



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 15, 2026 – 01:07 AM UTC

PDB ID : 7BY6 / pdb_00007by6
Title : Plasmodium vivax cytoplasmic Phenylalanyl-tRNA synthetase in complex with BRD1389
Authors : Malhotra, N.; Manmohan, S.; Harlos, K.; Melillo, B.; Schreiber, S.L.; Manickam, Y.; Sharma, S.
Deposited on : 2020-04-21
Resolution : 3.00 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

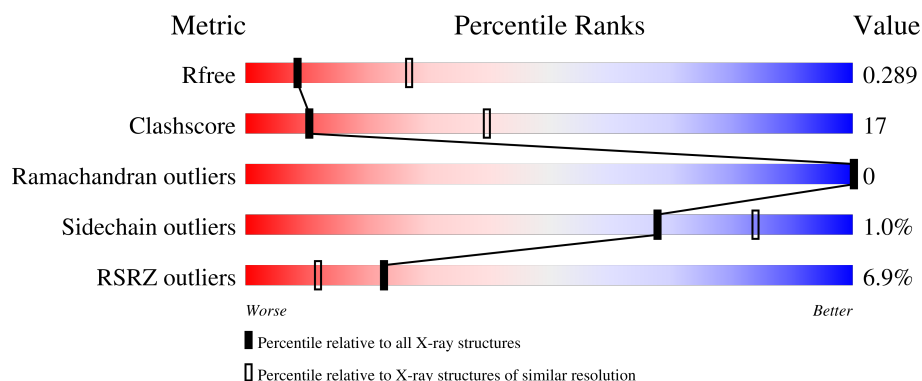
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	299	4% 72% 27%
2	B	618	8% 65% 31% ..

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6924 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 Phenylalanyl-tRNA synthetase beta chain, putative.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	299	Total	C	N	O	S	0	0	0
			2365	1521	405	428	11			

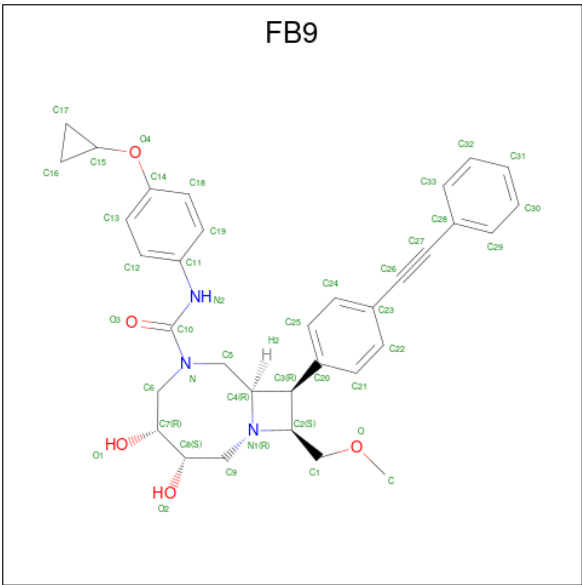
- Molecule 2 is a protein called Phenylalanyl-tRNA synthetase beta chain, putative.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	606	Total	C	N	O	S	0	1	0
			4516	2870	750	871	25			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	MET	-	expression tag	UNP A5K464
B	1	GLY	-	expression tag	UNP A5K464

- Molecule 3 is (3S,4R,8R,9R,10S)-N-(4-cyclopropyloxyphenyl)-10-(methoxymethyl)-3,4-bis(o xidanyl)-9-[4-(2-phenylethynyl)phenyl]-1,6-diazabicyclo[6.2.0]decane-6-carboxamide (CCD ID: FB9) (formula: C₃₄H₃₇N₃O₅) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			42	34	3	5		

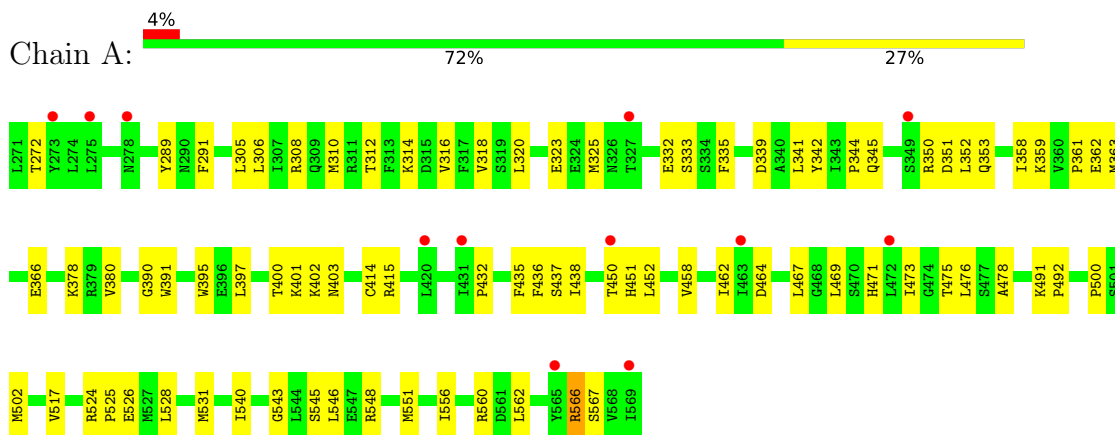
- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Mg	0	0
			1	1		

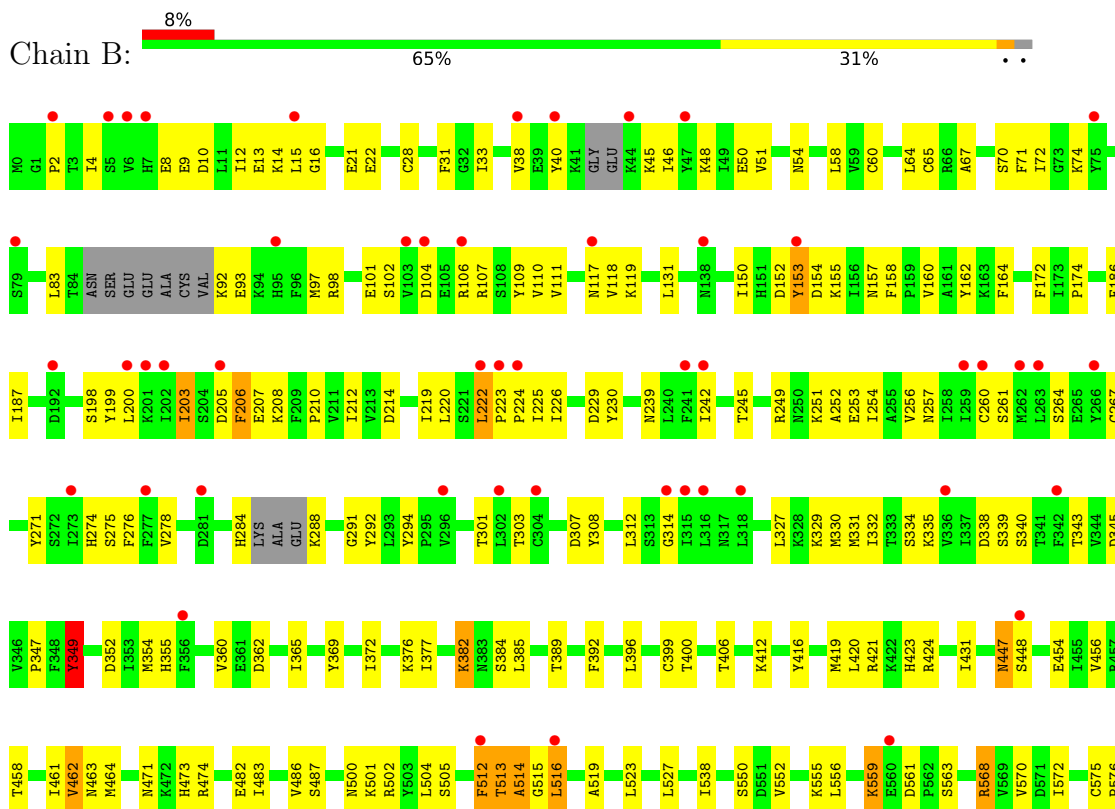
3 Residue-property plots [i](#)

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: Phenylalanyl-tRNA synthetase beta chain, putative



- Molecule 2: Phenylalanyl-tRNA synthetase beta chain, putative



F581	I586	V591	F595	V603	I604	I611	M612	L615	M616	M617
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4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	136.52Å 74.07Å 121.68Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	121.78 – 3.00 121.68 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.5 (121.78-3.00) 99.6 (121.68-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.90 (at 3.01Å)	Xtriage
Refinement program	PHENIX 1.15rc1_3423	Depositor
R, R_{free}	0.214 , 0.288 0.225 , 0.289	Depositor DCC
R_{free} test set	1297 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å ²)	75.2	Xtriage
Anisotropy	0.683	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 93.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6924	wwPDB-VP
Average B, all atoms (Å ²)	91.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.00% 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: FB9, MG

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	0.62	6/2430 (0.2%)	0.71	2/3289 (0.1%)
2	B	0.82	22/4600 (0.5%)	0.90	25/6254 (0.4%)
All	All	0.76	28/7030 (0.4%)	0.84	27/9543 (0.3%)

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	568	ARG	C-O	-7.75	1.15	1.24
2	B	198	SER	CA-C	-7.16	1.44	1.52
2	B	448	SER	C-O	-7.14	1.15	1.23
2	B	448	SER	CA-C	-7.10	1.43	1.52
2	B	153	TYR	CA-C	-6.77	1.44	1.52
2	B	14	LYS	CA-C	-6.67	1.43	1.52
1	A	351	ASP	C-O	-6.49	1.16	1.24
2	B	200	LEU	CA-C	-6.39	1.45	1.52
2	B	514	ALA	N-CA	-6.36	1.37	1.46
2	B	154	ASP	CA-C	-6.17	1.44	1.52
1	A	353	GLN	N-CA	-5.91	1.39	1.46
2	B	382	LYS	CA-C	-5.90	1.45	1.52
2	B	198	SER	C-O	-5.88	1.17	1.24
2	B	104	ASP	CA-C	-5.83	1.45	1.53
2	B	200	LEU	N-CA	-5.82	1.39	1.46
1	A	351	ASP	CA-C	-5.78	1.45	1.52
1	A	352	LEU	C-O	-5.65	1.17	1.24
2	B	515	GLY	C-O	-5.65	1.16	1.23
1	A	352	LEU	CA-C	-5.54	1.45	1.52
2	B	152	ASP	C-O	-5.50	1.17	1.23
2	B	226	ILE	N-CA	-5.44	1.39	1.46
2	B	205	ASP	C-O	-5.33	1.18	1.24
2	B	203	ILE	C-O	-5.32	1.18	1.24
2	B	102	SER	CA-C	-5.25	1.46	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	200	LEU	C-O	-5.23	1.18	1.24
2	B	512	PHE	CA-C	-5.17	1.45	1.52
2	B	559	LYS	C-O	-5.07	1.17	1.23
1	A	566	ARG	CA-C	-5.02	1.48	1.53

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	200	LEU	N-CA-C	10.87	122.82	110.97
2	B	222	LEU	CA-C-N	7.30	127.90	120.38
2	B	222	LEU	C-N-CA	7.30	127.90	120.38
1	A	352	LEU	N-CA-C	6.81	119.57	111.33
2	B	9	GLU	CA-C-N	6.81	129.74	120.54
2	B	9	GLU	C-N-CA	6.81	129.74	120.54
2	B	349	TYR	CB-CA-C	-6.72	100.60	110.96
2	B	514	ALA	N-CA-C	-6.64	98.43	109.46
2	B	14	LYS	CA-C-N	-6.15	112.07	120.44
2	B	14	LYS	C-N-CA	-6.15	112.07	120.44
1	A	450	THR	N-CA-C	-6.01	104.64	111.07
2	B	154	ASP	N-CA-C	6.00	119.86	112.54
2	B	512	PHE	N-CA-C	5.81	120.53	113.38
2	B	447	ASN	N-CA-C	5.72	122.99	110.80
2	B	516	LEU	N-CA-C	-5.68	106.20	113.01
2	B	104	ASP	N-CA-C	5.67	118.54	109.02
2	B	104	ASP	CB-CA-C	-5.65	108.43	115.89
2	B	225	ILE	N-CA-C	5.62	121.04	109.34
2	B	101	GLU	CA-C-N	5.62	127.81	120.28
2	B	101	GLU	C-N-CA	5.62	127.81	120.28
2	B	9	GLU	O-C-N	5.61	130.05	122.59
2	B	104	ASP	CA-C-N	-5.53	113.28	122.08
2	B	104	ASP	C-N-CA	-5.53	113.28	122.08
2	B	12	ILE	CA-C-N	5.18	127.65	120.29
2	B	12	ILE	C-N-CA	5.18	127.65	120.29
2	B	447	ASN	N-CA-CB	-5.18	101.73	110.49
2	B	153	TYR	CB-CA-C	-5.04	102.97	110.88

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	2365	0	2175	66	0
2	B	4516	0	4087	172	0
3	A	42	0	0	4	0
4	B	1	0	0	0	0
All	All	6924	0	6262	223	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 17.

All (223) 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:289:TYR:HB2	2:B:568:ARG:NH1	1.69	1.06
1:A:289:TYR:HB2	2:B:568:ARG:HH11	1.21	1.02
2:B:222:LEU:HB3	2:B:224:PRO:HD2	1.52	0.91
1:A:464:ASP:HB3	1:A:467:LEU:HD11	1.58	0.85
2:B:118:VAL:HA	2:B:267:CYS:HB2	1.61	0.83
1:A:272:THR:HG21	2:B:513:THR:HG22	1.62	0.81
2:B:389:THR:HG21	2:B:502:ARG:HD2	1.66	0.77
1:A:310:MET:HG3	1:A:438:ILE:HG21	1.65	0.77
2:B:334:SER:O	2:B:335:LYS:HG3	1.87	0.75
2:B:106:ARG:NH1	2:B:206:PHE:HB3	2.04	0.72
2:B:223:PRO:HD2	2:B:224:PRO:CD	2.20	0.72
2:B:464:MET:HE1	2:B:505:SER:OG	1.89	0.72
2:B:327:LEU:HB3	2:B:332:ILE:HB	1.70	0.72
2:B:106:ARG:HH11	2:B:206:PHE:CB	2.05	0.69
2:B:210:PRO:HG2	2:B:223:PRO:CG	2.22	0.69
2:B:257:ASN:O	2:B:261:SER:HB3	1.92	0.69
2:B:303:THR:HG22	2:B:343:THR:HA	1.75	0.68
2:B:419:MET:HE2	2:B:419:MET:HA	1.75	0.68
2:B:60:CYS:HB3	2:B:131:LEU:HD13	1.75	0.67
2:B:419:MET:HE3	2:B:595:PHE:CE1	2.29	0.67
2:B:251:LYS:HA	2:B:254:ILE:HD12	1.76	0.66
2:B:458:THR:HA	2:B:486:VAL:HG13	1.77	0.66
1:A:314:LYS:O	1:A:318:VAL:HG12	1.95	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:8:GLU:C	2:B:10:ASP:H	2.02	0.66
1:A:469:LEU:HD11	1:A:500:PRO:HG2	1.77	0.65
1:A:342:TYR:OH	2:B:354:MET:CE	2.45	0.65
2:B:223:PRO:CD	2:B:224:PRO:CD	2.75	0.65
2:B:106:ARG:NH1	2:B:206:PHE:CB	2.59	0.65
2:B:186:PHE:HE2	2:B:224:PRO:HG2	1.62	0.64
1:A:291:PHE:HE1	2:B:516:LEU:HD11	1.63	0.64
2:B:223:PRO:HB2	2:B:224:PRO:HD3	1.78	0.64
2:B:223:PRO:HD2	2:B:224:PRO:HD2	1.80	0.63
2:B:223:PRO:HD2	2:B:224:PRO:HD3	1.80	0.63
2:B:223:PRO:CB	2:B:224:PRO:HD3	2.29	0.63
2:B:223:PRO:CD	2:B:224:PRO:HD3	2.29	0.62
2:B:97:MET:HA	2:B:160:VAL:O	1.99	0.62
2:B:523:LEU:HD11	2:B:572:ILE:HD13	1.82	0.62
2:B:106:ARG:HH12	2:B:206:PHE:HA	1.64	0.61
2:B:385:LEU:O	2:B:389:THR:HG23	1.99	0.60
1:A:363:MET:HE2	1:A:402:LYS:HG2	1.81	0.60
2:B:40:TYR:HA	2:B:45:LYS:H	1.66	0.60
2:B:106:ARG:HH11	2:B:206:PHE:HB3	1.65	0.60
2:B:307:ASP:OD1	2:B:308:TYR:N	2.35	0.59
2:B:464:MET:HE1	2:B:505:SER:CB	2.32	0.59
2:B:164:PHE:CE1	2:B:210:PRO:HB3	2.38	0.58
2:B:97:MET:HG3	2:B:160:VAL:HG23	1.85	0.58
2:B:516:LEU:HD22	2:B:568:ARG:HD2	1.84	0.58
1:A:397:LEU:O	1:A:400:THR:HG22	2.04	0.58
2:B:362:ASP:HA	2:B:365:ILE:HD12	1.85	0.58
1:A:378:LYS:HG3	1:A:395:TRP:HB3	1.85	0.58
2:B:206:PHE:CD1	2:B:206:PHE:N	2.70	0.58
2:B:83:LEU:HD13	2:B:274:HIS:CD2	2.39	0.58
2:B:523:LEU:CD1	2:B:572:ILE:HD13	2.34	0.58
2:B:275:SER:HB2	2:B:294:TYR:O	2.04	0.57
2:B:284:HIS:HB3	2:B:288:LYS:N	2.19	0.57
2:B:347:PRO:HB2	2:B:349:TYR:HB2	1.86	0.57
2:B:447:ASN:OD1	2:B:447:ASN:N	2.36	0.57
2:B:71:PHE:O	2:B:329:LYS:HG2	2.04	0.57
1:A:272:THR:CG2	2:B:513:THR:HG22	2.33	0.57
1:A:305:LEU:HA	1:A:308:ARG:HD3	1.86	0.57
1:A:525:PRO:HG2	2:B:355:HIS:CE1	2.39	0.57
2:B:8:GLU:HB2	2:B:45:LYS:O	2.04	0.57
2:B:210:PRO:HG2	2:B:223:PRO:HG3	1.86	0.57
2:B:419:MET:HE3	2:B:595:PHE:HE1	1.70	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:432:PRO:HB3	1:A:464:ASP:CG	2.30	0.56
1:A:469:LEU:HD23	1:A:502:MET:HE3	1.87	0.56
1:A:291:PHE:CE1	2:B:516:LEU:HD11	2.40	0.56
2:B:150:ILE:HA	2:B:242:ILE:HD13	1.88	0.56
2:B:347:PRO:C	2:B:349:TYR:N	2.61	0.56
1:A:342:TYR:OH	2:B:354:MET:HE2	2.06	0.55
2:B:223:PRO:N	2:B:224:PRO:CD	2.69	0.55
1:A:306:LEU:HG	1:A:546:LEU:HD23	1.86	0.55
2:B:106:ARG:NH1	2:B:206:PHE:HA	2.20	0.55
2:B:8:GLU:C	2:B:10:ASP:N	2.59	0.55
2:B:21:GLU:HB2	2:B:22:GLU:OE2	2.06	0.55
1:A:469:LEU:O	1:A:473:ILE:HG13	2.07	0.54
2:B:212:ILE:HD12	2:B:220:LEU:HB2	1.89	0.54
2:B:556:LEU:HB3	2:B:570:VAL:HG11	1.89	0.54
1:A:325:MET:HE3	1:A:437:SER:HB3	1.87	0.54
1:A:478:ALA:HB2	2:B:377:ILE:CG2	2.37	0.54
3:A:601:FB9:C5	3:A:601:FB9:C21	2.85	0.54
1:A:528:LEU:O	1:A:531:MET:HB2	2.08	0.53
2:B:406:THR:HG21	2:B:463:ASN:HB3	1.89	0.53
1:A:462:ILE:HD12	1:A:540:ILE:HD11	1.89	0.53
2:B:347:PRO:C	2:B:349:TYR:H	2.15	0.53
2:B:8:GLU:HG3	2:B:45:LYS:HG2	1.90	0.53
2:B:471:ASN:HB3	2:B:474[B]:ARG:HG3	1.90	0.53
2:B:301:THR:HA	2:B:345:ASP:HA	1.90	0.53
1:A:344:PRO:HG3	1:A:391:TRP:CE2	2.43	0.53
2:B:4:ILE:O	2:B:48:LYS:HA	2.08	0.53
1:A:366:GLU:HG2	1:A:401:LYS:NZ	2.24	0.53
1:A:390:GLY:O	2:B:354:MET:HE3	2.09	0.53
2:B:155:LYS:HD3	2:B:230:TYR:CZ	2.44	0.53
2:B:212:ILE:CD1	2:B:220:LEU:HB2	2.38	0.53
3:A:601:FB9:C21	3:A:601:FB9:C1	2.85	0.52
1:A:316:VAL:O	1:A:320:LEU:HD12	2.10	0.52
1:A:341:LEU:HD21	1:A:414:CYS:HB2	1.92	0.52
1:A:289:TYR:CD2	2:B:516:LEU:HD23	2.46	0.51
2:B:71:PHE:C	2:B:329:LYS:HG2	2.35	0.51
2:B:519:ALA:O	2:B:604:ILE:HD11	2.10	0.51
2:B:119:LYS:H	2:B:267:CYS:HA	1.75	0.51
1:A:314:LYS:HG3	1:A:436:PHE:CZ	2.46	0.51
2:B:264:SER:OG	2:B:271:TYR:HA	2.11	0.50
2:B:249:ARG:O	2:B:253:GLU:HG3	2.11	0.50
2:B:369:TYR:HA	2:B:372:ILE:CD1	2.42	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:464:MET:HE1	2:B:505:SER:HB3	1.94	0.50
2:B:28:CYS:HB3	2:B:33:ILE:HB	1.94	0.50
2:B:38:VAL:HA	2:B:46:ILE:O	2.12	0.49
2:B:420:LEU:HG	2:B:591:VAL:HG12	1.94	0.49
2:B:312:LEU:HD23	2:B:360:VAL:HG11	1.93	0.49
1:A:545:SER:HB3	3:A:601:FB9:O3	2.12	0.49
2:B:33:ILE:HG23	2:B:50:GLU:O	2.12	0.49
2:B:252:ALA:O	2:B:256:VAL:HG22	2.13	0.49
2:B:456:VAL:HG11	2:B:487:SER:HB2	1.94	0.49
2:B:314:GLY:O	2:B:372:ILE:HG23	2.12	0.49
2:B:83:LEU:HD12	2:B:83:LEU:H	1.77	0.48
2:B:13:GLU:C	2:B:15:LEU:N	2.68	0.48
2:B:111:VAL:HG23	2:B:294:TYR:CG	2.49	0.48
2:B:223:PRO:CD	2:B:224:PRO:HD2	2.42	0.48
2:B:338:ASP:OD1	2:B:339:SER:N	2.46	0.48
1:A:524:ARG:HB3	1:A:526:GLU:OE2	2.14	0.48
2:B:72:ILE:HA	2:B:329:LYS:HE2	1.94	0.48
2:B:106:ARG:NH1	2:B:206:PHE:CA	2.77	0.48
2:B:512:PHE:CD1	2:B:512:PHE:N	2.82	0.48
1:A:478:ALA:HB2	2:B:377:ILE:HG23	1.96	0.48
2:B:13:GLU:O	2:B:16:GLY:N	2.47	0.48
2:B:206:PHE:HB2	2:B:208:LYS:O	2.14	0.48
2:B:117:ASN:N	2:B:239:ASN:HB3	2.29	0.47
2:B:172:PHE:O	2:B:174:PRO:HD3	2.14	0.47
2:B:199:TYR:O	2:B:203:ILE:HG13	2.14	0.47
1:A:344:PRO:HG3	1:A:391:TRP:CD2	2.49	0.47
2:B:392:PHE:O	2:B:396:LEU:HD12	2.15	0.47
2:B:561:ASP:OD1	2:B:563:SER:OG	2.33	0.46
2:B:157:ASN:O	2:B:214:ASP:HB2	2.15	0.46
1:A:560:ARG:O	1:A:566:ARG:O	2.34	0.46
2:B:174:PRO:O	2:B:229:ASP:HB2	2.15	0.46
2:B:187:ILE:HD12	2:B:187:ILE:H	1.79	0.46
1:A:310:MET:HB2	1:A:546:LEU:HD21	1.97	0.46
2:B:109:TYR:HE1	2:B:288:LYS:HE3	1.81	0.46
2:B:462:VAL:HG13	2:B:463:ASN:N	2.31	0.46
1:A:332:GLU:OE2	1:A:415:ARG:NH2	2.30	0.46
2:B:153:TYR:CE1	2:B:158:PHE:CE1	3.04	0.46
2:B:538:ILE:HD13	2:B:550:SER:HA	1.98	0.46
2:B:107:ARG:NH1	2:B:245:THR:HG23	2.30	0.46
2:B:416:TYR:CE1	2:B:423:HIS:HB3	2.51	0.46
2:B:552:VAL:HG12	2:B:576:PRO:HD2	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:467:LEU:N	1:A:467:LEU:HD12	2.31	0.46
2:B:70:SER:HA	2:B:74:LYS:O	2.16	0.46
2:B:482:GLU:O	2:B:504:LEU:HD12	2.16	0.46
1:A:358:ILE:CG2	1:A:361:PRO:HD2	2.46	0.45
2:B:424:ARG:HH22	2:B:431:ILE:HA	1.81	0.45
1:A:345:GLN:HA	1:A:350:ARG:HD3	1.98	0.45
2:B:164:PHE:HB3	2:B:208:LYS:HB2	1.99	0.45
2:B:384:SER:C	2:B:612:MET:HE1	2.42	0.45
1:A:548:ARG:NH1	1:A:551:MET:HG3	2.32	0.45
2:B:572:ILE:HG22	2:B:581:PHE:CZ	2.52	0.45
1:A:452:LEU:N	1:A:452:LEU:HD12	2.32	0.45
1:A:458:VAL:O	1:A:543:GLY:HA2	2.16	0.45
1:A:517:VAL:O	3:A:601:FB9:C12	2.64	0.45
2:B:172:PHE:HB3	2:B:219:ILE:HB	1.99	0.45
2:B:158:PHE:CE2	2:B:276:PHE:HZ	2.34	0.45
2:B:67:ALA:HA	2:B:331:MET:HG3	1.98	0.45
1:A:551:MET:CB	1:A:556:ILE:HB	2.47	0.45
2:B:92:LYS:HB2	2:B:92:LYS:HE3	1.72	0.44
2:B:399:CYS:O	2:B:400:THR:OG1	2.24	0.44
1:A:361:PRO:O	1:A:403:ASN:ND2	2.46	0.44
2:B:64:LEU:O	2:B:65:CYS:C	2.60	0.44
2:B:107:ARG:HG2	2:B:162:TYR:OH	2.17	0.44
2:B:58:LEU:CD1	2:B:64:LEU:HA	2.47	0.44
2:B:500:ASN:O	2:B:501:LYS:HD3	2.17	0.44
1:A:332:GLU:HG3	1:A:333:SER:O	2.17	0.44
2:B:2:PRO:HD2	2:B:51:VAL:O	2.17	0.44
2:B:482:GLU:HG3	2:B:483:ILE:N	2.33	0.44
2:B:611:ILE:O	2:B:615:LEU:HG	2.18	0.44
2:B:31:PHE:CE1	2:B:330:MET:HE1	2.53	0.44
2:B:514:ALA:HB1	2:B:586:ILE:HD13	2.00	0.44
2:B:31:PHE:HE1	2:B:330:MET:HE1	1.83	0.43
2:B:71:PHE:O	2:B:329:LYS:CG	2.66	0.43
2:B:222:LEU:HD23	2:B:223:PRO:HD3	2.01	0.43
2:B:222:LEU:HD23	2:B:223:PRO:CD	2.48	0.43
2:B:514:ALA:HB3	2:B:586:ILE:CD1	2.48	0.43
2:B:186:PHE:CE2	2:B:224:PRO:HG2	2.47	0.43
2:B:464:MET:HE3	2:B:464:MET:HB2	1.47	0.43
2:B:207:GLU:C	2:B:208:LYS:HD3	2.43	0.43
1:A:562:LEU:O	1:A:567:SER:CB	2.67	0.43
2:B:555:LYS:HG2	2:B:575:CYS:SG	2.59	0.43
1:A:289:TYR:CD2	2:B:516:LEU:CD2	3.02	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:471:HIS:O	1:A:475:THR:HG23	2.19	0.43
1:A:342:TYR:CE2	2:B:354:MET:HE1	2.54	0.42
1:A:335:PHE:HA	1:A:339:ASP:HB2	2.01	0.42
1:A:476:LEU:HD23	1:A:476:LEU:HA	1.84	0.42
2:B:210:PRO:O	2:B:222:LEU:HD23	2.19	0.42
1:A:342:TYR:CZ	1:A:526:GLU:HG2	2.53	0.42
2:B:376:LYS:HA	2:B:376:LYS:HD3	1.75	0.42
1:A:469:LEU:HD23	1:A:502:MET:CE	2.50	0.42
2:B:222:LEU:HB3	2:B:223:PRO:HD2	2.01	0.42
2:B:473:HIS:NE2	2:B:474[A]:ARG:HD2	2.35	0.42
2:B:97:MET:HG2	2:B:98:ARG:N	2.35	0.42
2:B:278:VAL:O	2:B:291:GLY:HA2	2.20	0.42
2:B:97:MET:SD	2:B:110:VAL:HG11	2.60	0.42
1:A:308:ARG:O	1:A:312:THR:HG23	2.20	0.41
2:B:419:MET:HB3	2:B:421:ARG:HD3	2.01	0.41
2:B:501:LYS:HD3	2:B:501:LYS:HA	1.79	0.41
1:A:359:LYS:O	1:A:362:GLU:HB2	2.20	0.41
2:B:340:SER:O	2:B:340:SER:OG	2.31	0.41
1:A:289:TYR:HB2	2:B:568:ARG:HH12	1.72	0.41
2:B:461:ILE:HG12	2:B:603:VAL:HG11	2.03	0.41
1:A:366:GLU:HG2	1:A:401:LYS:HZ1	1.85	0.41
2:B:372:ILE:H	2:B:372:ILE:HD12	1.86	0.41
1:A:342:TYR:CE2	2:B:354:MET:CE	3.03	0.41
1:A:491:LYS:HG2	1:A:492:PRO:O	2.21	0.41
1:A:323:GLU:O	1:A:435:PHE:HA	2.21	0.41
1:A:325:MET:HE3	1:A:325:MET:HB2	1.97	0.41
1:A:342:TYR:CE1	1:A:380:VAL:HG22	2.56	0.41
2:B:54:ASN:HD22	2:B:352:ASP:HB2	1.86	0.41
2:B:389:THR:HG21	2:B:502:ARG:CD	2.43	0.41
2:B:412:LYS:N	2:B:454:GLU:OE1	2.54	0.41
2:B:70:SER:C	2:B:72:ILE:N	2.79	0.40
2:B:292:TYR:HB2	2:B:294:TYR:CE2	2.56	0.40
2:B:527:LEU:HD23	2:B:527:LEU:HA	1.70	0.40
2:B:222:LEU:CD2	2:B:223:PRO:HD2	2.51	0.40
2:B:93:GLU:OE2	2:B:93:GLU:N	2.55	0.40
2:B:260:CYS:O	2:B:264:SER:HB3	2.22	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	297/299 (99%)	276 (93%)	21 (7%)	0	100	100
2	B	599/618 (97%)	545 (91%)	54 (9%)	0	100	100
All	All	896/917 (98%)	821 (92%)	75 (8%)	0	100	100

There are no Ramachandran outliers to report.

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	237/274 (86%)	236 (100%)	1 (0%)	84	90
2	B	442/572 (77%)	436 (99%)	6 (1%)	59	80
All	All	679/846 (80%)	672 (99%)	7 (1%)	68	84

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	451	HIS
2	B	206	PHE
2	B	349	TYR
2	B	382	LYS
2	B	462	VAL
2	B	513	THR
2	B	559	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	392	ASN
2	B	157	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	FB9	A	601	-	43,47,47	2.37	4 (9%)	50,66,66	0.78	1 (2%)

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	FB9	A	601	-	-	4/22/58/58	0/5/6/6

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	601	FB9	C6-N	-10.40	1.36	1.47
3	A	601	FB9	C5-N	-9.14	1.38	1.47
3	A	601	FB9	C9-N1	-2.91	1.41	1.47
3	A	601	FB9	C10-N2	-2.38	1.32	1.37

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	601	FB9	C20-C3-C2	-2.31	112.84	118.21

There are no chirality outliers.

All (4) torsion outliers are listed below:

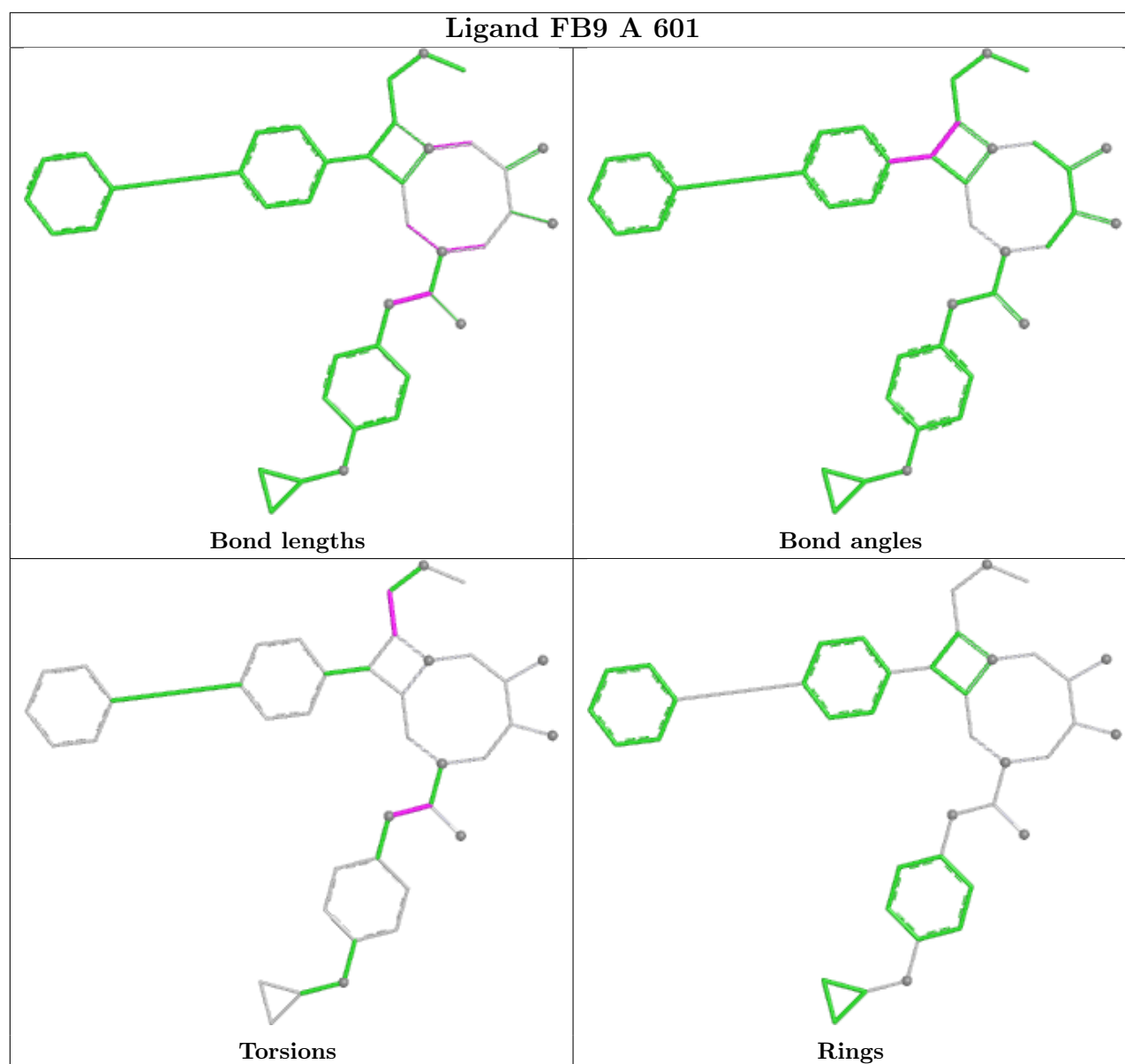
Mol	Chain	Res	Type	Atoms
3	A	601	FB9	O-C1-C2-N1
3	A	601	FB9	O-C1-C2-C3
3	A	601	FB9	N-C10-N2-C11
3	A	601	FB9	O3-C10-N2-C11

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	601	FB9	4	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	299/299 (100%)	0.22	12 (4%)	42 23	50, 81, 131, 148	0
2	B	606/618 (98%)	0.56	50 (8%)	17 9	45, 94, 153, 188	1 (0%)
All	All	905/917 (98%)	0.44	62 (6%)	23 12	45, 88, 148, 188	1 (0%)

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	516	LEU	7.0
1	A	565	TYR	5.3
2	B	201	LYS	4.4
2	B	7	HIS	4.4
2	B	277	PHE	3.8
2	B	304	CYS	3.8
1	A	273	TYR	3.7
2	B	47	TYR	3.6
2	B	318	LEU	3.6
1	A	278	ASN	3.4
2	B	260	CYS	3.3
2	B	205	ASP	3.0
2	B	273	ILE	3.0
2	B	95	HIS	3.0
2	B	296	VAL	3.0
2	B	316	LEU	2.8
2	B	242	ILE	2.7
2	B	192	ASP	2.7
2	B	281	ASP	2.6
2	B	259	ILE	2.6
2	B	223	PRO	2.6
2	B	138	ASN	2.6
2	B	512	PHE	2.5
1	A	569	ILE	2.5

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Mol	Chain	Res	Type	RSRZ
2	B	200	LEU	2.5
2	B	302	LEU	2.5
2	B	448	SER	2.5
2	B	202	ILE	2.4
2	B	106	ARG	2.4
2	B	117	ASN	2.4
2	B	44	LYS	2.4
2	B	315	ILE	2.4
2	B	103	VAL	2.3
2	B	15	LEU	2.3
2	B	222	LEU	2.3
2	B	153	TYR	2.3
2	B	104	ASP	2.3
2	B	314	GLY	2.3
1	A	472	LEU	2.2
2	B	342	PHE	2.2
2	B	75	TYR	2.2
2	B	263	LEU	2.2
1	A	327	THR	2.2
1	A	349	SER	2.2
2	B	2	PRO	2.2
2	B	262	MET	2.2
2	B	266	TYR	2.2
2	B	560	GLU	2.2
1	A	463	ILE	2.1
2	B	40	TYR	2.1
2	B	79	SER	2.1
1	A	431	ILE	2.1
2	B	6	VAL	2.1
2	B	38	VAL	2.1
2	B	241	PHE	2.1
2	B	356	PHE	2.1
1	A	275	LEU	2.0
1	A	450	THR	2.0
2	B	5	SER	2.0
2	B	336	VAL	2.0
2	B	224	PRO	2.0
1	A	420	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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](#)

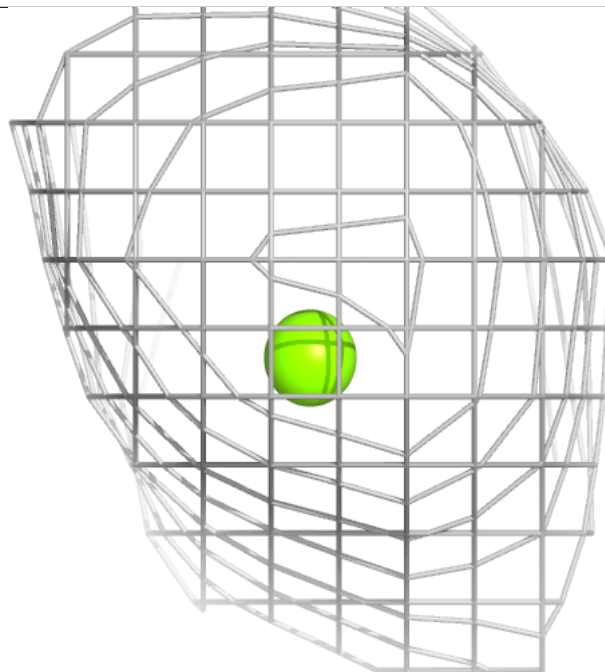
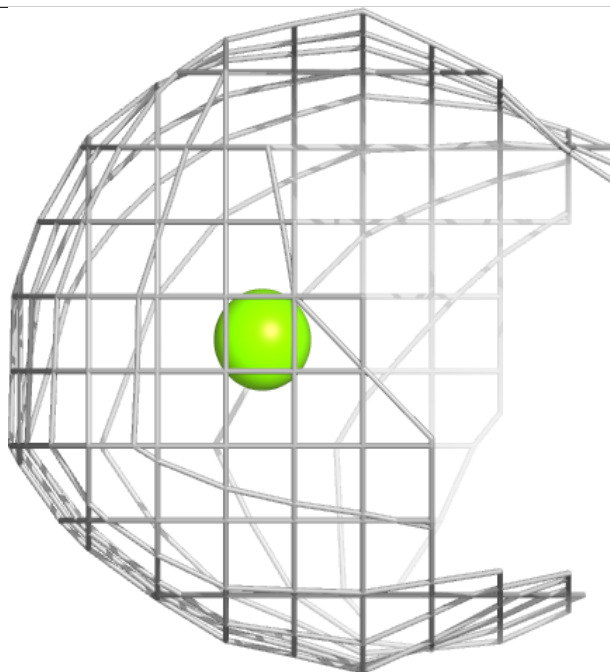
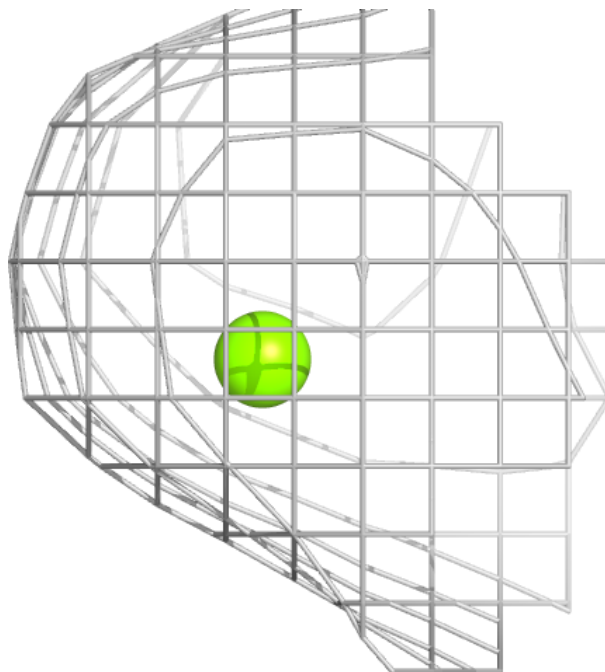
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

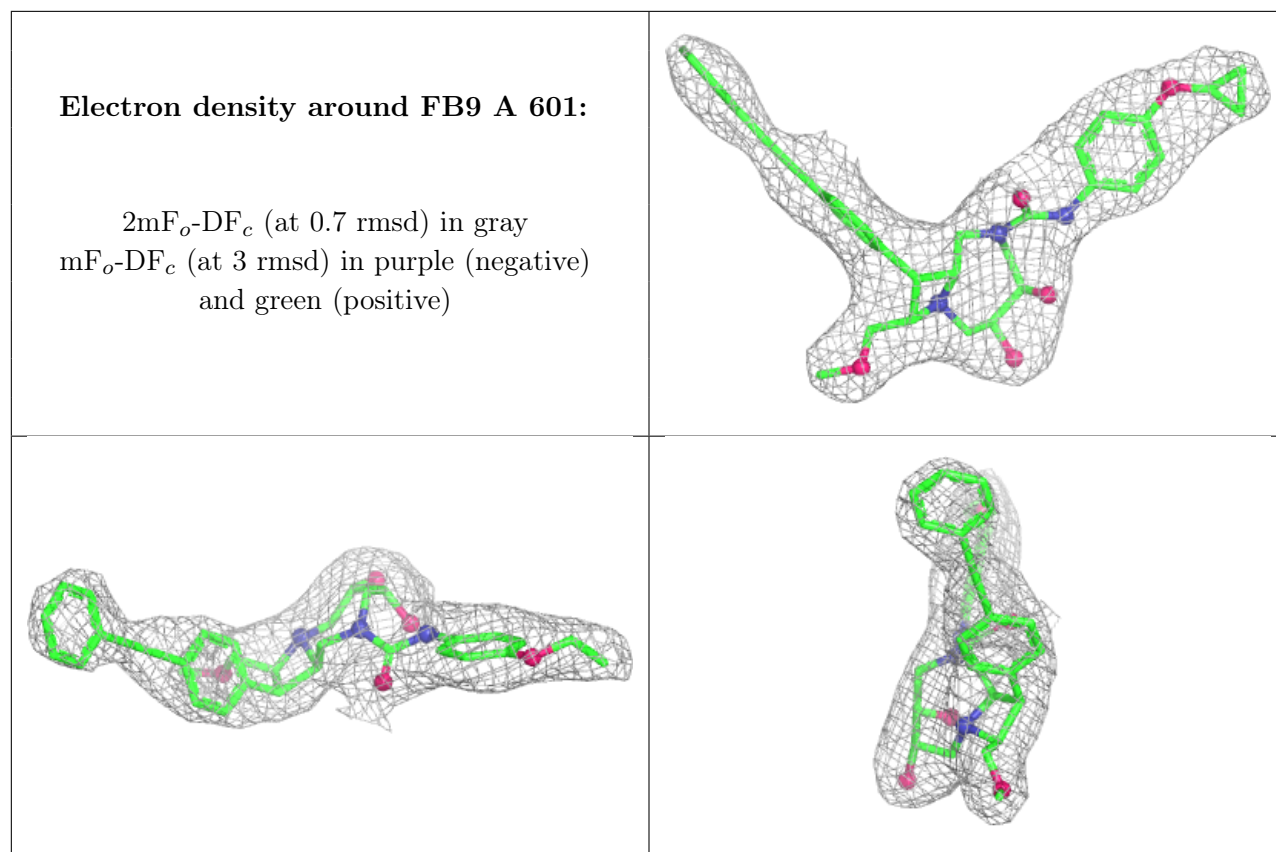
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	MG	B	701	1/1	0.91	0.16	83,83,83,83	0
3	FB9	A	601	42/42	0.95	0.12	52,62,73,76	42

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.

Electron density around MG B 701:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





6.5 Other polymers [i](#)

There are no such residues in this entry.