



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 04:27 am BST

PDB ID : 5X33
Title : Leukotriene B4 receptor BLT1 in complex with BIIL260
Authors : Hori, T.; Hirata, K.; Yamashita, K.; Kawano, Y.; Yamamoto, M.; Yokoyama, S.
Deposited on : 2017-02-03
Resolution : 3.70 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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

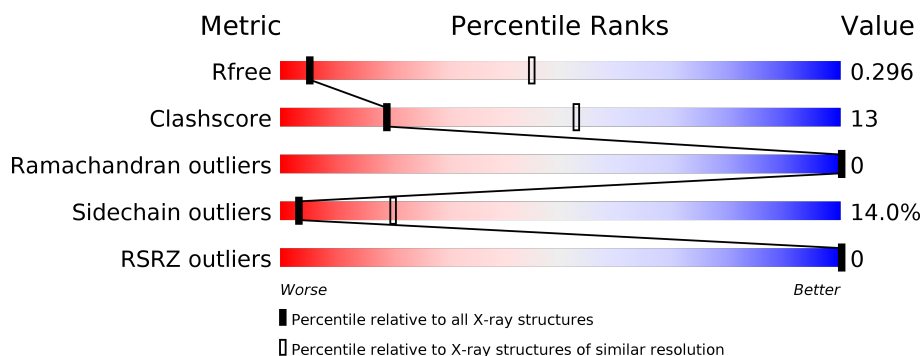
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.70 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}	130704	1049 (3.88-3.52)
Clashscore	141614	1027 (3.86-3.54)
Ramachandran outliers	138981	1069 (3.88-3.52)
Sidechain outliers	138945	1065 (3.88-3.52)
RSRZ outliers	127900	1578 (3.90-3.50)

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 ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	523	<div> <div></div> <div>55%</div> <div>25%</div> <div>•</div> <div>16%</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3375 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 LTB4 receptor, Lysozyme, LTB4 receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	439	Total	C	N	O	S	0	0	0
			3340	2163	582	579	16			

There are 35 discrepancies between the modelled and reference sequences:

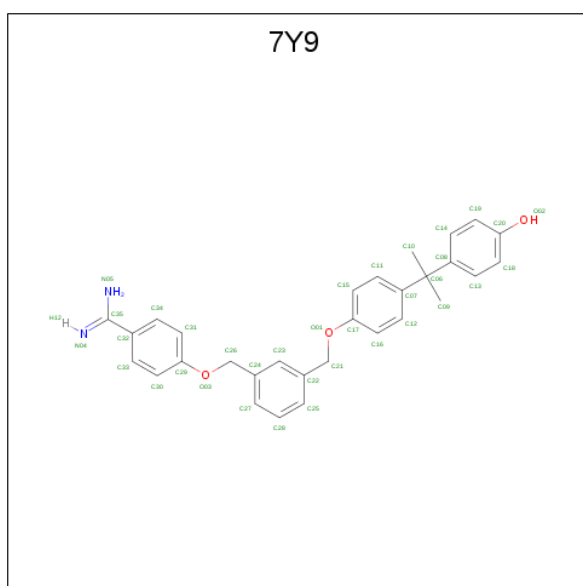
Chain	Residue	Modelled	Actual	Comment	Reference
A	2	GLU	-	expression tag	UNP Q9WTK1
A	3	LEU	-	expression tag	UNP Q9WTK1
A	4	MET	-	expression tag	UNP Q9WTK1
A	5	ASP	-	expression tag	UNP Q9WTK1
A	6	TYR	-	expression tag	UNP Q9WTK1
A	7	LYS	-	expression tag	UNP Q9WTK1
A	8	ASP	-	expression tag	UNP Q9WTK1
A	9	ASP	-	expression tag	UNP Q9WTK1
A	10	ASP	-	expression tag	UNP Q9WTK1
A	11	ASP	-	expression tag	UNP Q9WTK1
A	12	LYS	-	expression tag	UNP Q9WTK1
A	13	GLU	-	expression tag	UNP Q9WTK1
A	14	PHE	-	expression tag	UNP Q9WTK1
A	83	GLY	HIS	engineered mutation	UNP Q9WTK1
A	88	GLY	LYS	engineered mutation	UNP Q9WTK1
A	212	ALA	VAL	engineered mutation	UNP Q9WTK1
A	900	GLY	-	linker	UNP Q9WTK1
A	901	SER	-	linker	UNP Q9WTK1
A	902	GLY	-	linker	UNP Q9WTK1
A	903	SER	-	linker	UNP Q9WTK1
A	1054	THR	CYS	engineered mutation	UNP A0A097J792
A	1097	ALA	CYS	engineered mutation	UNP A0A097J792
A	1200	GLY	-	linker	UNP A0A097J792
A	1201	SER	-	linker	UNP A0A097J792
A	1202	GLY	-	linker	UNP A0A097J792
A	1203	SER	-	linker	UNP A0A097J792
A	309	ALA	SER	engineered mutation	UNP Q9WTK1

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Chain	Residue	Modelled	Actual	Comment	Reference
A	349	GLU	-	expression tag	UNP Q9WTK1
A	350	PHE	-	expression tag	UNP Q9WTK1
A	351	LEU	-	expression tag	UNP Q9WTK1
A	352	GLU	-	expression tag	UNP Q9WTK1
A	353	VAL	-	expression tag	UNP Q9WTK1
A	354	LEU	-	expression tag	UNP Q9WTK1
A	355	PHE	-	expression tag	UNP Q9WTK1
A	356	GLN	-	expression tag	UNP Q9WTK1

- Molecule 2 is 4-[[3-[[4-[2-(4-hydroxyphenyl)propan-2-yl]phenoxy]methyl]phenyl]methoxy]benzenecarboximidamide (three-letter code: 7Y9) (formula: C₃₀H₃₀N₂O₃).

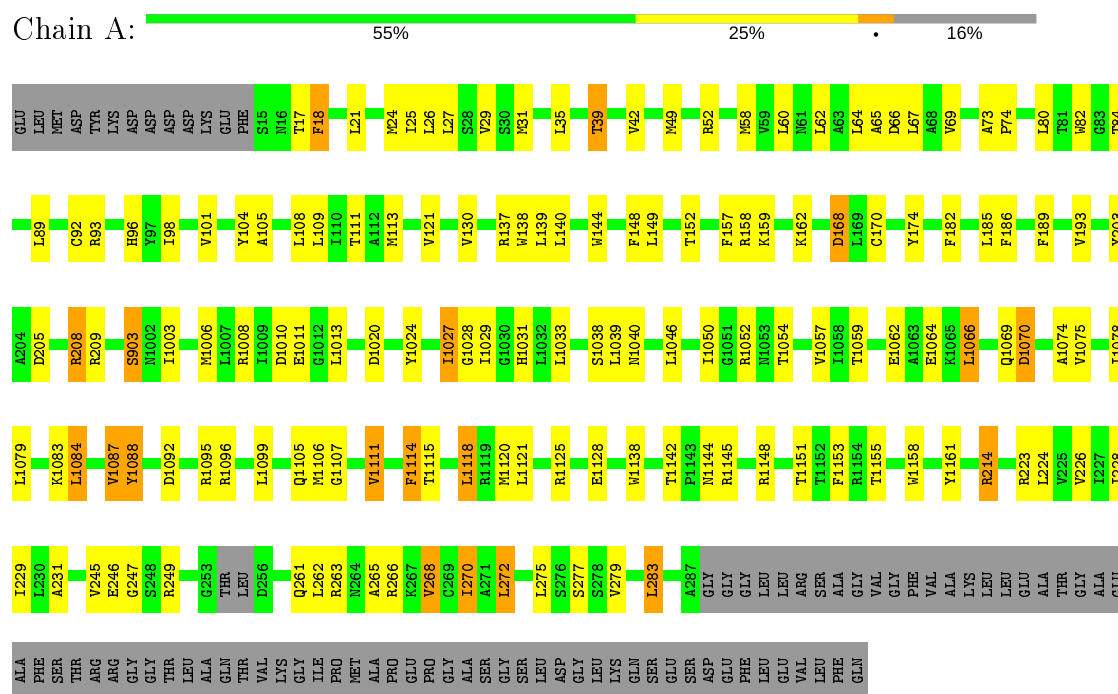


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			35	30	2	3		

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.

- Molecule 1: LTB4 receptor, Lysozyme, LTB4 receptor



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, α , β , γ	69.56Å 77.62Å 135.53Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.54 – 3.70 48.54 – 3.70	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.54-3.70) 89.4 (48.54-3.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.77 (at 3.67Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.248 , 0.296 0.248 , 0.296	Depositor DCC
R_{free} test set	414 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	49.3	Xtriage
Anisotropy	0.767	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 35.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.85	EDS
Total number of atoms	3375	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.30% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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: 7Y9

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	0.30	0/3409	0.51	0/4638

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	3340	0	3375	88	0
2	A	35	0	0	2	0
All	All	3375	0	3375	88	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 13.

All (88) 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:1010:ASP:HB3	1:A:1145:ARG:HD2	1.37	1.05
1:A:189:PHE:HA	1:A:193:VAL:HB	1.64	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1120:MET:HG2	1:A:1125:ARG:HD2	1.67	0.75
1:A:35:LEU:O	1:A:39:THR:OG1	2.03	0.74
1:A:1105:GLN:HB2	1:A:1145:ARG:NH2	2.07	0.69
1:A:903:SER:OG	1:A:1064:GLU:OE1	2.12	0.67
1:A:113:MET:HE1	1:A:228:ILE:HD11	1.78	0.65
1:A:262:LEU:O	1:A:266:ARG:N	2.25	0.65
1:A:1092:ASP:OD2	1:A:1095:ARG:NH1	2.29	0.65
1:A:1078:ILE:HG23	1:A:1084:LEU:HD12	1.81	0.63
1:A:228:ILE:HG22	1:A:283:LEU:HD23	1.80	0.62
1:A:1105:GLN:HB2	1:A:1145:ARG:HH22	1.65	0.61
1:A:1046:LEU:HG	1:A:1050:ILE:HD13	1.81	0.61
1:A:101:VAL:HG21	1:A:148:PHE:HD1	1.66	0.60
1:A:1038:SER:OG	1:A:1040:ASN:OD1	2.19	0.60
1:A:214:ARG:HB3	1:A:214:ARG:HH11	1.67	0.60
1:A:1087:VAL:HG11	1:A:1118:LEU:HD12	1.84	0.60
1:A:1010:ASP:CB	1:A:1145:ARG:HD2	2.24	0.59
1:A:149:LEU:O	1:A:152:THR:OG1	2.19	0.59
1:A:96:HIS:ND1	1:A:158:ARG:HD3	2.18	0.59
1:A:182:PHE:O	1:A:186:PHE:N	2.36	0.55
1:A:66:ASP:OD1	2:A:2001:7Y9:N04	2.38	0.55
1:A:1033:LEU:HD13	1:A:1046:LEU:HB2	1.89	0.55
1:A:1125:ARG:NH1	1:A:1128:GLU:OE2	2.40	0.54
1:A:1031:HIS:ND1	1:A:1070:ASP:OD2	2.41	0.53
1:A:60:LEU:O	1:A:64:LEU:N	2.42	0.53
1:A:25:ILE:O	1:A:29:VAL:HG13	2.08	0.53
1:A:157:PHE:HD2	1:A:174:TYR:CD1	2.27	0.53
1:A:185:LEU:HD13	1:A:247:GLY:HA3	1.92	0.52
1:A:1083:LYS:HD3	1:A:1115:THR:HG22	1.91	0.52
1:A:168:ASP:OD1	1:A:168:ASP:N	2.43	0.52
1:A:157:PHE:HD2	1:A:174:TYR:HD1	1.59	0.51
1:A:65:ALA:HB2	1:A:144:TRP:HH2	1.74	0.51
1:A:162:LYS:HA	1:A:168:ASP:HB3	1.92	0.51
1:A:1105:GLN:OE1	1:A:1138:TRP:NE1	2.43	0.51
1:A:1027:ILE:HG22	1:A:1033:LEU:HD21	1.91	0.51
1:A:1151:THR:O	1:A:1155:THR:OG1	2.29	0.51
1:A:266:ARG:O	1:A:270:ILE:HB	2.10	0.50
1:A:1144:ASN:O	1:A:1148:ARG:NH1	2.45	0.50
1:A:89:LEU:O	1:A:93:ARG:HB2	2.12	0.49
1:A:208:ARG:HB3	1:A:208:ARG:CZ	2.42	0.49
1:A:96:HIS:CG	1:A:158:ARG:HD3	2.48	0.48
1:A:214:ARG:HB3	1:A:214:ARG:NH1	2.29	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:65:ALA:HB2	1:A:144:TRP:CH2	2.47	0.48
1:A:58:MET:O	1:A:62:LEU:HB3	2.13	0.48
1:A:49:MET:HE3	1:A:52:ARG:HA	1.96	0.47
1:A:26:LEU:HB3	1:A:275:LEU:HD11	1.97	0.47
1:A:1011:GLU:HG2	1:A:1145:ARG:NH1	2.29	0.47
1:A:1008:ARG:HE	1:A:1013:LEU:HD22	1.80	0.47
1:A:84:THR:OG1	1:A:168:ASP:OD1	2.31	0.47
1:A:1088:TYR:HE1	1:A:1096:ARG:HD3	1.80	0.47
1:A:268:VAL:O	1:A:272:LEU:N	2.33	0.47
1:A:1006:MET:HG3	1:A:1161:TYR:CE2	2.50	0.46
1:A:1020:ASP:N	1:A:1024:TYR:O	2.38	0.46
1:A:1107:GLY:O	1:A:1111:VAL:HG12	2.15	0.46
1:A:31:MET:HE3	1:A:74:PRO:HD3	1.98	0.46
1:A:1092:ASP:O	1:A:1096:ARG:HG3	2.16	0.46
1:A:1054:THR:HB	1:A:1057:VAL:O	2.15	0.46
1:A:1027:ILE:HG12	1:A:1028:GLY:N	2.31	0.45
1:A:1105:GLN:CD	1:A:1145:ARG:HH21	2.19	0.45
1:A:261:GLN:O	1:A:265:ALA:N	2.42	0.45
1:A:69:VAL:HG21	2:A:2001:7Y9:C34	2.46	0.45
1:A:1121:LEU:HD22	1:A:1153:PHE:HD1	1.81	0.45
1:A:1050:ILE:HG22	1:A:1052:ARG:HG2	1.98	0.45
1:A:1062:GLU:O	1:A:1066:LEU:HB2	2.17	0.45
1:A:1069:GLN:OE1	1:A:1069:GLN:N	2.50	0.45
1:A:1114:PHE:O	1:A:1118:LEU:HB2	2.17	0.44
1:A:27:LEU:O	1:A:31:MET:HG3	2.17	0.44
1:A:24:MET:HB3	1:A:80:LEU:HD23	1.99	0.44
1:A:1158:TRP:HB3	1:A:214:ARG:HH21	1.82	0.44
1:A:42:VAL:HG21	1:A:64:LEU:HD13	1.99	0.44
1:A:223:ARG:O	1:A:226:VAL:HG12	2.18	0.43
1:A:231:ALA:HB1	1:A:279:VAL:HG21	2.00	0.43
1:A:1121:LEU:HD13	1:A:1153:PHE:HE1	1.83	0.43
1:A:21:LEU:O	1:A:24:MET:HG2	2.18	0.43
1:A:1074:ALA:O	1:A:1078:ILE:HG13	2.19	0.42
1:A:1114:PHE:CD1	1:A:1114:PHE:N	2.86	0.42
1:A:104:TYR:O	1:A:108:LEU:HG	2.20	0.42
1:A:1011:GLU:HB2	1:A:1029:ILE:HG23	2.00	0.42
1:A:1011:GLU:HG2	1:A:1145:ARG:HH12	1.85	0.41
1:A:1121:LEU:HD13	1:A:1153:PHE:CE1	2.55	0.41
1:A:66:ASP:O	1:A:277:SER:OG	2.30	0.41
1:A:1114:PHE:O	1:A:1118:LEU:N	2.54	0.41
1:A:18:PHE:HA	1:A:21:LEU:HB3	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:159:LYS:HB2	1:A:159:LYS:HZ2	1.86	0.40
1:A:246:GLU:HG3	1:A:262:LEU:HD22	2.03	0.40
1:A:105:ALA:O	1:A:109:LEU:HG	2.22	0.40
1:A:31:MET:HE3	1:A:73:ALA:HB3	2.04	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	435/523 (83%)	421 (97%)	14 (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	342/429 (80%)	294 (86%)	48 (14%)	3	20

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	THR
1	A	18	PHE

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Mol	Chain	Res	Type
1	A	39	THR
1	A	67	LEU
1	A	82	TRP
1	A	92	CYS
1	A	98	ILE
1	A	111	THR
1	A	121	VAL
1	A	130	VAL
1	A	137	ARG
1	A	138	TRP
1	A	139	LEU
1	A	140	LEU
1	A	168	ASP
1	A	170	CYS
1	A	203	TYR
1	A	205	ASP
1	A	208	ARG
1	A	209	ARG
1	A	903	SER
1	A	1003	ILE
1	A	1027	ILE
1	A	1039	LEU
1	A	1059	THR
1	A	1066	LEU
1	A	1070	ASP
1	A	1075	VAL
1	A	1079	LEU
1	A	1084	LEU
1	A	1087	VAL
1	A	1088	TYR
1	A	1099	LEU
1	A	1106	MET
1	A	1111	VAL
1	A	1114	PHE
1	A	1118	LEU
1	A	1142	THR
1	A	214	ARG
1	A	224	LEU
1	A	229	ILE
1	A	245	VAL
1	A	249	ARG
1	A	263	ARG

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Mol	Chain	Res	Type
1	A	268	VAL
1	A	270	ILE
1	A	272	LEU
1	A	283	LEU

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	1123	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](#)

1 ligand is 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	7Y9	A	2001	-	38,38,38	1.10	3 (7%)	49,53,53	0.96	3 (6%)

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	7Y9	A	2001	-	-	2/26/26/26	0/4/4/4

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	2001	7Y9	C35-N04	4.00	1.45	1.28
2	A	2001	7Y9	C35-N05	-3.22	1.25	1.33
2	A	2001	7Y9	C32-C35	2.25	1.51	1.47

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	2001	7Y9	O01-C21-C22	2.51	116.68	109.16
2	A	2001	7Y9	C12-C07-C06	-2.38	115.73	121.52
2	A	2001	7Y9	C34-C32-C33	2.14	121.64	118.59

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	2001	7Y9	C15-C17-O01-C21
2	A	2001	7Y9	C16-C17-O01-C21

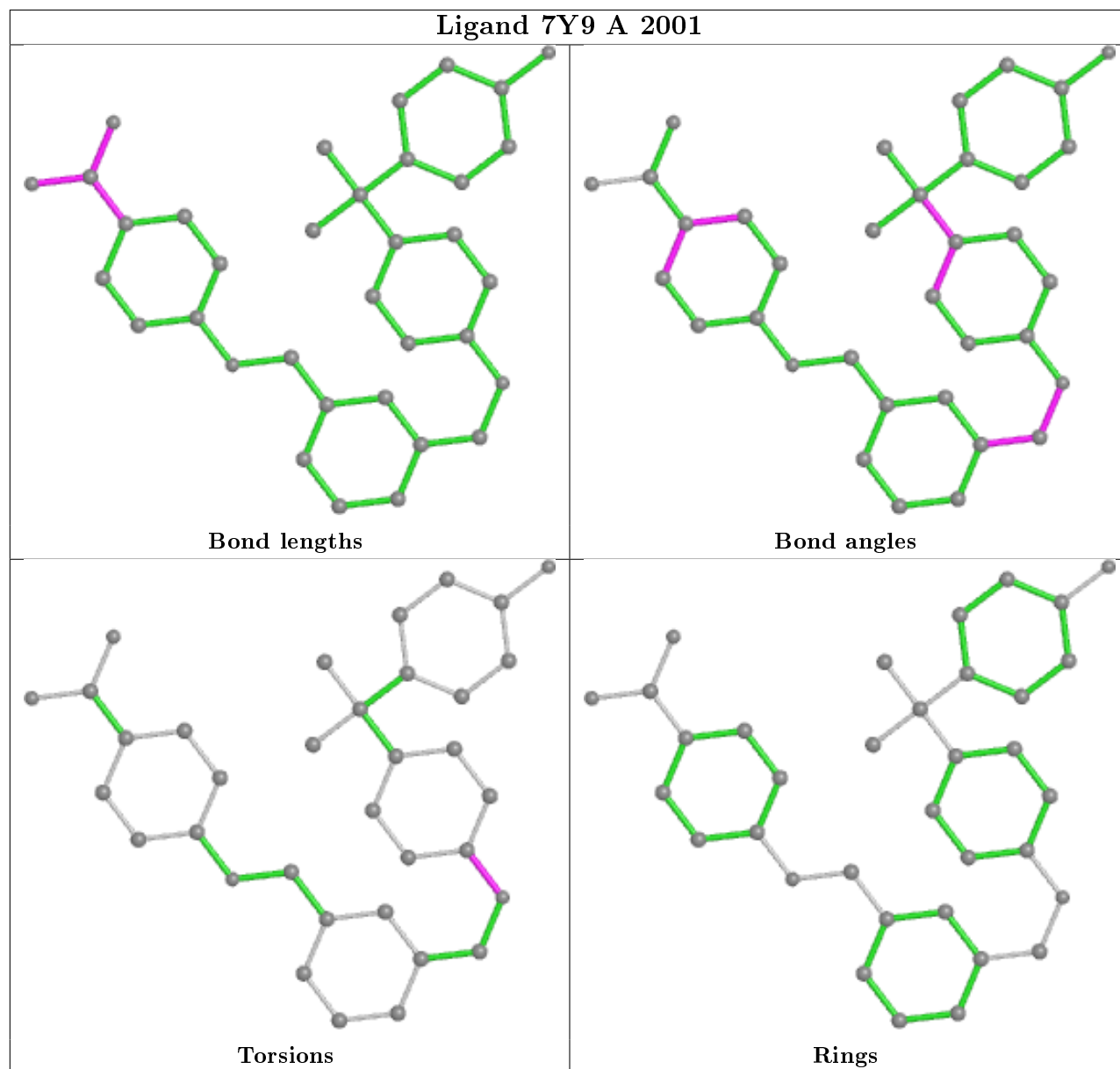
There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	2001	7Y9	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and

any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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, 95th 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 ²)	Q<0.9
1	A	439/523 (83%)	-0.45	0 100 100	45, 67, 89, 118	0

There are no RSRZ outliers to report.

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](#)

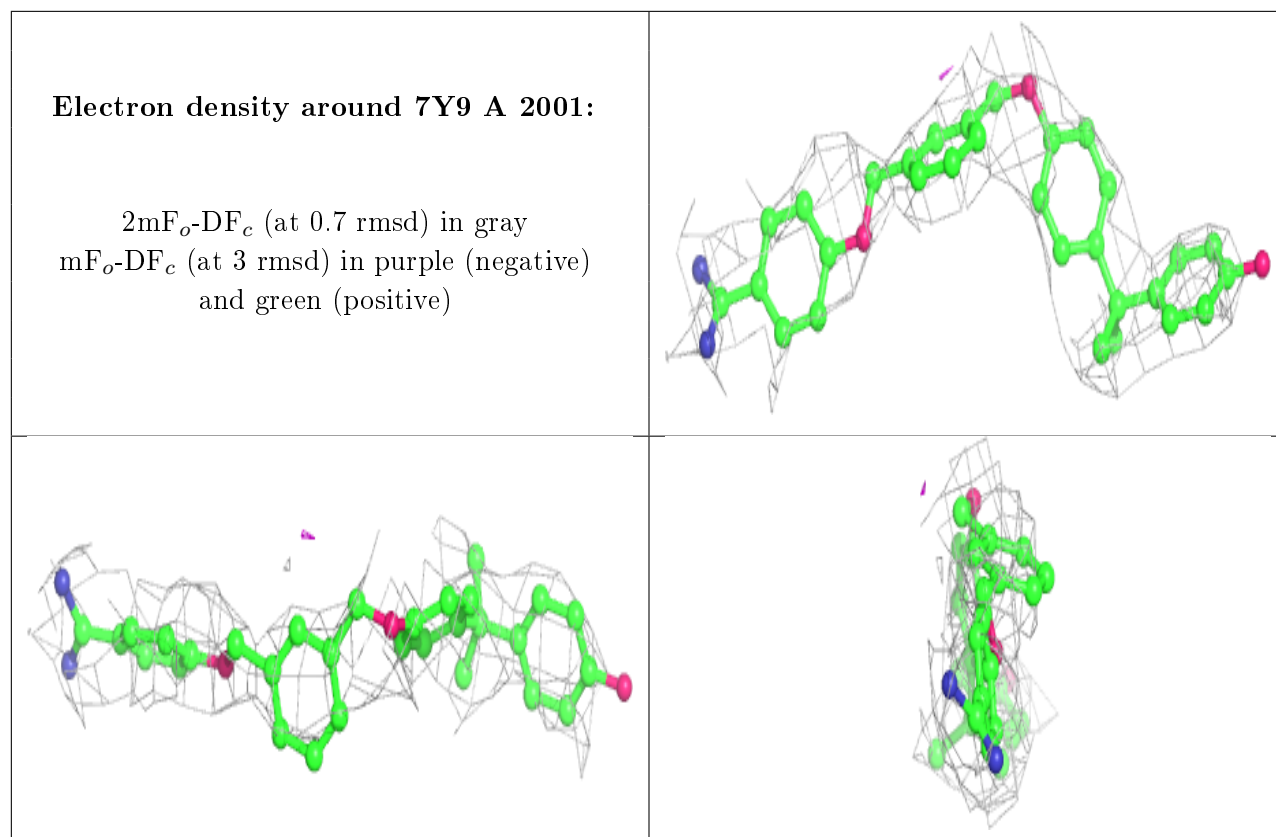
There are no carbohydrates in this entry.

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. The B-factors column lists the minimum, median, 95th 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	B-factors(Å ²)	Q<0.9
2	7Y9	A	2001	35/35	0.90	0.33	60,77,92,94	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers [i](#)

There are no such residues in this entry.