



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:11 AM UTC

PDB ID : 6ZLY / pdb_00006zly
Title : Crystal structure of the complex between PPARgamma LBD and the ligand NV1362 (7a)
Authors : Pochetti, G.; Montanari, R.; Capelli, D.
Deposited on : 2020-07-01
Resolution : 1.79 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

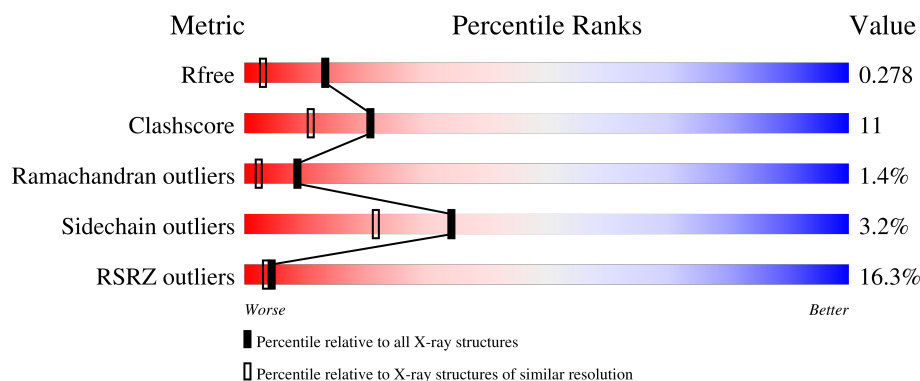
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.79 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}	180053	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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 of 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	300	10% 64% 18% • 16%
1	B	300	17% 65% 14% • 19%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4161 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 Peroxisome proliferator-activated receptor gamma.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	252	Total	C	N	O	S	0	1	0
			2020	1306	328	376	10			
1	B	244	Total	C	N	O	S	0	0	0
			1954	1264	319	362	9			

There are 34 discrepancies between the modelled and reference sequences:

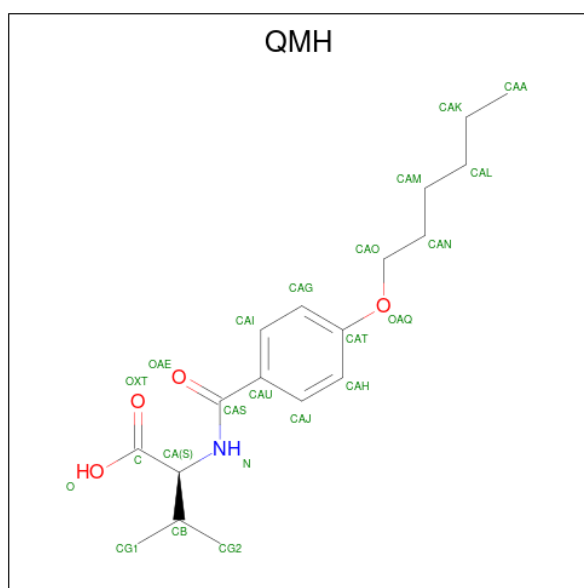
Chain	Residue	Modelled	Actual	Comment	Reference
A	178	HIS	-	expression tag	UNP P37231
A	179	HIS	-	expression tag	UNP P37231
A	180	HIS	-	expression tag	UNP P37231
A	181	HIS	-	expression tag	UNP P37231
A	182	HIS	-	expression tag	UNP P37231
A	183	HIS	-	expression tag	UNP P37231
A	184	SER	-	expression tag	UNP P37231
A	185	SER	-	expression tag	UNP P37231
A	186	GLY	-	expression tag	UNP P37231
A	187	LEU	-	expression tag	UNP P37231
A	188	VAL	-	expression tag	UNP P37231
A	189	PRO	-	expression tag	UNP P37231
A	190	ARG	-	expression tag	UNP P37231
A	191	GLY	-	expression tag	UNP P37231
A	192	SER	-	expression tag	UNP P37231
A	193	HIS	-	expression tag	UNP P37231
A	194	MET	-	expression tag	UNP P37231
B	178	HIS	-	expression tag	UNP P37231
B	179	HIS	-	expression tag	UNP P37231
B	180	HIS	-	expression tag	UNP P37231
B	181	HIS	-	expression tag	UNP P37231
B	182	HIS	-	expression tag	UNP P37231
B	183	HIS	-	expression tag	UNP P37231
B	184	SER	-	expression tag	UNP P37231
B	185	SER	-	expression tag	UNP P37231

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Chain	Residue	Modelled	Actual	Comment	Reference
B	186	GLY	-	expression tag	UNP P37231
B	187	LEU	-	expression tag	UNP P37231
B	188	VAL	-	expression tag	UNP P37231
B	189	PRO	-	expression tag	UNP P37231
B	190	ARG	-	expression tag	UNP P37231
B	191	GLY	-	expression tag	UNP P37231
B	192	SER	-	expression tag	UNP P37231
B	193	HIS	-	expression tag	UNP P37231
B	194	MET	-	expression tag	UNP P37231

- Molecule 2 is (2 {S})-2-[(4-hexoxyphenyl)carbonylamino]-3-methyl-butanoic acid (CCD ID: QMH) (formula: C₁₈H₂₇NO₄) (labeled as "Ligand of Interest" by depositor).

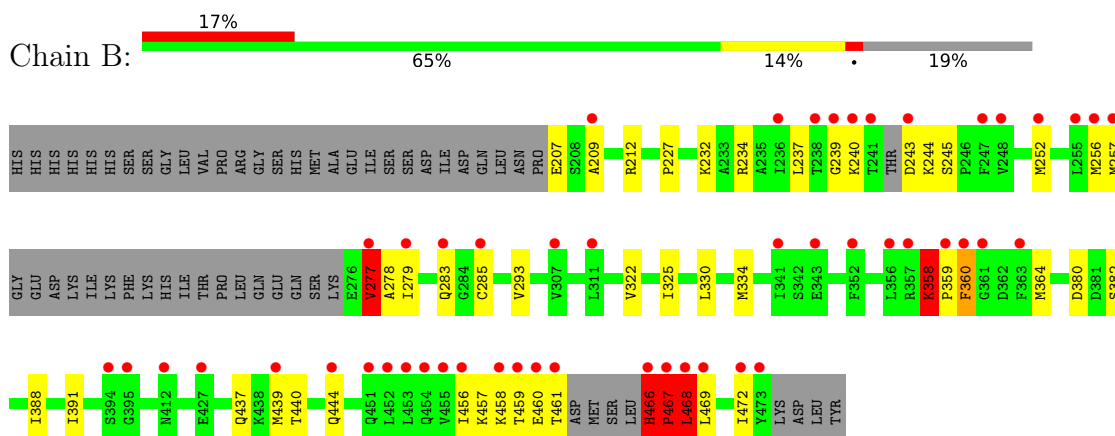


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	N	O	
			23	18	1	4	

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	94	Total	O		
			94	94	0	0
3	B	70	Total	O		
			70	70	0	0

- Molecule 1: Peroxisome proliferator-activated receptor gamma



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	92.89Å 60.07Å 118.02Å 90.00° 103.34° 90.00°	Depositor
Resolution (Å)	57.42 – 1.79 57.42 – 1.79	Depositor EDS
% Data completeness (in resolution range)	96.2 (57.42-1.79) 96.2 (57.42-1.79)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.57 (at 1.80Å)	Xtriage
Refinement program	PHENIX 1.8.4_1496	Depositor
R, R_{free}	0.234 , 0.280 0.236 , 0.278	Depositor DCC
R_{free} test set	2911 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å ²)	32.6	Xtriage
Anisotropy	0.479	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 36.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4161	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.59% 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 ⓘ

5.1 Standard geometry ⓘ

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

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	0.56	0/2058	0.88	4/2773 (0.1%)
1	B	0.58	0/1985	0.96	8/2673 (0.3%)
All	All	0.57	0/4043	0.92	12/5446 (0.2%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	466	HIS	CA-C-N	10.91	133.48	119.84
1	B	466	HIS	C-N-CA	10.91	133.48	119.84
1	A	358	LYS	CA-C-N	9.27	131.42	119.84
1	A	358	LYS	C-N-CA	9.27	131.42	119.84
1	B	468	LEU	N-CA-C	-8.75	102.59	113.18
1	A	365	GLU	CA-C-N	-6.35	113.14	119.56
1	A	365	GLU	C-N-CA	-6.35	113.14	119.56
1	B	360	PHE	N-CA-C	5.96	119.81	112.54
1	B	277	VAL	N-CA-C	5.93	121.67	109.34
1	B	239	GLY	N-CA-C	-5.75	103.73	111.54
1	B	467	PRO	CB-CA-C	-5.57	102.37	111.56
1	B	444	GLN	CA-CB-CG	-5.22	103.66	114.10

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	2020	0	2075	60	0
1	B	1954	0	2010	34	0
2	A	23	0	0	1	0
3	A	94	0	0	7	0
3	B	70	0	0	4	0
All	All	4161	0	4085	90	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 11.

All (90) 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:358:LYS:HB3	1:A:359:PRO:CD	1.15	1.53
1:A:358:LYS:CB	1:A:359:PRO:CD	1.98	1.32
1:A:358:LYS:CB	1:A:359:PRO:HD3	1.58	1.28
1:A:358:LYS:HB3	1:A:359:PRO:HD2	1.29	1.09
1:A:358:LYS:CB	1:A:359:PRO:HD2	1.82	1.06
1:B:358:LYS:HB3	1:B:359:PRO:CD	1.86	1.03
1:B:467:PRO:HG2	1:B:468:LEU:H	1.23	1.00
1:B:358:LYS:HB3	1:B:359:PRO:HD3	1.43	0.99
1:B:277:VAL:HG23	3:B:502:HOH:O	1.63	0.96
1:B:325:ILE:HD12	1:B:388:ILE:HG23	1.63	0.78
1:B:293:VAL:HG22	1:B:322:VAL:HG21	1.66	0.78
1:A:358:LYS:HB2	1:A:359:PRO:HD2	1.66	0.77
1:A:290:VAL:HG23	1:A:468:LEU:HD23	1.70	0.73
1:B:467:PRO:CG	1:B:468:LEU:H	1.94	0.70
1:A:358:LYS:HB3	1:A:359:PRO:HD3	0.67	0.67
1:B:277:VAL:HG23	1:B:278:ALA:H	1.59	0.66
1:A:249:ILE:HD12	1:A:255:LEU:HA	1.78	0.65
1:B:330:LEU:HD21	1:B:364:MET:HE1	1.80	0.64
1:B:358:LYS:CB	1:B:359:PRO:CD	2.64	0.63
1:A:349:THR:HG21	3:A:681:HOH:O	1.99	0.63
1:A:363[B]:PHE:HE1	1:A:449:HIS:HE1	1.48	0.62
1:A:430:GLN:HG3	1:A:433:ALA:HB3	1.83	0.60
1:B:466:HIS:HB3	3:B:563:HOH:O	2.01	0.59
1:A:279:ILE:N	3:A:601:HOH:O	2.36	0.59
1:B:467:PRO:HG2	1:B:468:LEU:N	2.05	0.59
1:B:279:ILE:O	1:B:283:GLN:HG3	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:439:MET:HE1	1:B:437:GLN:HG2	1.86	0.58
1:A:349:THR:HG22	1:A:351:GLU:H	1.69	0.57
1:B:227:PRO:O	1:B:232:LYS:NZ	2.22	0.57
1:A:363[B]:PHE:CE1	1:A:449:HIS:HE1	2.23	0.56
1:A:319:LYS:NZ	1:A:471:GLU:O	2.39	0.56
1:A:207:GLU:HG3	1:A:209:ALA:H	1.70	0.56
1:A:235:ALA:HA	1:A:240:LYS:HG2	1.86	0.56
1:B:252:MET:SD	1:B:277:VAL:HG11	2.45	0.56
1:A:276:GLU:HG2	1:A:278:ALA:H	1.72	0.55
1:A:325:ILE:HD11	1:A:392:ILE:HG13	1.87	0.55
1:A:279:ILE:O	1:A:283:GLN:HG2	2.08	0.54
1:A:243:ASP:OD1	1:A:244:LYS:N	2.41	0.54
1:A:443:ARG:NH1	3:A:607:HOH:O	2.40	0.54
1:A:436:LEU:HA	1:A:439:MET:HE2	1.90	0.54
1:B:467:PRO:CG	1:B:468:LEU:N	2.60	0.54
1:A:370:PHE:HB2	1:A:445:ILE:HD11	1.88	0.54
1:A:287:PHE:O	1:A:290:VAL:HG12	2.08	0.53
1:A:370:PHE:CB	1:A:445:ILE:HD11	2.40	0.51
1:A:349:THR:HG22	1:A:351:GLU:N	2.26	0.50
1:A:437:GLN:O	1:A:440:THR:HG22	2.11	0.50
1:A:363[A]:PHE:HE2	1:A:456:ILE:HG12	1.77	0.49
1:B:466:HIS:HB2	1:B:469:LEU:HD23	1.95	0.48
1:A:252:MET:O	1:A:256:MET:HG2	2.13	0.48
1:B:457:LYS:C	1:B:458:LYS:HD3	2.39	0.48
1:A:357:ARG:HG3	1:A:358:LYS:HB2	1.94	0.48
1:A:360:PHE:HA	1:A:363[A]:PHE:HD2	1.80	0.47
1:A:437:GLN:HG2	1:B:439:MET:HE1	1.96	0.47
1:B:459:THR:HG23	1:B:460:GLU:N	2.29	0.47
1:A:379:LEU:HD11	1:A:435:LEU:HD13	1.96	0.47
1:A:216:LYS:NZ	3:A:605:HOH:O	2.30	0.47
1:B:359:PRO:HG2	1:B:360:PHE:CD2	2.49	0.47
1:A:241:THR:HB	3:A:678:HOH:O	2.14	0.47
1:B:207:GLU:HG3	1:B:209:ALA:H	1.81	0.46
1:A:252:MET:SD	1:A:277:VAL:HG11	2.55	0.46
1:A:325:ILE:HD11	1:A:392:ILE:CG1	2.46	0.46
1:B:459:THR:HG23	1:B:460:GLU:H	1.80	0.46
1:A:325:ILE:HG23	1:A:388:ILE:HD12	1.96	0.46
1:A:465:LEU:HG	1:A:470:GLN:HG2	1.97	0.46
1:B:456:ILE:O	1:B:459:THR:HG22	2.16	0.45
1:A:293:VAL:HG22	1:A:322:VAL:HG11	1.97	0.45
1:A:259:GLU:HB2	3:A:669:HOH:O	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:357:ARG:HD2	1:A:358:LYS:HG3	1.99	0.45
1:A:341:ILE:HD13	1:A:348:MET:HE3	1.99	0.45
1:B:252:MET:O	1:B:256:MET:HG2	2.18	0.44
1:A:439:MET:CE	1:B:437:GLN:HG2	2.46	0.44
1:B:388:ILE:HG13	3:B:524:HOH:O	2.17	0.44
1:A:252:MET:HE2	3:A:652:HOH:O	2.18	0.44
1:B:234:ARG:NH2	1:B:334:MET:O	2.35	0.44
1:B:325:ILE:HD11	1:B:391:ILE:HB	2.00	0.43
1:A:276:GLU:CG	1:A:277:VAL:N	2.81	0.43
1:A:363[B]:PHE:CD2	1:A:364:MET:HG2	2.53	0.43
1:A:363[B]:PHE:HD2	1:A:364:MET:HG2	1.82	0.43
1:A:207:GLU:HG3	1:A:209:ALA:N	2.34	0.42
1:A:240:LYS:HG3	1:A:241:THR:H	1.84	0.42
1:B:468:LEU:O	1:B:472:ILE:HG13	2.18	0.42
1:A:443:ARG:NH2	1:B:440:THR:OG1	2.52	0.42
1:A:252:MET:SD	1:A:277:VAL:HG21	2.60	0.41
1:A:241:THR:HG23	1:A:241:THR:O	2.21	0.41
1:B:285:CYS:SG	3:B:567:HOH:O	2.62	0.41
1:A:238:THR:O	1:A:239:GLY:O	2.39	0.41
1:A:363[A]:PHE:CE2	1:A:456:ILE:HG12	2.54	0.41
1:B:380:ASP:OD1	1:B:382:SER:OG	2.32	0.40
1:A:438:LYS:HD3	1:A:438:LYS:HA	1.92	0.40
1:A:286:GLN:HG2	2:A:501:QMH:CG2	2.52	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	249/300 (83%)	239 (96%)	7 (3%)	3 (1%)	10 3
1	B	236/300 (79%)	226 (96%)	6 (2%)	4 (2%)	7 2

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	485/600 (81%)	465 (96%)	13 (3%)	7 (1%)	9 2

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	358	LYS
1	B	277	VAL
1	B	358	LYS
1	A	239	GLY
1	B	237	LEU
1	B	467	PRO
1	A	359	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	226/270 (84%)	222 (98%)	4 (2%)	51 43
1	B	218/270 (81%)	208 (95%)	10 (5%)	24 12
All	All	444/540 (82%)	430 (97%)	14 (3%)	34 22

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

Mol	Chain	Res	Type
1	A	207	GLU
1	A	240	LYS
1	A	322	VAL
1	A	463	MET
1	B	212	ARG
1	B	240	LYS
1	B	243	ASP
1	B	244	LYS
1	B	245	SER
1	B	257	MET
1	B	358	LYS

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Mol	Chain	Res	Type
1	B	461	THR
1	B	466	HIS
1	B	468	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	283	GLN
1	A	410	GLN
1	B	253	ASN
1	B	286	GLN
1	B	437	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 oligosaccharides 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	QMH	A	501	-	23,23,23	1.15	1 (4%)	29,29,29	0.83	1 (3%)

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	QMH	A	501	-	-	4/23/23/23	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	QMH	CAU-CAS	-4.50	1.40	1.50

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	QMH	CAO-OAQ-CAT	-2.77	110.72	117.93

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	QMH	CAN-CAO-OAQ-CAT
2	A	501	QMH	CAL-CAM-CAN-CAO
2	A	501	QMH	N-CA-CB-CG1
2	A	501	QMH	N-CA-CB-CG2

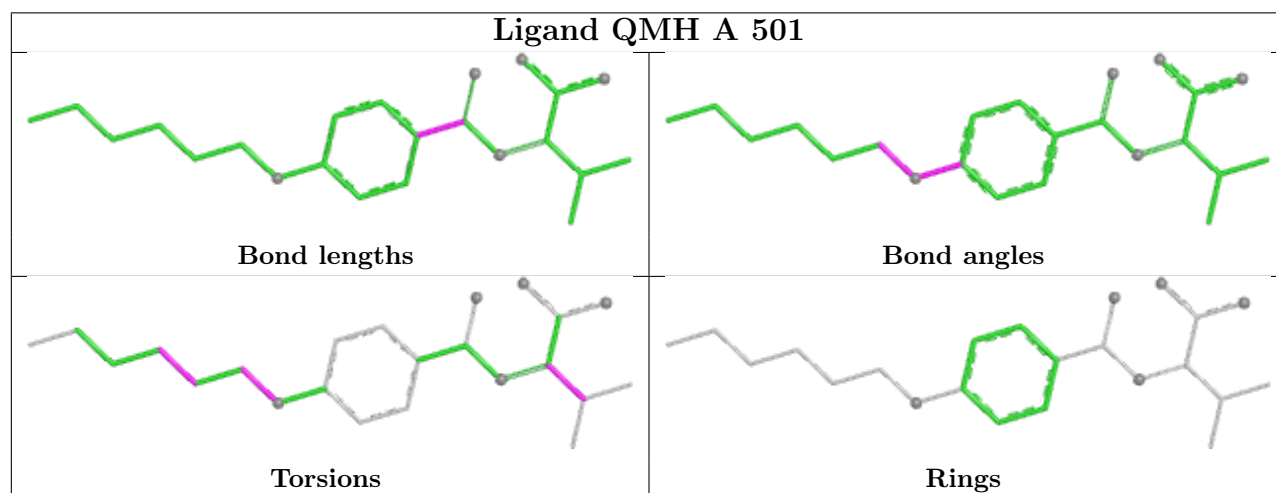
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	QMH	1	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

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, 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	252/300 (84%)	1.05	31 (12%)	8 6	24, 36, 59, 76	1 (0%)
1	B	244/300 (81%)	1.26	50 (20%)	2 2	26, 40, 72, 87	0
All	All	496/600 (82%)	1.15	81 (16%)	4 3	24, 38, 67, 87	1 (0%)

All (81) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	238	THR	8.3
1	A	241	THR	6.8
1	A	239	GLY	5.8
1	B	467	PRO	5.3
1	A	243	ASP	5.3
1	A	240	LYS	4.7
1	A	282	PHE	4.6
1	A	474	LYS	4.6
1	B	363	PHE	4.4
1	B	257	MET	4.4
1	B	240	LYS	4.0
1	B	459	THR	3.9
1	A	242	THR	3.9
1	B	456	ILE	3.8
1	B	241	THR	3.5
1	B	359	PRO	3.5
1	B	454	GLN	3.4
1	B	473	TYR	3.4
1	A	279	ILE	3.4
1	A	363[A]	PHE	3.4
1	B	209	ALA	3.4
1	B	236	ILE	3.2
1	A	464	SER	3.2
1	A	287	PHE	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	277	VAL	3.1
1	B	472	ILE	3.0
1	A	244	LYS	3.0
1	B	452	LEU	3.0
1	B	239	GLY	3.0
1	B	360	PHE	3.0
1	B	455	VAL	2.9
1	B	256	MET	2.9
1	B	469	LEU	2.9
1	A	359	PRO	2.8
1	B	283	GLN	2.8
1	A	277	VAL	2.8
1	A	290	VAL	2.8
1	B	279	ILE	2.7
1	A	465	LEU	2.7
1	B	458	LYS	2.7
1	B	394	SER	2.7
1	A	256	MET	2.6
1	A	463	MET	2.6
1	B	357	ARG	2.6
1	B	247	PHE	2.6
1	A	473	TYR	2.6
1	B	361	GLY	2.5
1	A	360	PHE	2.5
1	A	285	CYS	2.5
1	A	255	LEU	2.5
1	B	468	LEU	2.5
1	B	341	ILE	2.5
1	B	461	THR	2.5
1	B	451	GLN	2.4
1	B	352	PHE	2.4
1	B	466	HIS	2.4
1	B	453	LEU	2.3
1	B	285	CYS	2.3
1	A	333	LEU	2.3
1	B	243	ASP	2.2
1	B	439	MET	2.2
1	A	472	ILE	2.2
1	A	412	ASN	2.2
1	B	395	GLY	2.2
1	B	356	LEU	2.2
1	A	276	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	444	GLN	2.1
1	B	343	GLU	2.1
1	A	278	ALA	2.1
1	A	455	VAL	2.1
1	B	307	VAL	2.1
1	B	412	ASN	2.1
1	A	343	GLU	2.1
1	A	411	ASP	2.1
1	B	255	LEU	2.1
1	B	311	LEU	2.1
1	A	234	ARG	2.0
1	B	427	GLU	2.0
1	B	460	GLU	2.0
1	B	252	MET	2.0
1	B	248	VAL	2.0

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](#)

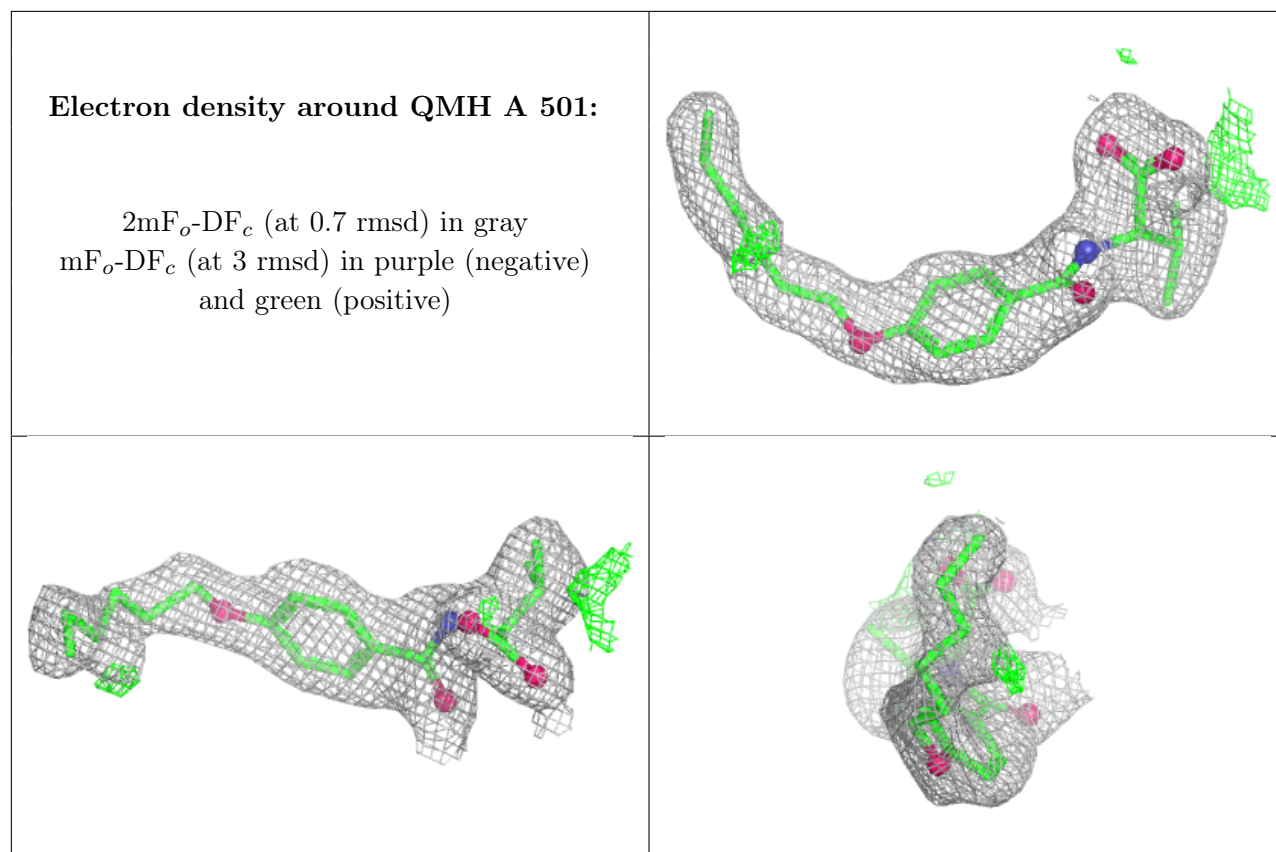
There are no oligosaccharides in this entry.

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($\text{\AA}^2$)	Q<0.9
2	QMH	A	501	23/23	0.86	0.13	37,48,51,53	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.