



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 11:46 AM UTC

PDB ID : 4EMA / pdb_00004ema
Title : Human peroxisome proliferator-activated receptor gamma in complex with
rosiglitazone
Authors : Liberato, M.V.; Nascimento, A.S.; Polikarpov, I.
Deposited on : 2012-04-11
Resolution : 2.54 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

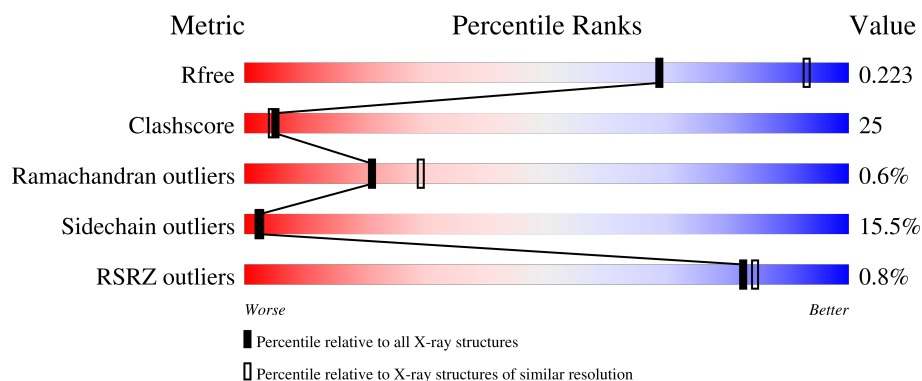
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.54 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	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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	275	60% 27% 6% • 7%
1	B	275	46% 29% 12% • 11%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4102 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.

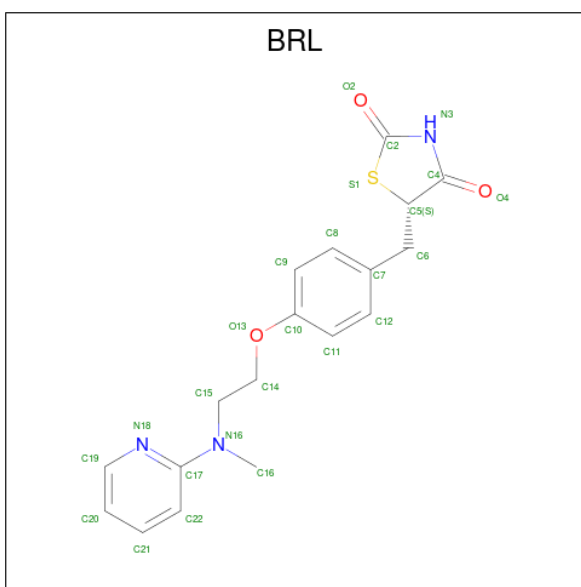
- Molecule 1 is a protein called Peroxisome proliferator-activated receptor gamma.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	256	Total	C	N	O	S	1	0	0
			2049	1322	332	385	10			
1	B	244	Total	C	N	O	S	6	0	0
			1949	1258	317	364	10			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	203	GLY	-	expression tag	UNP P37231
A	204	SER	-	expression tag	UNP P37231
A	205	HIS	-	expression tag	UNP P37231
A	206	MET	-	expression tag	UNP P37231
B	203	GLY	-	expression tag	UNP P37231
B	204	SER	-	expression tag	UNP P37231
B	205	HIS	-	expression tag	UNP P37231
B	206	MET	-	expression tag	UNP P37231

- Molecule 2 is 2,4-THIAZOLIDINEDIONE, 5-[[4-[2-(METHYL-2-PYRIDINYLAMINO)ETHOXY]PHENYL]METHYL]-(9CL) (CCD ID: BRL) (formula: C₁₈H₁₉N₃O₃S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			25	18	3	3	1		

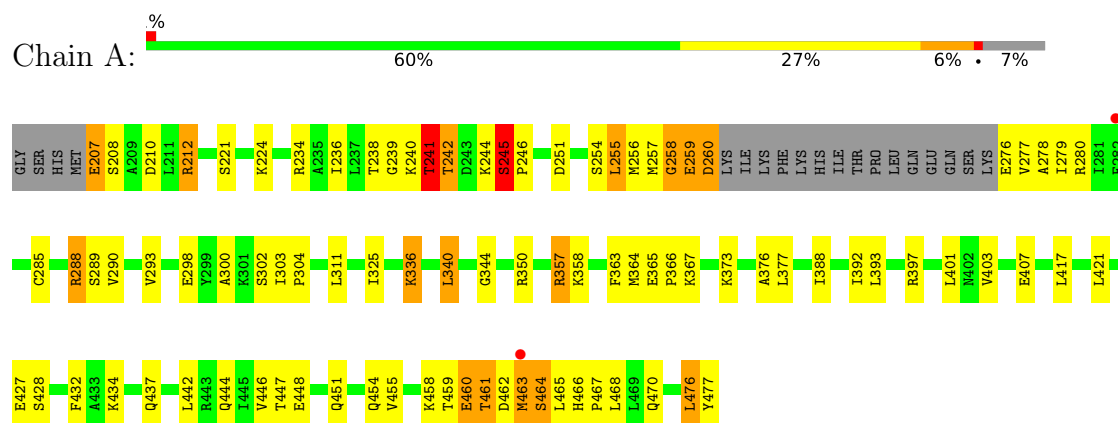
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	43	Total O 43 43	0	0
3	B	36	Total O 36 36	0	0

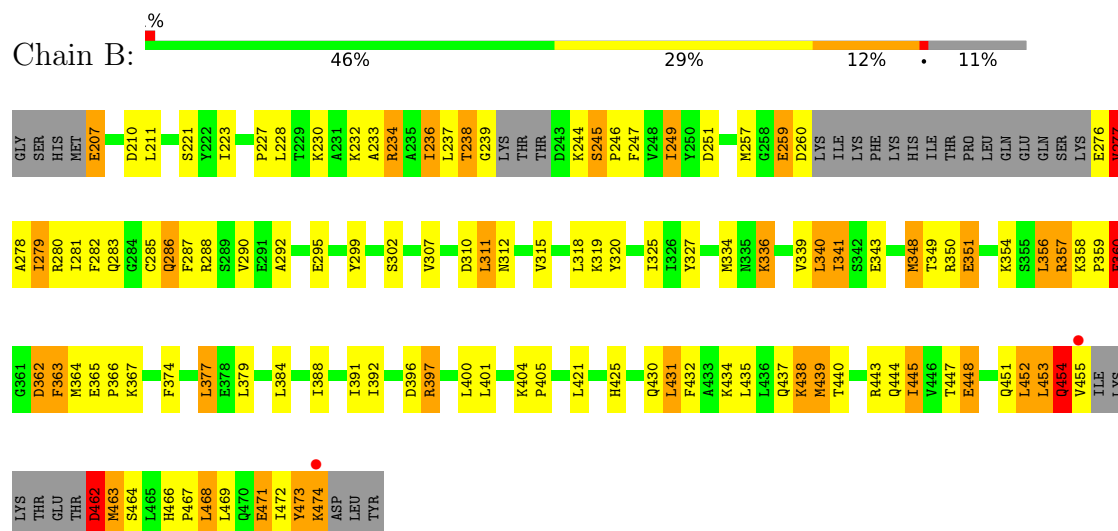
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: Peroxisome proliferator-activated receptor gamma



- 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	90.51Å 61.96Å 117.60Å 90.00° 100.70° 90.00°	Depositor
Resolution (Å)	28.89 – 2.54 28.89 – 2.54	Depositor EDS
% Data completeness (in resolution range)	98.2 (28.89-2.54) 98.1 (28.89-2.54)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.31 (at 2.54Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.180 , 0.225 0.180 , 0.223	Depositor DCC
R_{free} test set	1071 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	48.1	Xtriage
Anisotropy	0.341	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 80.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4102	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.49% 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: BRL

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	0.60	0/2083	1.06	5/2808 (0.2%)
1	B	0.63	0/1980	1.15	19/2666 (0.7%)
All	All	0.62	0/4063	1.10	24/5474 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

There are no bond length outliers.

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	245	SER	CA-C-N	9.33	129.45	120.31
1	A	245	SER	C-N-CA	9.33	129.45	120.31
1	B	237	LEU	N-CA-C	-9.14	102.12	113.18
1	A	241	THR	N-CA-C	8.85	122.53	109.07
1	B	238	THR	N-CA-C	8.28	123.21	112.12
1	A	397	ARG	CA-C-N	7.95	129.21	119.98
1	A	397	ARG	C-N-CA	7.95	129.21	119.98
1	B	207	GLU	CB-CA-C	7.34	124.04	110.10
1	B	360	PHE	N-CA-C	6.31	118.16	111.28
1	B	245	SER	CA-C-N	6.28	126.78	119.99
1	B	245	SER	C-N-CA	6.28	126.78	119.99
1	B	239	GLY	N-CA-C	-6.23	95.23	113.30
1	B	425	HIS	CA-C-N	6.21	126.67	119.47
1	B	425	HIS	C-N-CA	6.21	126.67	119.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	363	PHE	N-CA-C	5.94	117.84	111.36
1	B	236	ILE	CB-CA-C	-5.79	102.88	112.26
1	B	277	VAL	N-CA-C	-5.65	107.38	112.12
1	B	473	TYR	N-CA-C	5.24	117.07	108.52
1	B	249	ILE	CB-CA-C	-5.20	103.96	111.19
1	B	473	TYR	CB-CA-C	5.09	118.54	110.19
1	B	454	GLN	CA-C-N	5.08	130.84	121.70
1	B	454	GLN	C-N-CA	5.08	130.84	121.70
1	B	285	CYS	N-CA-C	-5.05	106.62	112.89
1	B	281	ILE	CB-CA-C	-5.03	105.45	112.04

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	462	ASP	Peptide

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	2049	0	2098	79	0
1	B	1949	0	1988	123	0
2	A	25	0	19	2	0
3	A	43	0	0	0	0
3	B	36	0	0	5	0
All	All	4102	0	4105	200	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 25.

All (200) 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:462:ASP:OD1	1:B:463:MET:HB3	1.33	1.25

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:454:GLN:HB2	1:B:455:VAL:HB	1.27	1.15
1:B:468:LEU:O	1:B:471:GLU:HB2	1.58	1.03
1:B:454:GLN:CB	1:B:455:VAL:HB	1.90	1.00
1:B:463:MET:O	1:B:463:MET:HG3	1.59	0.99
1:A:460:GLU:HB3	1:A:463:MET:HG2	1.46	0.97
1:B:451:GLN:HG3	3:B:535:HOH:O	1.68	0.92
1:B:453:LEU:HD12	1:B:474:LYS:NZ	1.85	0.92
1:A:460:GLU:HB3	1:A:463:MET:CG	2.01	0.89
1:B:466:HIS:HD2	1:B:468:LEU:H	1.20	0.87
1:B:277:VAL:HG23	1:B:280:ARG:HH21	1.42	0.85
1:A:241:THR:OG1	1:A:244:LYS:HB2	1.77	0.84
1:A:357:ARG:CG	1:A:357:ARG:HH11	1.91	0.83
1:B:453:LEU:CD1	1:B:474:LYS:HZ1	1.91	0.82
1:B:228:LEU:HD12	1:B:232:LYS:HD2	1.61	0.81
1:B:278:ALA:HB1	1:B:360:PHE:CD2	2.17	0.80
1:B:454:GLN:HA	1:B:455:VAL:HG23	1.61	0.80
1:B:466:HIS:CD2	1:B:468:LEU:H	1.99	0.79
1:B:360:PHE:C	1:B:362:ASP:H	1.90	0.79
1:B:453:LEU:CD1	1:B:474:LYS:NZ	2.45	0.78
1:A:444:GLN:O	1:A:448:GLU:HG3	1.83	0.77
1:B:350:ARG:O	1:B:354:LYS:HG3	1.84	0.77
1:B:325:ILE:HD11	1:B:392:ILE:HG13	1.64	0.77
1:A:357:ARG:HH11	1:A:357:ARG:HG3	1.49	0.76
1:B:462:ASP:CG	1:B:463:MET:HB3	2.11	0.75
1:B:278:ALA:HB1	1:B:360:PHE:HD2	1.52	0.74
1:B:466:HIS:HD2	1:B:468:LEU:N	1.86	0.74
1:B:454:GLN:HA	1:B:455:VAL:CG2	2.18	0.73
1:B:430:GLN:O	1:B:434:LYS:HG3	1.90	0.72
1:B:364:MET:HA	1:B:364:MET:HE2	1.71	0.72
1:A:276:GLU:HG3	1:A:280:ARG:NH1	2.05	0.72
1:B:453:LEU:HD12	1:B:474:LYS:HZ3	1.55	0.72
1:B:454:GLN:CA	1:B:455:VAL:HB	2.20	0.71
1:B:462:ASP:OD1	1:B:462:ASP:C	2.34	0.71
1:B:444:GLN:O	1:B:448:GLU:HG2	1.91	0.70
1:B:207:GLU:O	1:B:211:LEU:HG	1.93	0.69
1:A:212:ARG:HG3	1:A:212:ARG:HH11	1.58	0.69
1:A:241:THR:HG21	1:A:244:LYS:CE	2.23	0.69
1:A:403:VAL:CG1	1:A:407:GLU:OE2	2.41	0.69
1:A:455:VAL:O	1:A:459:THR:HB	1.95	0.67
1:A:241:THR:CB	1:A:244:LYS:HB2	2.26	0.66
1:B:391:ILE:HG12	1:B:439:MET:HE3	1.77	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:454:GLN:HB2	1:B:455:VAL:CB	2.17	0.65
1:A:461:THR:O	1:A:462:ASP:OD1	2.15	0.65
1:B:360:PHE:O	1:B:362:ASP:N	2.30	0.64
1:B:360:PHE:N	1:B:360:PHE:HD1	1.95	0.64
1:A:212:ARG:HG3	1:A:212:ARG:NH1	2.12	0.63
1:B:364:MET:HE2	1:B:367:LYS:HG3	1.81	0.63
1:B:360:PHE:N	1:B:360:PHE:CD1	2.65	0.62
1:A:300:ALA:HA	1:A:303:ILE:HD12	1.81	0.61
1:A:277:VAL:HG23	1:A:278:ALA:H	1.64	0.61
1:A:255:LEU:HD12	1:A:255:LEU:O	2.00	0.61
1:B:299:TYR:HE2	1:B:388:ILE:HG22	1.66	0.61
1:B:360:PHE:C	1:B:362:ASP:N	2.56	0.60
1:B:421:LEU:HD22	1:B:431:LEU:HD13	1.83	0.60
1:A:241:THR:OG1	1:A:244:LYS:HD2	2.02	0.60
1:A:460:GLU:HB3	1:A:463:MET:HG3	1.83	0.60
1:B:364:MET:HE1	1:B:367:LYS:HD2	1.85	0.59
1:B:454:GLN:CA	1:B:455:VAL:CB	2.81	0.59
1:B:327:TYR:CE1	1:B:367:LYS:HE3	2.38	0.59
1:B:453:LEU:HD12	1:B:474:LYS:HZ1	1.49	0.58
1:A:373:LYS:NZ	1:B:396:ASP:HB3	2.19	0.57
1:A:276:GLU:CG	1:A:280:ARG:NH1	2.67	0.57
1:B:334:MET:HG3	1:B:339:VAL:HG23	1.87	0.57
1:B:360:PHE:O	1:B:363:PHE:HD2	1.87	0.57
1:B:341:ILE:CD1	1:B:348:MET:HE2	2.35	0.56
1:A:357:ARG:NH2	1:A:460:GLU:OE2	2.29	0.56
1:B:454:GLN:HA	1:B:455:VAL:CB	2.35	0.56
1:A:476:LEU:O	1:A:477:TYR:HB2	2.05	0.56
1:B:464:SER:C	3:B:531:HOH:O	2.49	0.56
1:B:249:ILE:O	1:B:349:THR:HG23	2.06	0.55
1:B:290:VAL:HG21	1:B:473:TYR:CD1	2.41	0.55
1:A:239:GLY:C	1:A:241:THR:H	2.13	0.55
1:A:241:THR:HG21	1:A:244:LYS:CD	2.37	0.55
1:A:241:THR:HG1	1:A:244:LYS:HB2	1.72	0.55
1:B:310:ASP:OD1	1:B:312:ASN:ND2	2.36	0.55
1:B:359:PRO:O	1:B:362:ASP:HB2	2.07	0.55
1:B:236:ILE:CG2	1:B:246:PRO:HG3	2.37	0.54
1:A:238:THR:OG1	1:A:240:LYS:HE3	2.07	0.54
1:B:341:ILE:HD12	1:B:348:MET:CE	2.37	0.54
1:B:365:GLU:HB3	1:B:366:PRO:HD3	1.89	0.54
1:B:278:ALA:CB	1:B:360:PHE:CD2	2.89	0.54
1:B:341:ILE:HD12	1:B:348:MET:HE2	1.88	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:276:GLU:HG3	1:B:276:GLU:O	2.06	0.53
1:A:403:VAL:HG12	1:A:407:GLU:OE2	2.09	0.53
1:A:236:ILE:CG2	1:A:246:PRO:HG3	2.40	0.52
1:A:340:LEU:HG	1:A:344:GLY:HA2	1.91	0.52
1:B:327:TYR:CZ	1:B:367:LYS:HE3	2.44	0.52
1:A:357:ARG:HG3	1:A:357:ARG:NH1	2.22	0.52
1:A:434:LYS:HA	1:A:437:GLN:HE21	1.76	0.51
1:B:466:HIS:CD2	1:B:466:HIS:C	2.89	0.50
1:A:367:LYS:H	1:A:367:LYS:CD	2.25	0.50
1:A:277:VAL:HG23	1:A:278:ALA:N	2.26	0.50
1:A:373:LYS:O	1:A:376:ALA:HB3	2.11	0.50
1:B:462:ASP:OD1	1:B:462:ASP:O	2.30	0.50
1:A:259:GLU:O	1:A:260:ASP:O	2.29	0.49
1:A:357:ARG:HH11	1:A:357:ARG:HG2	1.74	0.49
1:A:241:THR:HB	1:A:244:LYS:HB2	1.94	0.49
1:B:360:PHE:O	1:B:363:PHE:CD2	2.64	0.49
1:B:443:ARG:O	1:B:447:THR:HG23	2.12	0.49
1:B:259:GLU:HG2	1:B:259:GLU:O	2.13	0.49
1:A:357:ARG:CG	1:A:357:ARG:NH1	2.61	0.48
1:B:290:VAL:HG21	1:B:473:TYR:CE1	2.48	0.48
1:A:357:ARG:HG2	1:A:357:ARG:O	2.12	0.48
1:A:288:ARG:HA	1:A:288:ARG:HD2	1.62	0.48
1:B:236:ILE:HG22	1:B:236:ILE:O	2.12	0.48
1:A:239:GLY:C	1:A:241:THR:N	2.72	0.48
1:A:373:LYS:HZ3	1:B:396:ASP:HB3	1.78	0.48
1:A:224:LYS:HE2	1:A:298:GLU:OE2	2.14	0.48
1:B:260:ASP:HB3	3:B:526:HOH:O	2.12	0.48
1:B:357:ARG:HE	1:B:357:ARG:HB2	1.46	0.47
1:A:302:SER:O	1:A:304:PRO:HD3	2.13	0.47
1:B:452:LEU:C	1:B:453:LEU:HD23	2.39	0.47
1:B:276:GLU:O	1:B:276:GLU:CG	2.62	0.47
1:B:432:PHE:HB3	3:B:501:HOH:O	2.15	0.47
1:A:212:ARG:NE	1:A:212:ARG:HA	2.29	0.47
1:B:260:ASP:OD1	1:B:260:ASP:C	2.58	0.47
1:B:307:VAL:HG12	1:B:307:VAL:O	2.13	0.47
1:A:212:ARG:HH11	1:A:212:ARG:CG	2.26	0.47
1:A:365:GLU:N	1:A:366:PRO:CD	2.78	0.47
1:B:299:TYR:CE2	1:B:388:ILE:HG22	2.48	0.47
1:B:282:PHE:C	1:B:282:PHE:CD1	2.92	0.47
1:A:285:CYS:SG	1:A:363:PHE:HZ	2.37	0.47
1:A:421:LEU:CD1	1:A:432:PHE:HA	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:454:GLN:O	1:A:458:LYS:HG3	2.14	0.46
1:B:292:ALA:HA	1:B:295:GLU:HG3	1.97	0.46
1:A:325:ILE:HG12	1:A:388:ILE:HB	1.96	0.46
1:A:290:VAL:HG21	1:A:466:HIS:CD2	2.51	0.46
1:B:207:GLU:HB3	1:B:210:ASP:OD2	2.15	0.46
1:A:325:ILE:HD11	1:A:392:ILE:HG13	1.96	0.46
1:A:336:LYS:NZ	1:A:350:ARG:HH12	2.14	0.46
1:B:318:LEU:O	1:B:319:LYS:C	2.58	0.46
1:A:364:MET:SD	2:A:601:BRL:H11	2.56	0.46
1:B:454:GLN:HA	1:B:455:VAL:HB	1.93	0.46
1:A:476:LEU:O	1:A:477:TYR:CB	2.65	0.45
1:B:374:PHE:O	1:B:377:LEU:HB2	2.15	0.45
1:B:236:ILE:C	1:B:238:THR:N	2.74	0.45
1:B:282:PHE:CD1	1:B:463:MET:HE1	2.52	0.45
1:B:251:ASP:OD1	1:B:251:ASP:C	2.59	0.45
1:A:245:SER:HA	1:A:246:PRO:HD3	1.58	0.45
1:A:241:THR:HG21	1:A:244:LYS:HE3	1.99	0.45
1:B:466:HIS:CG	1:B:467:PRO:HD2	2.53	0.44
1:A:293:VAL:HG11	1:A:468:LEU:HD21	1.98	0.44
1:B:340:LEU:HD12	1:B:340:LEU:HA	1.83	0.44
1:B:453:LEU:HD23	1:B:453:LEU:N	2.31	0.44
1:B:469:LEU:HA	1:B:472:ILE:HD12	1.99	0.44
1:A:207:GLU:O	1:A:210:ASP:HB2	2.18	0.44
1:A:254:SER:O	1:A:258:GLY:N	2.50	0.44
1:A:365:GLU:N	1:A:366:PRO:HD2	2.33	0.44
1:B:334:MET:CG	1:B:339:VAL:HG23	2.47	0.44
1:B:286:GLN:O	1:B:287:PHE:C	2.58	0.44
1:A:242:THR:C	1:A:244:LYS:H	2.26	0.43
1:A:462:ASP:O	1:A:464:SER:N	2.51	0.43
1:B:365:GLU:N	1:B:366:PRO:CD	2.81	0.43
1:B:236:ILE:C	1:B:238:THR:H	2.22	0.43
1:A:255:LEU:HD12	1:A:255:LEU:C	2.43	0.43
1:B:364:MET:CE	1:B:367:LYS:HE2	2.48	0.43
1:A:462:ASP:OD1	1:A:462:ASP:C	2.60	0.43
1:A:417:LEU:HD23	1:A:432:PHE:HE1	1.82	0.43
1:B:404:LYS:HB3	1:B:405:PRO:HD3	2.01	0.43
1:B:363:PHE:CE1	1:B:364:MET:HG2	2.54	0.43
1:B:351:GLU:OE1	1:B:354:LYS:HD3	2.18	0.43
1:B:374:PHE:HD1	1:B:438:LYS:HG2	1.84	0.43
1:A:259:GLU:O	1:A:260:ASP:C	2.62	0.42
1:A:276:GLU:HG2	1:A:280:ARG:HH12	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:444:GLN:O	1:B:448:GLU:CG	2.65	0.42
1:B:471:GLU:C	1:B:473:TYR:N	2.77	0.42
1:A:367:LYS:H	1:A:367:LYS:HD2	1.84	0.42
1:A:467:PRO:O	1:A:470:GLN:HB2	2.19	0.42
1:B:356:LEU:O	1:B:357:ARG:C	2.63	0.42
1:A:466:HIS:O	1:A:467:PRO:C	2.62	0.42
1:B:279:ILE:HD12	1:B:463:MET:HA	2.01	0.42
1:B:379:LEU:HD11	1:B:435:LEU:HD13	2.00	0.42
1:B:452:LEU:HD23	1:B:452:LEU:HA	1.57	0.42
1:B:286:GLN:C	1:B:288:ARG:N	2.74	0.42
1:B:336:LYS:O	1:B:350:ARG:HD2	2.19	0.42
1:B:365:GLU:N	1:B:366:PRO:HD2	2.34	0.42
1:B:448:GLU:HG2	1:B:448:GLU:H	1.57	0.42
1:B:227:PRO:O	1:B:232:LYS:CE	2.67	0.42
1:B:311:LEU:HD23	1:B:315:VAL:HG23	2.00	0.42
1:A:357:ARG:O	1:A:358:LYS:C	2.63	0.41
1:B:283:GLN:O	1:B:286:GLN:HB2	2.20	0.41
1:A:289:SER:OG	2:A:601:BRL:H5	2.20	0.41
1:B:232:LYS:O	1:B:233:ALA:C	2.63	0.41
1:A:392:ILE:HG22	1:A:393:LEU:HD22	2.01	0.41
1:B:468:LEU:HA	1:B:468:LEU:HD12	1.76	0.41
1:A:251:ASP:OD1	1:A:251:ASP:C	2.63	0.41
1:B:230:LYS:O	1:B:234:ARG:HB2	2.20	0.41
1:A:276:GLU:CG	1:A:280:ARG:HH12	2.32	0.41
1:B:445:ILE:HD13	1:B:445:ILE:HA	1.84	0.41
1:B:286:GLN:OE1	1:B:286:GLN:CA	2.68	0.41
1:B:320:TYR:HB2	1:B:397:ARG:CG	2.51	0.41
1:B:430:GLN:NE2	3:B:501:HOH:O	2.54	0.41
1:B:437:GLN:O	1:B:440:THR:HB	2.21	0.41
1:B:466:HIS:CD2	1:B:467:PRO:HD2	2.56	0.40
1:A:277:VAL:C	1:A:279:ILE:N	2.79	0.40
1:B:247:PHE:C	1:B:247:PHE:CD1	2.99	0.40
1:B:358:LYS:HG3	1:B:359:PRO:HA	2.02	0.40
1:B:223:ILE:HD13	1:B:223:ILE:HA	1.86	0.40
1:B:466:HIS:O	1:B:467:PRO:C	2.65	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	252/275 (92%)	238 (94%)	12 (5%)	2 (1%)	16	22
1	B	236/275 (86%)	218 (92%)	17 (7%)	1 (0%)	30	39
All	All	488/550 (89%)	456 (93%)	29 (6%)	3 (1%)	21	29

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	336	LYS
1	A	463	MET
1	A	258	GLY

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	229/247 (93%)	198 (86%)	31 (14%)	4	3
1	B	217/247 (88%)	179 (82%)	38 (18%)	2	1
All	All	446/494 (90%)	377 (84%)	69 (16%)	2	2

All (69) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	207	GLU
1	A	208	SER
1	A	212	ARG

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Mol	Chain	Res	Type
1	A	221	SER
1	A	234	ARG
1	A	241	THR
1	A	242	THR
1	A	245	SER
1	A	255	LEU
1	A	256	MET
1	A	257	MET
1	A	259	GLU
1	A	260	ASP
1	A	288	ARG
1	A	311	LEU
1	A	336	LYS
1	A	340	LEU
1	A	357	ARG
1	A	377	LEU
1	A	401	LEU
1	A	427	GLU
1	A	428	SER
1	A	442	LEU
1	A	446	VAL
1	A	447	THR
1	A	451	GLN
1	A	460	GLU
1	A	461	THR
1	A	464	SER
1	A	465	LEU
1	A	476	LEU
1	B	221	SER
1	B	234	ARG
1	B	244	LYS
1	B	245	SER
1	B	257	MET
1	B	259	GLU
1	B	277	VAL
1	B	279	ILE
1	B	286	GLN
1	B	302	SER
1	B	311	LEU
1	B	340	LEU
1	B	341	ILE
1	B	343	GLU

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Mol	Chain	Res	Type
1	B	348	MET
1	B	351	GLU
1	B	356	LEU
1	B	357	ARG
1	B	360	PHE
1	B	362	ASP
1	B	377	LEU
1	B	384	LEU
1	B	397	ARG
1	B	400	LEU
1	B	401	LEU
1	B	431	LEU
1	B	438	LYS
1	B	439	MET
1	B	445	ILE
1	B	448	GLU
1	B	452	LEU
1	B	453	LEU
1	B	454	GLN
1	B	462	ASP
1	B	463	MET
1	B	468	LEU
1	B	471	GLU
1	B	474	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	286	GLN
1	A	294	GLN
1	A	308	ASN
1	A	314	GLN
1	A	345	GLN
1	A	415	GLN
1	A	454	GLN
1	A	470	GLN
1	B	294	GLN
1	B	314	GLN
1	B	345	GLN
1	B	410	GLN
1	B	412	ASN
1	B	415	GLN

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Mol	Chain	Res	Type
1	B	430	GLN
1	B	444	GLN
1	B	454	GLN
1	B	466	HIS

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	BRL	A	601	-	27,27,27	2.85	6 (22%)	36,36,36	2.69	10 (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	BRL	A	601	-	-	2/14/26/26	0/3/3/3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	BRL	C2-S1	-12.83	1.65	1.76
2	A	601	BRL	C4-N3	-3.58	1.33	1.37
2	A	601	BRL	C2-N3	-3.47	1.32	1.36
2	A	601	BRL	C17-N16	-2.15	1.31	1.37
2	A	601	BRL	O13-C10	-2.09	1.32	1.37
2	A	601	BRL	O13-C14	-2.07	1.36	1.43

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	BRL	C4-C5-S1	-8.37	98.93	105.93
2	A	601	BRL	C5-S1-C2	7.05	97.21	92.77
2	A	601	BRL	C5-C4-N3	6.22	117.25	111.99
2	A	601	BRL	C4-N3-C2	-4.39	115.27	118.20
2	A	601	BRL	C19-N18-C17	3.82	122.37	116.81
2	A	601	BRL	C16-N16-C15	3.08	123.73	115.56
2	A	601	BRL	O4-C4-N3	-2.96	121.44	124.93
2	A	601	BRL	C14-O13-C10	-2.85	110.53	117.93
2	A	601	BRL	N18-C17-N16	2.66	120.83	116.62
2	A	601	BRL	C15-N16-C17	-2.42	117.24	120.71

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	601	BRL	C9-C10-O13-C14
2	A	601	BRL	C11-C10-O13-C14

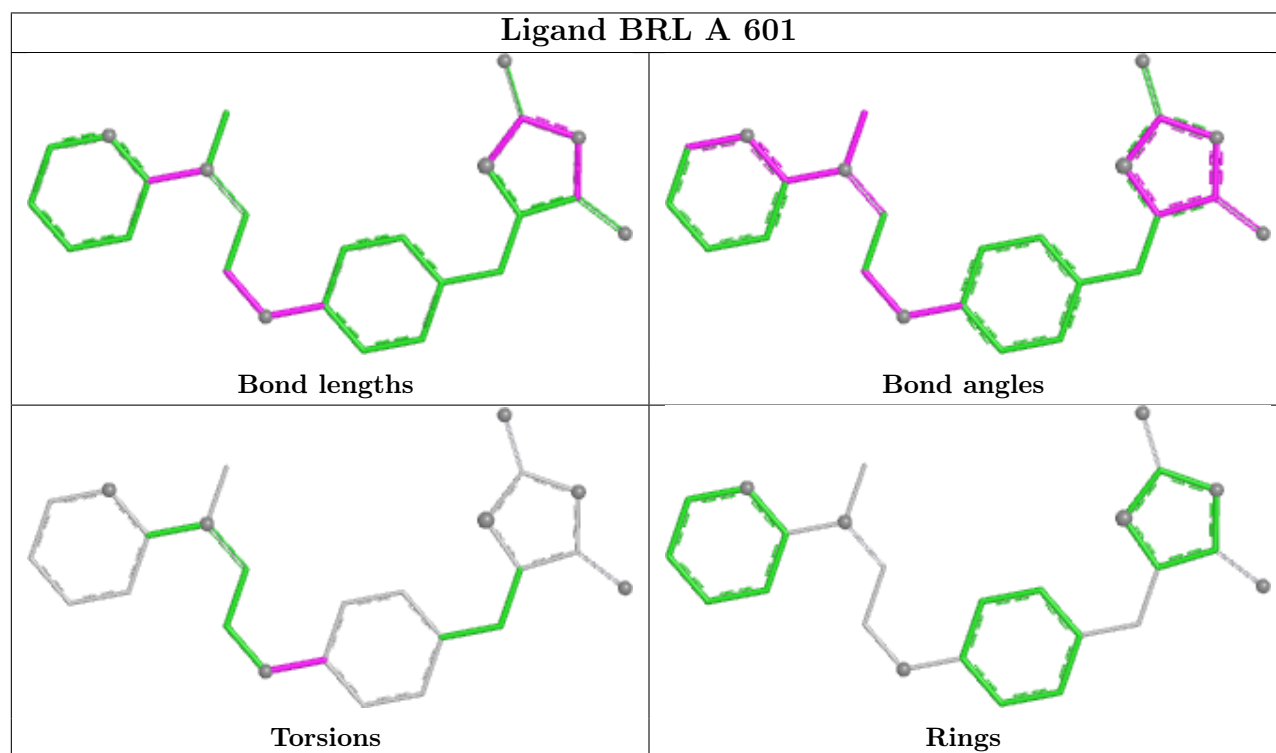
There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	BRL	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	256/275 (93%)	-0.38	2 (0%) 82 84	31, 53, 89, 148	6 (2%)
1	B	244/275 (88%)	-0.27	2 (0%) 82 84	29, 57, 100, 134	11 (4%)
All	All	500/550 (90%)	-0.33	4 (0%) 82 84	29, 54, 97, 148	17 (3%)

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	455	VAL	3.0
1	A	463	MET	2.4
1	A	282	PHE	2.4
1	B	474	LYS	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](#)

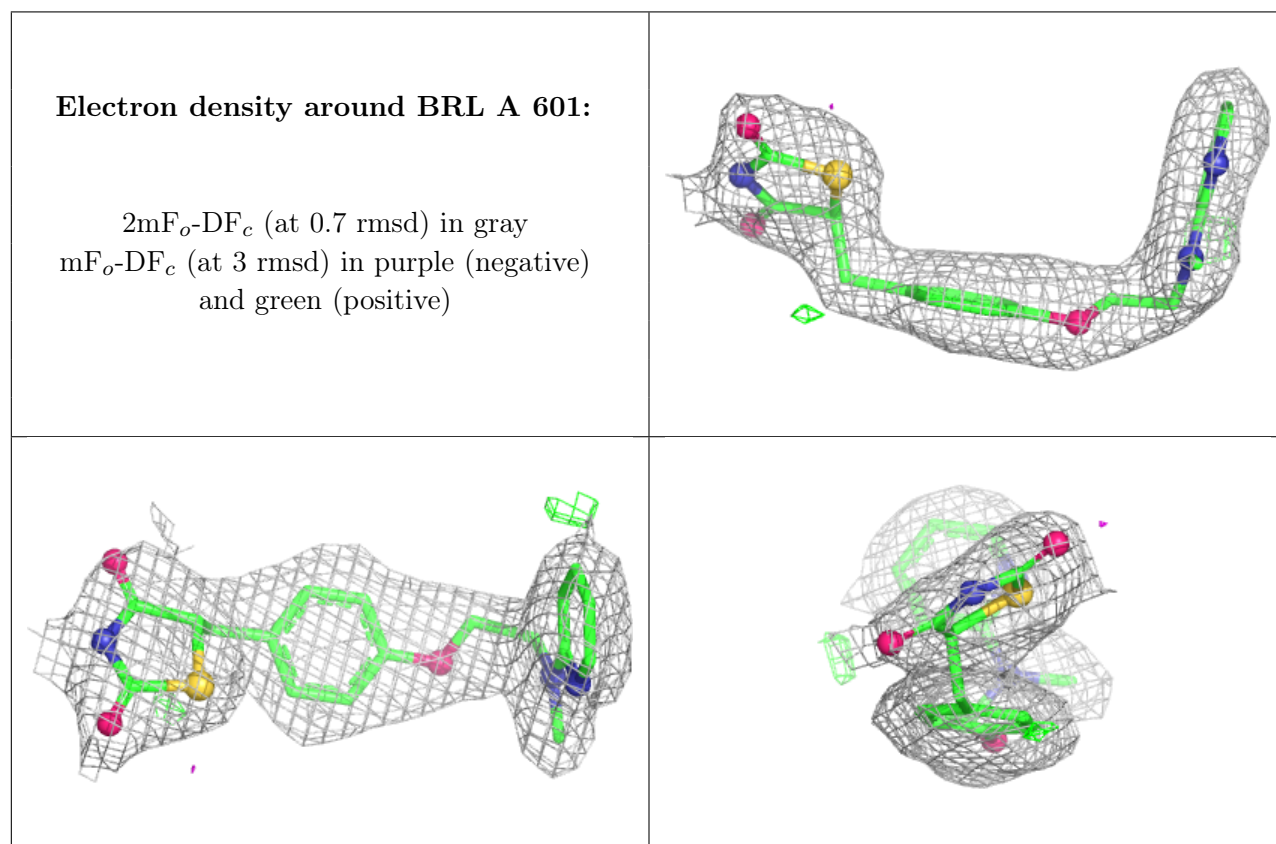
There are no oligosaccharides in this entry.

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(Å ²)	Q<0.9
2	BRL	A	601	25/25	0.93	0.09	52,64,84,85	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.