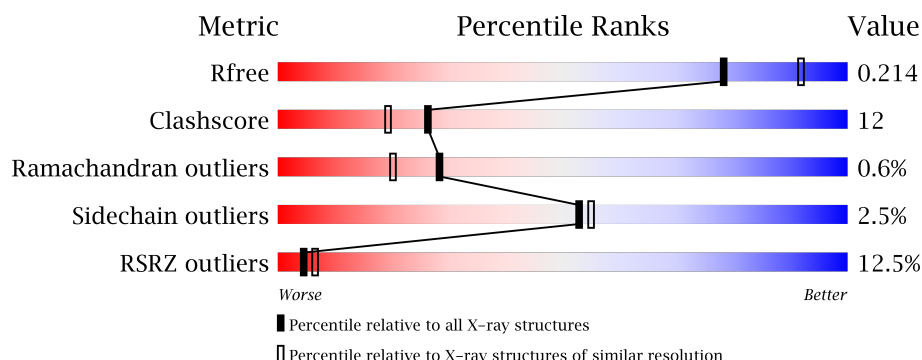
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.15 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	1170 (2.16-2.16)
Clashscore	112137	1278 (2.16-2.16)
Ramachandran outliers	110173	1256 (2.16-2.16)
Sidechain outliers	110143	1255 (2.16-2.16)
RSRZ outliers	101464	1175 (2.16-2.16)

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	257	<div> <div>10%</div> <div> <div></div> <div>77%</div> <div>18%</div> <div>• •</div> </div> </div>
1	B	257	<div> <div>15%</div> <div> <div></div> <div>74%</div> <div>21%</div> <div>• 5%</div> </div> </div>
2	D	11	<div> <div></div> <div> <div>27%</div> <div>55%</div> <div>9%</div> <div>9%</div> </div> </div>
2	H	11	<div> <div></div> <div> <div>82%</div> <div>18%</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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	GOL	A	778	-	-	-	X

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4330 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 Glucocorticoid receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	248	Total	C	N	O	S	0	0	0
			2003	1300	328	358	17			
1	B	245	Total	C	N	O	S	0	0	0
			1965	1282	317	349	17			

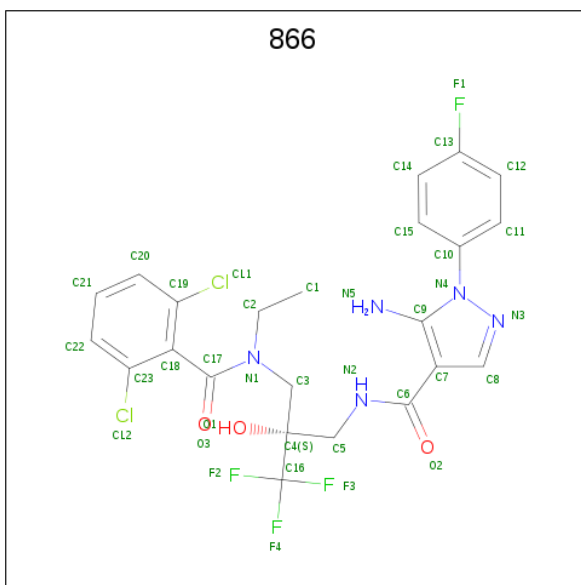
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	602	TYR	PHE	ENGINEERED	UNP P04150
A	638	GLY	CYS	ENGINEERED	UNP P04150
B	602	TYR	PHE	ENGINEERED	UNP P04150
B	638	GLY	CYS	ENGINEERED	UNP P04150

- Molecule 2 is a protein called Nuclear receptor coactivator 2.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	H	11	Total	C	N	O	0	0	0
			91	58	15	18			
2	D	10	Total	C	N	O	0	0	0
			86	56	15	15			

- Molecule 3 is 5-amino-N-[(2S)-2-({[(2,6-dichlorophenyl)carbonyl](ethyl)amino}methyl)-3,3,3-trifluoro-2-hydroxypropyl]-1-(4-fluorophenyl)-1H-pyrazole-4-carboxamide (three-letter code: 866) (formula: C<sub>23</sub>H<sub>21</sub>Cl<sub>2</sub>F<sub>4</sub>N<sub>5</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	Cl	F	N	O	0	0
			37	23	2	4	5	3		
3	B	1	Total	C	Cl	F	N	O	0	0
			37	23	2	4	5	3		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			6	3	3		

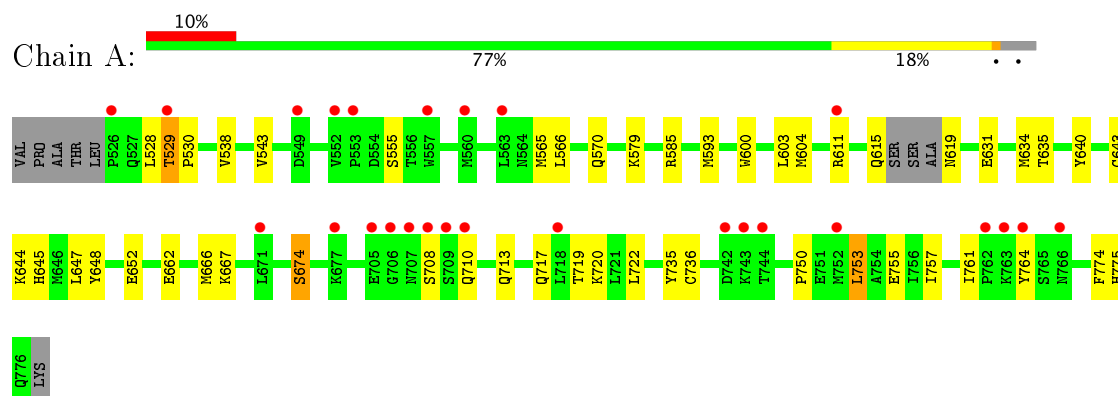
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	39	Total	O	0	0
			39	39		
5	H	2	Total	O	0	0
			2	2		
5	B	52	Total	O	0	0
			52	52		

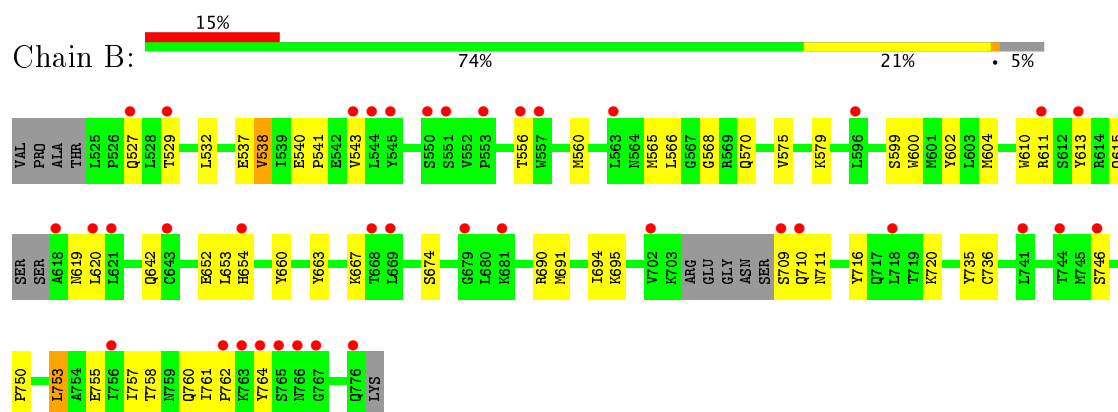
### 3 Residue-property plots [i](#)

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.

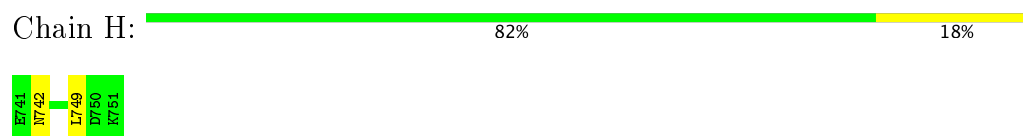
#### • Molecule 1: Glucocorticoid receptor



#### • Molecule 1: Glucocorticoid receptor



#### • Molecule 2: Nuclear receptor coactivator 2



#### • Molecule 2: Nuclear receptor coactivator 2



GLU
N742
N743
L744
L745
N746
N747
L748
L749
D750
K751



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	126.65Å 126.65Å 78.99Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.91 – 2.15 19.91 – 2.15	Depositor EDS
% Data completeness (in resolution range)	85.5 (19.91-2.15) 96.2 (19.91-2.15)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.34 (at 2.15Å)	Xtriage
Refinement program	PHENIX	Depositor
R, $R_{free}$	0.210 , 0.266 0.210 , 0.214	Depositor DCC
$R_{free}$ test set	2674 reflections (7.09%)	DCC
Wilson B-factor (Å <sup>2</sup> )	44.7	Xtriage
Anisotropy	0.739	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 58.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.032 for h,-h-k,-l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	4330	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	77.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.13% 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: GOL, 866

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	0.36	0/2048	0.50	0/2773
1	B	0.35	0/2009	0.51	0/2723
2	D	0.38	0/86	0.51	0/113
2	H	0.37	0/91	0.51	0/121
All	All	0.36	0/4234	0.50	0/5730

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	2003	0	1999	42	0
1	B	1965	0	1956	53	0
2	D	86	0	93	7	0
2	H	91	0	88	4	0
3	A	37	0	21	5	0
3	B	37	0	21	7	0
4	A	6	0	8	2	0
4	B	12	0	16	3	0
5	A	39	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	52	0	0	1	0
5	H	2	0	0	1	0
All	All	4330	0	4202	99	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 12.

All (99) 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:B:735:TYR:O	3:B:2:866:CL2	2.10	1.05
1:B:568:GLY:HA2	1:B:753:LEU:CD2	2.00	0.91
1:A:667:LYS:HD2	4:A:778:GOL:H11	1.62	0.82
1:A:634:MET:HE1	1:A:643:CYS:HB2	1.65	0.77
1:A:603:LEU:HB3	4:A:778:GOL:H12	1.68	0.76
1:B:736:CYS:HA	3:B:2:866:CL2	2.24	0.75
1:B:642:GLN:HB2	1:B:735:TYR:HD2	1.51	0.74
1:B:735:TYR:C	3:B:2:866:CL2	2.65	0.72
1:B:642:GLN:HB2	1:B:735:TYR:CD2	2.27	0.69
3:A:1:866:CL2	3:A:1:866:H1A	2.29	0.69
1:B:568:GLY:HA2	1:B:753:LEU:HD23	1.76	0.67
1:B:667:LYS:HZ3	4:B:778:GOL:H12	1.60	0.66
1:A:674:SER:HB2	5:A:818:HOH:O	1.96	0.65
1:A:631:GLU:O	1:A:635:THR:HG23	1.97	0.64
1:B:642:GLN:OE1	1:B:735:TYR:HB3	1.98	0.64
1:B:568:GLY:CA	1:B:753:LEU:HD23	2.28	0.64
1:B:579:LYS:HD3	2:D:749:LEU:HA	1.81	0.63
1:B:568:GLY:HA2	1:B:753:LEU:HD22	1.79	0.63
1:A:615:GLN:OE1	1:A:615:GLN:HA	1.98	0.63
1:A:593:MET:HG3	2:H:749:LEU:HD12	1.80	0.63
1:B:543:VAL:HG22	1:B:611:ARG:NH1	2.14	0.62
1:B:543:VAL:HG22	1:B:611:ARG:CZ	2.31	0.59
1:A:736:CYS:HA	3:A:1:866:CL2	2.39	0.59
2:D:747:TYR:O	2:D:751:LYS:HG3	2.02	0.59
1:B:568:GLY:CA	1:B:753:LEU:CD2	2.78	0.58
1:B:691:MET:HG3	1:B:695:LYS:HE3	1.87	0.56
1:A:543:VAL:HG22	1:A:611:ARG:NH2	2.20	0.56
1:B:566:LEU:O	1:B:570:GLN:HG2	2.04	0.56
1:B:667:LYS:NZ	4:B:778:GOL:H12	2.20	0.56
1:A:735:TYR:O	3:A:1:866:CL2	2.62	0.55
1:A:662:GLU:O	1:A:666:MET:HG3	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:528:LEU:O	1:A:530:PRO:HD3	2.09	0.53
1:B:755:GLU:OE1	2:D:742:ASN:HA	2.09	0.53
1:B:709:SER:C	1:B:711:ASN:H	2.12	0.53
1:A:611:ARG:O	1:A:615:GLN:HG2	2.08	0.53
1:B:652:GLU:HA	1:B:652:GLU:OE1	2.08	0.52
1:B:660:TYR:O	1:B:663:TYR:HB3	2.09	0.52
1:A:710:GLN:O	1:A:710:GLN:HG2	2.07	0.52
1:A:757:ILE:O	1:A:761:ILE:HG13	2.10	0.52
1:A:652:GLU:HA	1:A:652:GLU:OE1	2.10	0.52
1:A:565:MET:HE2	1:A:750:PRO:HD3	1.92	0.51
1:A:543:VAL:HA	1:A:611:ARG:NH2	2.26	0.50
1:A:713:GLN:O	1:A:717:GLN:HG3	2.11	0.50
2:D:744:LEU:HD12	2:D:745:LEU:H	1.77	0.50
2:D:748:LEU:HA	2:D:751:LYS:HD3	1.91	0.49
1:B:543:VAL:HA	1:B:611:ARG:NH1	2.28	0.49
1:A:565:MET:CE	1:A:750:PRO:HD3	2.42	0.49
1:A:750:PRO:HG2	1:A:753:LEU:HB2	1.95	0.49
1:B:755:GLU:CD	2:D:742:ASN:HA	2.33	0.49
3:A:1:866:H1A	3:A:1:866:C23	2.43	0.49
1:B:541:PRO:HG2	4:B:778:GOL:H2	1.94	0.49
1:B:579:LYS:HD3	2:D:749:LEU:HD23	1.94	0.48
1:B:735:TYR:CZ	3:B:2:866:H22	2.48	0.48
1:B:565:MET:HA	1:B:750:PRO:HG3	1.96	0.48
1:A:566:LEU:O	1:A:570:GLN:HG2	2.14	0.48
1:A:538:VAL:CG2	1:B:538:VAL:HG11	2.44	0.47
1:A:579:LYS:O	1:A:585:ARG:HG3	2.14	0.47
1:A:634:MET:HE1	1:A:640:TYR:HA	1.97	0.47
1:B:527:GLN:C	1:B:529:THR:N	2.66	0.47
1:A:644:LYS:HE2	1:A:645:HIS:CE1	2.51	0.46
1:B:556:THR:O	1:B:560:MET:HG3	2.16	0.46
1:B:736:CYS:CA	3:B:2:866:CL2	2.99	0.45
1:B:760:GLN:O	1:B:764:TYR:HD2	1.99	0.45
1:B:610:TRP:CZ2	1:B:660:TYR:HD1	2.35	0.45
1:A:735:TYR:CE2	3:A:1:866:H22	2.51	0.45
1:A:530:PRO:HD3	1:B:532:LEU:HD13	1.99	0.45
1:B:600:TRP:O	1:B:604:MET:HG2	2.15	0.45
1:B:610:TRP:HE3	1:B:653:LEU:HD13	1.82	0.44
1:B:761:ILE:HB	1:B:762:PRO:HD3	2.00	0.44
1:A:579:LYS:HD2	2:H:749:LEU:HA	1.99	0.44
1:B:575:VAL:O	1:B:579:LYS:HG3	2.17	0.44
1:B:619:ASN:C	1:B:620:LEU:HD23	2.38	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:757:ILE:O	1:B:761:ILE:HG12	2.17	0.44
2:H:742:ASN:HB2	5:H:36:HOH:O	2.17	0.44
1:B:527:GLN:C	1:B:529:THR:H	2.20	0.44
1:B:716:TYR:O	1:B:720:LYS:HB2	2.18	0.44
1:B:690:ARG:O	1:B:694:ILE:HG13	2.18	0.44
1:B:540:GLU:HA	1:B:541:PRO:HD3	1.90	0.43
1:A:634:MET:CE	1:A:634:MET:HA	2.49	0.42
3:B:2:866:CL2	3:B:2:866:H1A	2.56	0.42
1:A:755:GLU:OE1	2:H:742:ASN:OD1	2.37	0.42
1:B:613:TYR:CE1	1:B:654:HIS:HA	2.54	0.42
1:B:663:TYR:OH	1:B:667:LYS:HE2	2.19	0.42
1:A:719:THR:HB	1:A:774:PHE:CG	2.55	0.42
1:B:599:SER:HA	1:B:602:TYR:HD1	1.84	0.42
1:B:537:GLU:O	1:B:538:VAL:C	2.57	0.42
1:A:600:TRP:O	1:A:604:MET:HG2	2.20	0.42
1:A:753:LEU:O	1:A:757:ILE:HG13	2.20	0.42
1:B:674:SER:HB2	5:B:11:HOH:O	2.19	0.42
1:A:648:TYR:CZ	1:A:652:GLU:HG3	2.55	0.42
1:A:720:LYS:HD2	1:A:775:HIS:CD2	2.55	0.41
1:A:666:MET:HB3	1:A:722:LEU:HD21	2.01	0.41
1:A:543:VAL:HG22	1:A:611:ARG:CZ	2.50	0.41
1:B:619:ASN:O	1:B:620:LEU:HD23	2.19	0.41
1:A:543:VAL:HG22	1:A:611:ARG:HH21	1.83	0.41
1:B:735:TYR:CE2	3:B:2:866:H22	2.55	0.41
1:A:529:THR:HA	1:A:530:PRO:HD2	1.76	0.40
1:A:643:CYS:O	1:A:647:LEU:HG	2.21	0.40
1:A:538:VAL:HG23	1:B:538:VAL:HG11	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	244/257 (95%)	235 (96%)	9 (4%)	0	100	100
1	B	239/257 (93%)	225 (94%)	11 (5%)	3 (1%)	14	7
2	D	8/11 (73%)	7 (88%)	1 (12%)	0	100	100
2	H	9/11 (82%)	8 (89%)	1 (11%)	0	100	100
All	All	500/536 (93%)	475 (95%)	22 (4%)	3 (1%)	28	20

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	710	GLN
1	B	746	SER
1	B	538	VAL

### 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	217/235 (92%)	210 (97%)	7 (3%)	44	43
1	B	211/235 (90%)	208 (99%)	3 (1%)	71	77
2	D	9/10 (90%)	8 (89%)	1 (11%)	7	3
2	H	9/10 (90%)	9 (100%)	0	100	100
All	All	446/490 (91%)	435 (98%)	11 (2%)	53	55

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	529	THR
1	A	555	SER
1	A	619	ASN
1	A	674	SER
1	A	708	SER
1	A	753	LEU
1	A	764	TYR
1	B	615	GLN

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Mol	Chain	Res	Type
1	B	753	LEU
1	B	758	THR
2	D	744	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	760	GLN
1	B	654	HIS
1	B	759	ASN
1	B	776	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](#)

5 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	$\# Z  > 2$
3	866	A	1	-	36,39,39	1.60	6 (16%)	47,58,58	1.14	5 (10%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	GOL	A	778	-	5,5,5	0.36	0	5,5,5	0.51	0
3	866	B	2	-	36,39,39	1.53	5 (13%)	47,58,58	1.23	4 (8%)
4	GOL	B	3	-	5,5,5	0.32	0	5,5,5	0.35	0
4	GOL	B	778	-	5,5,5	0.36	0	5,5,5	0.17	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
3	866	A	1	-	-	0/35/38/38	0/3/3/3
4	GOL	A	778	-	-	0/4/4/4	0/0/0/0
3	866	B	2	-	-	0/35/38/38	0/3/3/3
4	GOL	B	3	-	-	0/4/4/4	0/0/0/0
4	GOL	B	778	-	-	0/4/4/4	0/0/0/0

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1	866	C3-C4	-4.08	1.51	1.53
3	B	2	866	C7-C6	-3.90	1.41	1.50
3	A	1	866	C7-C6	-3.77	1.42	1.50
3	B	2	866	C3-C4	-3.53	1.51	1.53
3	A	1	866	C18-C17	-3.44	1.46	1.51
3	B	2	866	C18-C17	-3.30	1.46	1.51
3	B	2	866	N3-N4	-3.10	1.34	1.39
3	A	1	866	N3-N4	-2.96	1.34	1.39
3	B	2	866	C10-N4	-2.85	1.38	1.44
3	A	1	866	C16-C4	-2.63	1.50	1.53
3	A	1	866	C10-N4	-2.58	1.38	1.44

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	2	866	C5-N2-C6	-3.46	117.47	123.02
3	B	2	866	F4-C16-C4	-3.08	110.00	112.22
3	A	1	866	C5-N2-C6	-2.63	118.81	123.02
3	A	1	866	C7-C6-N2	-2.40	112.68	117.45
3	A	1	866	F4-C16-C4	-2.37	110.51	112.22
3	A	1	866	O3-C17-N1	2.14	126.05	122.36
3	B	2	866	C23-C18-C19	2.44	120.02	116.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1	866	C23-C18-C19	2.46	120.05	116.72
3	B	2	866	C18-C23-CL2	3.08	123.77	119.62

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1	866	5	0
4	A	778	GOL	2	0
3	B	2	866	7	0
4	B	778	GOL	3	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 ⓘ

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	248/257 (96%)	0.62	26 (10%) 7 10	40, 70, 133, 162	2 (0%)
1	B	245/257 (95%)	0.75	38 (15%) 2 3	39, 74, 128, 170	1 (0%)
2	D	10/11 (90%)	0.48	0 100 100	53, 66, 88, 91	0
2	H	11/11 (100%)	0.24	0 100 100	56, 67, 101, 131	0
All	All	514/536 (95%)	0.67	64 (12%) 4 6	39, 72, 131, 170	3 (0%)

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	557	TRP	6.2
1	B	620	LEU	6.0
1	A	764	TYR	5.7
1	A	526	PRO	5.6
1	B	764	TYR	5.5
1	B	710	GLN	5.2
1	B	529	THR	4.6
1	A	707	ASN	4.4
1	B	551	SER	4.4
1	B	563	LEU	4.2
1	A	710	GLN	4.1
1	B	702	VAL	3.9
1	B	557	TRP	3.8
1	A	563	LEU	3.7
1	A	706	GLY	3.6
1	A	529	THR	3.5
1	B	765	SER	3.5
1	A	552	VAL	3.4
1	B	611	ARG	3.3
1	B	762	PRO	3.0
1	B	709	SER	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	763	LYS	2.9
1	B	744	THR	2.8
1	B	553	PRO	2.8
1	A	705	GLU	2.8
1	A	709	SER	2.7
1	A	708	SER	2.7
1	A	677	LYS	2.7
1	B	618	ALA	2.7
1	A	763	LYS	2.7
1	B	621	LEU	2.7
1	A	549	ASP	2.7
1	B	746	SER	2.6
1	A	553	PRO	2.6
1	A	611	ARG	2.6
1	B	679	GLY	2.6
1	B	596	LEU	2.6
1	B	550	SER	2.6
1	B	767	GLY	2.6
1	A	743	LYS	2.5
1	B	766	ASN	2.5
1	B	776	GLN	2.5
1	B	741	LEU	2.5
1	B	718	LEU	2.5
1	A	671	LEU	2.5
1	A	762	PRO	2.4
1	B	669	LEU	2.4
1	B	654	HIS	2.4
1	A	766	ASN	2.4
1	A	742	ASP	2.4
1	B	681	LYS	2.3
1	A	560	MET	2.3
1	A	718	LEU	2.3
1	B	545	TYR	2.3
1	B	527	GLN	2.3
1	B	756	ILE	2.3
1	B	544	LEU	2.2
1	A	752	MET	2.2
1	B	668	THR	2.2
1	B	613	TYR	2.2
1	B	556	THR	2.2
1	B	543	VAL	2.2
1	A	744	THR	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	643	CYS	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
4	GOL	A	778	6/6	0.88	0.23	2.83	92,96,99,102	0
4	GOL	B	778	6/6	0.94	0.18	1.22	74,89,94,99	0
4	GOL	B	3	6/6	0.63	0.24	1.04	88,93,95,95	0
3	866	A	1	37/37	0.94	0.13	-0.55	41,52,65,77	0
3	866	B	2	37/37	0.92	0.13	-0.63	35,51,67,83	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.