Full wwPDB X-ray Structure Validation Report (i)

Nov 2, 2026 – 03:34 pm GMT

PDB ID	:	30EH / pdb_000030eh
Title	:	2-heptyl-4(1H)-quinolone synthase PqsBC in complex with compound 35
Authors	:	Borgert, S.R.; Blankenfeldt, W.
Deposited on	:	2026-04-21
Resolution	:	1.95 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 (i) 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 (i)) were used in the production of this report:

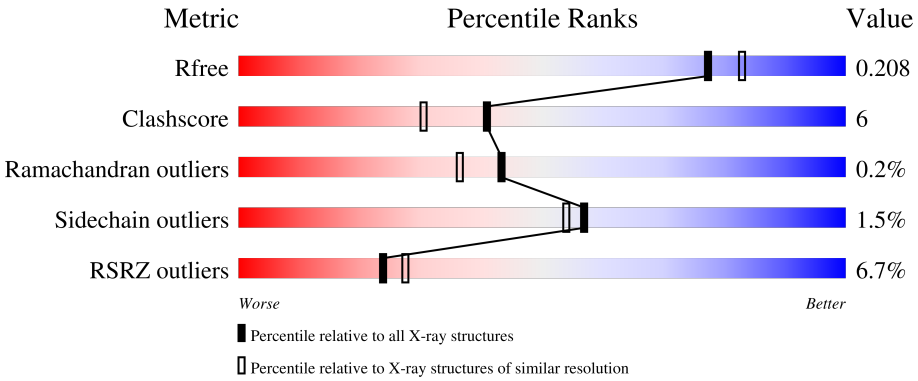
MolProbity	:	4-5-2 with Phenix2.2
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	2.2
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.51

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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.



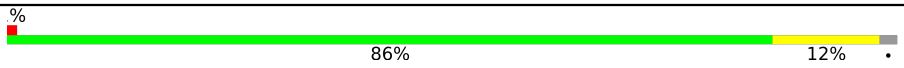
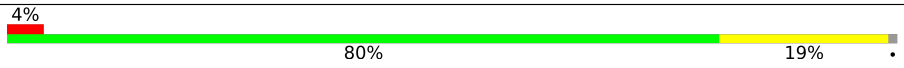
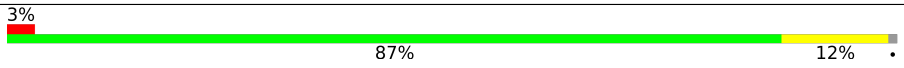
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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 <=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	365	8% 77% 16% • 5%
1	C	365	7% 81% 13% • •
1	E	365	10% 82% 13% •
1	G	365	16% 78% 16% • •
2	B	283	86% 12% •

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Mol	Chain	Length	Quality of chain
2	D	283	
2	F	283	
2	H	283	

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 20390 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 PqsC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	345	Total	C	N	O	S	0	4	0
			2681	1698	462	504	17			
1	C	349	Total	C	N	O	S	0	3	0
			2713	1723	468	504	18			
1	E	350	Total	C	N	O	S	0	5	0
			2719	1729	466	506	18			
1	G	350	Total	C	N	O	S	0	5	0
			2691	1713	457	503	18			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-16	MET	-	initiating methionine	UNP Q9I4X1
A	-15	GLY	-	expression tag	UNP Q9I4X1
A	-14	HIS	-	expression tag	UNP Q9I4X1
A	-13	HIS	-	expression tag	UNP Q9I4X1
A	-12	HIS	-	expression tag	UNP Q9I4X1
A	-11	HIS	-	expression tag	UNP Q9I4X1
A	-10	HIS	-	expression tag	UNP Q9I4X1
A	-9	HIS	-	expression tag	UNP Q9I4X1
A	-8	ALA	-	expression tag	UNP Q9I4X1
A	-7	GLU	-	expression tag	UNP Q9I4X1
A	-6	ASN	-	expression tag	UNP Q9I4X1
A	-5	LEU	-	expression tag	UNP Q9I4X1
A	-4	TYR	-	expression tag	UNP Q9I4X1
A	-3	PHE	-	expression tag	UNP Q9I4X1
A	-2	GLN	-	expression tag	UNP Q9I4X1
A	-1	GLY	-	expression tag	UNP Q9I4X1
A	0	HIS	-	expression tag	UNP Q9I4X1
A	129A	ALA	CYS	engineered mutation	UNP Q9I4X1
C	-16	MET	-	initiating methionine	UNP Q9I4X1
C	-15	GLY	-	expression tag	UNP Q9I4X1
C	-14	HIS	-	expression tag	UNP Q9I4X1

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-13	HIS	-	expression tag	UNP Q9I4X1
C	-12	HIS	-	expression tag	UNP Q9I4X1
C	-11	HIS	-	expression tag	UNP Q9I4X1
C	-10	HIS	-	expression tag	UNP Q9I4X1
C	-9	HIS	-	expression tag	UNP Q9I4X1
C	-8	ALA	-	expression tag	UNP Q9I4X1
C	-7	GLU	-	expression tag	UNP Q9I4X1
C	-6	ASN	-	expression tag	UNP Q9I4X1
C	-5	LEU	-	expression tag	UNP Q9I4X1
C	-4	TYR	-	expression tag	UNP Q9I4X1
C	-3	PHE	-	expression tag	UNP Q9I4X1
C	-2	GLN	-	expression tag	UNP Q9I4X1
C	-1	GLY	-	expression tag	UNP Q9I4X1
C	0	HIS	-	expression tag	UNP Q9I4X1
C	129A	ALA	CYS	engineered mutation	UNP Q9I4X1
E	-16	MET	-	initiating methionine	UNP Q9I4X1
E	-15	GLY	-	expression tag	UNP Q9I4X1
E	-14	HIS	-	expression tag	UNP Q9I4X1
E	-13	HIS	-	expression tag	UNP Q9I4X1
E	-12	HIS	-	expression tag	UNP Q9I4X1
E	-11	HIS	-	expression tag	UNP Q9I4X1
E	-10	HIS	-	expression tag	UNP Q9I4X1
E	-9	HIS	-	expression tag	UNP Q9I4X1
E	-8	ALA	-	expression tag	UNP Q9I4X1
E	-7	GLU	-	expression tag	UNP Q9I4X1
E	-6	ASN	-	expression tag	UNP Q9I4X1
E	-5	LEU	-	expression tag	UNP Q9I4X1
E	-4	TYR	-	expression tag	UNP Q9I4X1
E	-3	PHE	-	expression tag	UNP Q9I4X1
E	-2	GLN	-	expression tag	UNP Q9I4X1
E	-1	GLY	-	expression tag	UNP Q9I4X1
E	0	HIS	-	expression tag	UNP Q9I4X1
E	129A	ALA	CYS	engineered mutation	UNP Q9I4X1
G	-16	MET	-	initiating methionine	UNP Q9I4X1
G	-15	GLY	-	expression tag	UNP Q9I4X1
G	-14	HIS	-	expression tag	UNP Q9I4X1
G	-13	HIS	-	expression tag	UNP Q9I4X1
G	-12	HIS	-	expression tag	UNP Q9I4X1
G	-11	HIS	-	expression tag	UNP Q9I4X1
G	-10	HIS	-	expression tag	UNP Q9I4X1
G	-9	HIS	-	expression tag	UNP Q9I4X1
G	-8	ALA	-	expression tag	UNP Q9I4X1

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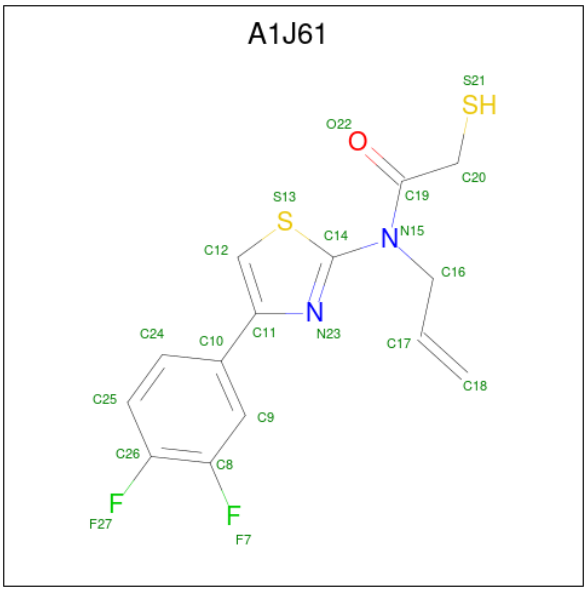
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Chain	Residue	Modelled	Actual	Comment	Reference
G	-7	GLU	-	expression tag	UNP Q9I4X1
G	-6	ASN	-	expression tag	UNP Q9I4X1
G	-5	LEU	-	expression tag	UNP Q9I4X1
G	-4	TYR	-	expression tag	UNP Q9I4X1
G	-3	PHE	-	expression tag	UNP Q9I4X1
G	-2	GLN	-	expression tag	UNP Q9I4X1
G	-1	GLY	-	expression tag	UNP Q9I4X1
G	0	HIS	-	expression tag	UNP Q9I4X1
G	129A	ALA	CYS	engineered mutation	UNP Q9I4X1

- Molecule 2 is a protein called 2-heptyl-4(1H)-quinolone synthase subunit PqsB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	279	Total	C	N	O	S	0	9	0
			2170	1374	384	403	9			
2	D	278	Total	C	N	O	S	0	11	0
			2169	1378	381	402	8			
2	F	279	Total	C	N	O	S	0	10	0
			2167	1375	379	405	8			
2	H	279	Total	C	N	O	S	0	9	0
			2159	1368	378	405	8			

- Molecule 3 is {N}-[4-[3,4-bis(fluoranyl)phenyl]-1,3-thiazol-2-yl]- {N}-prop-2-enyl-2-sulfanyl -ethanamide (CCD ID: A1J61) (formula: $C_{14}H_{12}F_2N_2OS_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	S	0	0
			21	14	2	2	1	2		
3	C	1	Total	C	F	N	O	S	0	0
			21	14	2	2	1	2		
3	E	1	Total	C	F	N	O	S	0	0
			21	14	2	2	1	2		
3	G	1	Total	C	F	N	O	S	0	0
			21	14	2	2	1	2		

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Cl	0	0
			1	1		
4	C	1	Total	Cl	0	0
			1	1		

- Molecule 5 is water.

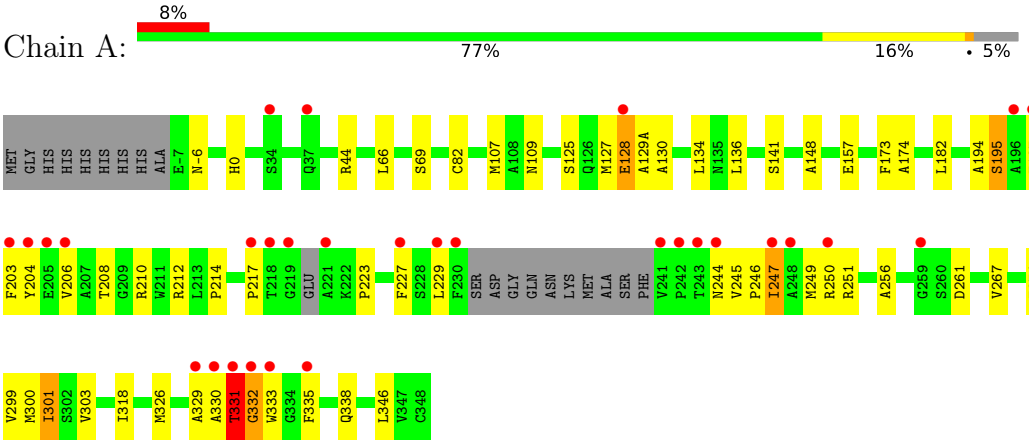
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	167	Total	O	0	0
			167	167		
5	B	193	Total	O	0	0
			193	193		
5	C	92	Total	O	0	0
			92	92		
5	D	135	Total	O	0	0
			135	135		
5	E	53	Total	O	0	0
			53	53		
5	F	61	Total	O	0	0
			61	61		
5	G	26	Total	O	0	0
			26	26		
5	H	108	Total	O	0	0
			108	108		

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.

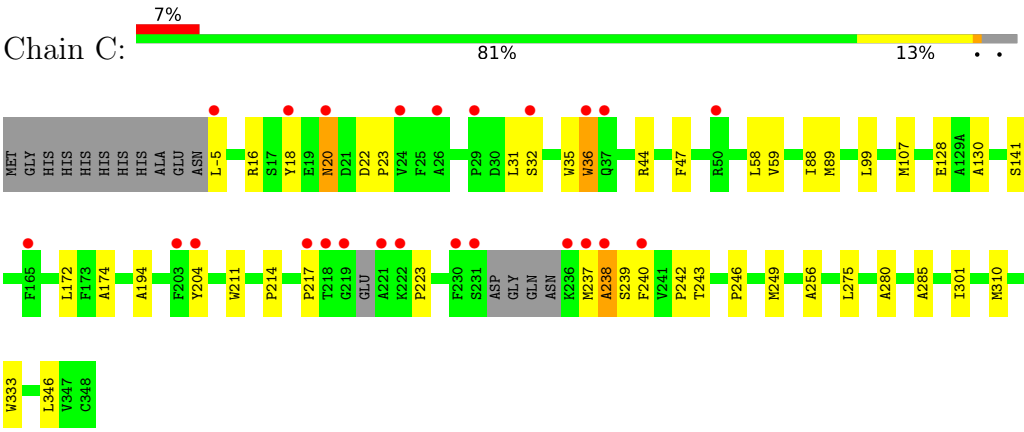
• Molecule 1: PqsC

Chain A:



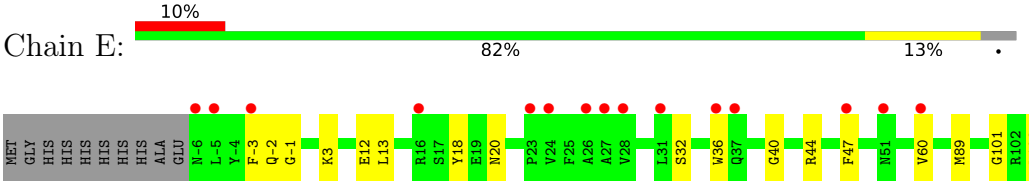
• Molecule 1: PqsC

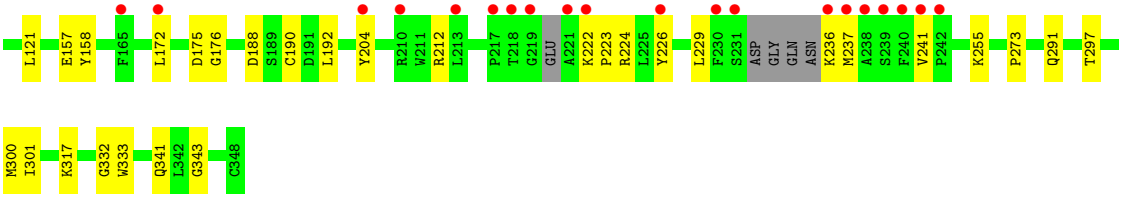
Chain C:



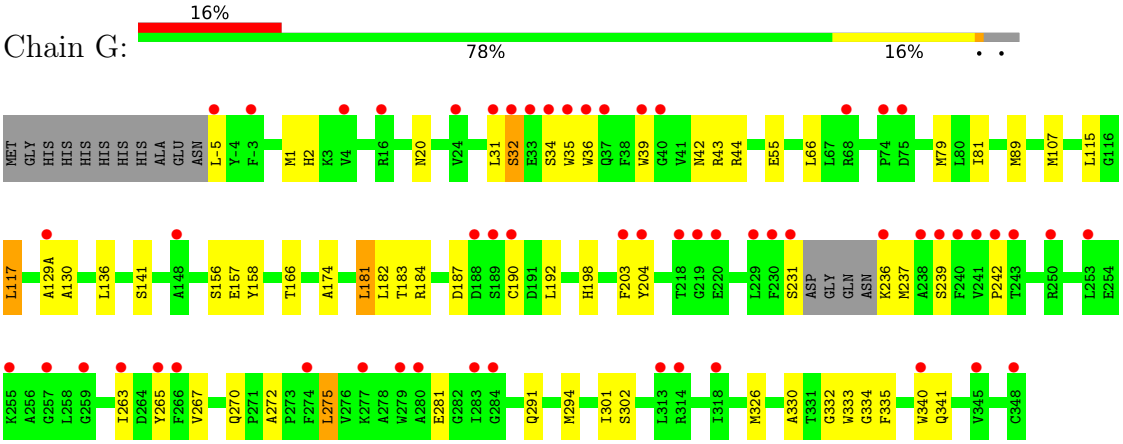
• Molecule 1: PqsC

Chain E:

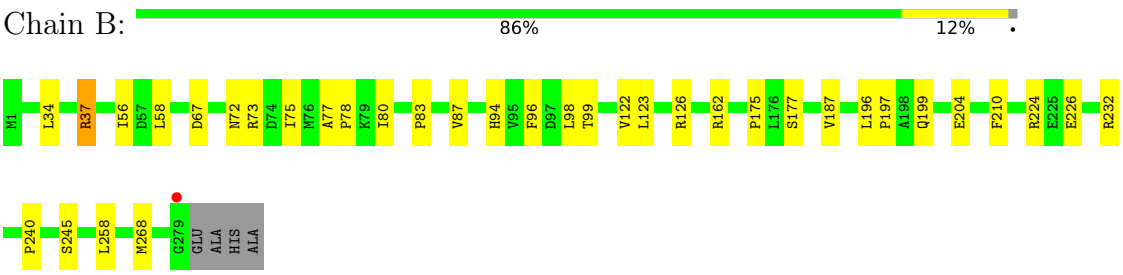




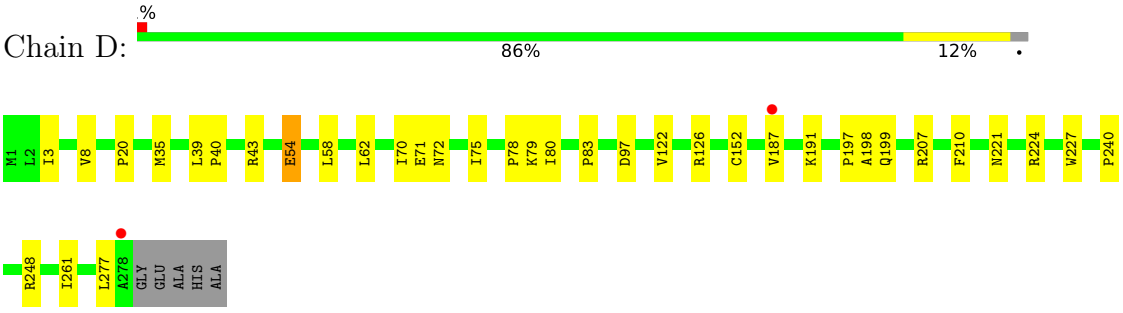
• Molecule 1: PqsC



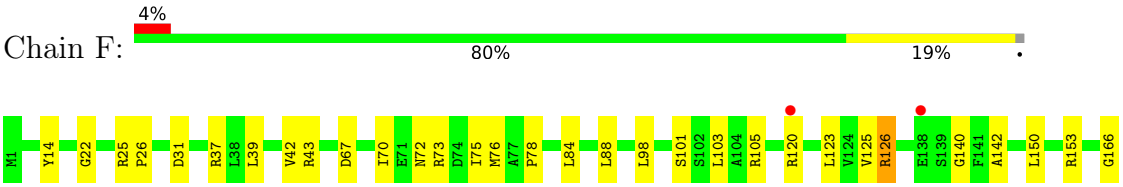
• Molecule 2: 2-heptyl-4(1H)-quinolone synthase subunit PqsB

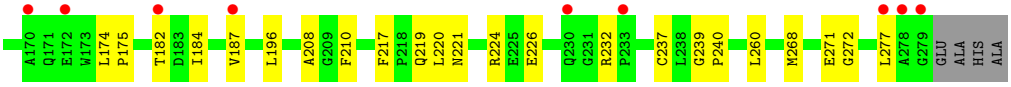


• Molecule 2: 2-heptyl-4(1H)-quinolone synthase subunit PqsB

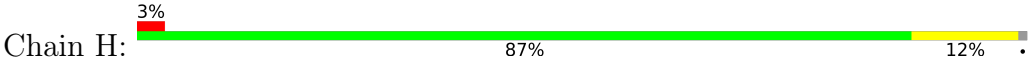


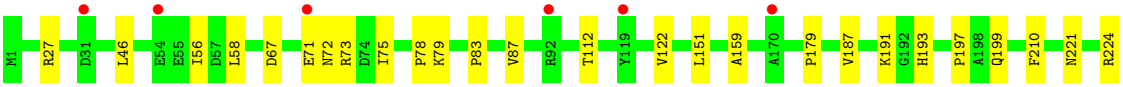
• Molecule 2: 2-heptyl-4(1H)-quinolone synthase subunit PqsB

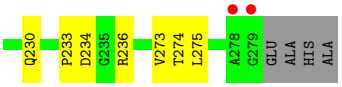




• Molecule 2: 2-heptyl-4(1H)-quinolone synthase subunit PqsB







4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	161.28Å 222.89Å 134.62Å 90.00° 111.62° 90.00°	Depositor
Resolution (Å)	83.23 – 1.95 83.23 – 1.95	Depositor EDS
% Data completeness (in resolution range)	54.9 (83.23-1.95) 54.9 (83.23-1.95)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.67 (at 1.82Å)	Xtriage
Refinement program	PHENIX (2.0_5936: ???)	Depositor
R, R_{free}	0.174 , 0.210 0.173 , 0.208	Depositor DCC
R_{free} test set	8792 reflections (2.23%)	wwPDB-VP
Wilson B-factor (Å^2)	27.1	Xtriage
Anisotropy	0.092	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 51.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	20390	wwPDB-VP
Average B, all atoms (Å^2)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.69% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: A1J61, CL

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	0.87	2/2751 (0.1%)	1.01	7/3733 (0.2%)
1	C	0.73	2/2782 (0.1%)	0.92	2/3773 (0.1%)
1	E	0.59	1/2794 (0.0%)	0.80	1/3789 (0.0%)
1	G	0.63	0/2766	0.84	1/3757 (0.0%)
2	B	0.95	0/2242	1.06	0/3044
2	D	0.76	0/2245	0.92	2/3049 (0.1%)
2	F	0.63	0/2239	0.80	2/3041 (0.1%)
2	H	0.69	0/2227	0.85	2/3026 (0.1%)
All	All	0.74	5/20046 (0.0%)	0.90	17/27212 (0.1%)

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
2	F	0	1
2	H	0	1
All	All	0	2

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	197	GLU	CA-CB	7.81	1.68	1.53
1	C	20	ASN	CB-CG	-6.61	1.35	1.52
1	E	241	VAL	C-N	5.37	1.40	1.33
1	A	128	GLU	CA-CB	-5.25	1.44	1.53
1	C	36	TRP	CG-CD2	-5.14	1.34	1.43

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	141	SER	CA-CB-OG	-6.89	97.33	111.10
1	E	343	GLY	N-CA-C	-6.26	103.74	112.13
1	C	141	SER	CA-CB-OG	-6.21	98.67	111.10
1	A	331	THR	N-CA-CB	5.75	118.59	110.36
1	C	20	ASN	CB-CG-ND2	5.70	124.95	116.40
1	A	125	SER	CA-CB-OG	-5.70	99.71	111.10
1	G	187	ASP	CB-CG-OD2	5.36	130.72	118.40
2	H	233	PRO	CA-C-N	5.23	131.14	121.52
2	H	233	PRO	C-N-CA	5.23	131.14	121.52
2	D	221	ASN	CA-C-N	-5.19	112.77	121.75
2	D	221	ASN	C-N-CA	-5.19	112.77	121.75
1	A	332	GLY	CA-C-N	5.15	130.97	121.70
1	A	332	GLY	C-N-CA	5.15	130.97	121.70
2	F	221	ASN	CA-C-N	-5.05	113.02	121.75
2	F	221	ASN	C-N-CA	-5.05	113.02	121.75
1	A	148	ALA	CA-C-N	-5.03	114.35	122.24
1	A	148	ALA	C-N-CA	-5.03	114.35	122.24

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	F	126	ARG	Sidechain
2	H	27	ARG	Sidechain

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	2681	0	2648	47	0
1	C	2713	0	2689	43	0
1	E	2719	0	2691	34	0
1	G	2691	0	2650	42	0
2	B	2170	0	2207	27	0
2	D	2169	0	2215	26	0
2	F	2167	0	2204	34	0
2	H	2159	0	2186	23	0
3	A	21	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	21	0	0	1	0
3	E	21	0	0	2	0
3	G	21	0	0	0	0
4	A	1	0	0	0	0
4	C	1	0	0	0	0
5	A	167	0	0	4	0
5	B	193	0	0	5	0
5	C	92	0	0	1	0
5	D	135	0	0	1	0
5	E	53	0	0	1	0
5	F	61	0	0	1	0
5	G	26	0	0	2	0
5	H	108	0	0	1	0
All	All	20390	0	19490	253	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 6.

All (253) 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:C:20:ASN:ND2	1:C:44:ARG:HH12	1.36	1.23
1:C:20:ASN:ND2	1:C:44:ARG:NH1	2.10	0.97
1:C:20:ASN:HD22	1:C:44:ARG:NH2	1.69	0.89
1:C:20:ASN:HD22	1:C:44:ARG:HH22	1.20	0.87
1:C:20:ASN:ND2	1:C:44:ARG:HH22	1.79	0.79
1:A:127:MET:HE3	1:A:134:LEU:HD22	1.67	0.76
1:C:20:ASN:HD22	1:C:44:ARG:CZ	1.98	0.76
1:C:20:ASN:HD22	1:C:44:ARG:NH1	1.82	0.76
1:C:44:ARG:HB2	1:C:174:ALA:HB2	1.71	0.72
2:B:245:SER:OG	1:C:-5:LEU:HD21	1.91	0.71
1:C:214:PRO:HB2	1:C:217:PRO:HA	1.73	0.70
1:E:13:LEU:HD22	1:E:300:MET:HE1	1.74	0.70
2:F:219:GLN:HG2	1:G:1:MET:HE1	1.72	0.70
1:C:20:ASN:ND2	1:C:44:ARG:NH2	2.39	0.68
1:A:127:MET:HE2	2:B:96:PHE:HB3	1.76	0.68
1:E:121:LEU:HD11	2:F:105:ARG:HD2	1.76	0.67
2:F:226[B]:GLU:HG2	2:F:232:ARG:HD3	1.78	0.65
1:C:239:SER:HA	1:C:242:PRO:HG2	1.78	0.65
1:C:20:ASN:ND2	1:C:44:ARG:CZ	2.60	0.64
2:D:72:ASN:HB3	2:D:75:ILE:HD12	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:162:ARG:HD3	5:B:433:HOH:O	1.99	0.63
2:F:210:PHE:CD2	2:F:224:ARG:HD2	2.34	0.63
1:A:208:THR:HA	2:B:78:PRO:HD3	1.79	0.62
2:B:37:ARG:NH1	5:B:301:HOH:O	2.33	0.60
1:G:44:ARG:HB2	1:G:174:ALA:HB2	1.83	0.59
2:F:42:VAL:HG11	2:F:88:LEU:HD11	1.84	0.59
2:B:56[B]:ILE:HD13	2:B:123[B]:LEU:HD12	1.85	0.59
1:E:47:PHE:CE2	1:E:175:ASP:HB3	2.38	0.58
2:F:232:ARG:HH12	2:F:237:CYS:HA	1.69	0.58
1:E:223:PRO:HB3	2:F:187:VAL:HG12	1.85	0.58
1:G:129(A):ALA:O	1:G:302:SER:HA	2.04	0.58
1:E:89:MET:HE3	1:E:106:ARG:CZ	2.34	0.58
1:C:128:GLU:HB2	5:C:578:HOH:O	2.04	0.57
2:H:234:ASP:HB3	2:H:236:ARG:H	1.68	0.57
1:A:245:VAL:HG12	1:A:249:MET:HE2	1.87	0.57
1:E:222:LYS:HD2	2:F:184:ILE:O	2.04	0.57
1:E:20:ASN:HB3	1:E:36:TRP:CE2	2.39	0.57
1:E:332:GLY:N	1:E:333:TRP:HA	2.20	0.57
2:D:58:LEU:HB3	2:D:122[B]:VAL:HG12	1.86	0.57
2:F:14:TYR:CE2	2:F:26:PRO:HG3	2.40	0.57
2:F:123[A]:LEU:HG	2:F:125:VAL:HG13	1.88	0.56
2:B:34:LEU:HD23	2:B:80:ILE:HD13	1.87	0.56
1:G:270:GLN:HE21	1:G:270:GLN:HA	1.69	0.56
2:B:72:ASN:HB3	2:B:75:ILE:HD12	1.88	0.55
1:C:172:LEU:HD12	3:C:401:A1J61:C9	2.36	0.55
2:F:217:PHE:HB3	2:F:220:LEU:HD12	1.86	0.55
1:E:291:GLN:NE2	1:E:317:LYS:HE2	2.22	0.55
2:H:72:ASN:HB3	2:H:75:ILE:HD12	1.89	0.54
1:A:130:ALA:H	1:A:330:ALA:HB2	1.71	0.54
2:B:58:LEU:HD12	2:B:94:HIS:O	2.07	0.54
1:G:66:LEU:HD21	1:G:181:LEU:HB2	1.89	0.54
1:E:236:LYS:HD3	1:E:237:MET:H	1.72	0.54
2:F:39:LEU:HD21	2:F:43:ARG:NH2	2.23	0.53
2:D:3:ILE:HB	2:D:277:LEU:HD23	1.90	0.53
1:A:331:THR:C	1:A:333:TRP:HA	2.33	0.53
1:G:20:ASN:N	1:G:20:ASN:OD1	2.38	0.53
1:A:204:TYR:CE1	2:B:83:PRO:HD3	2.42	0.53
1:C:332:GLY:N	1:C:333:TRP:HA	2.24	0.53
1:E:212:ARG:HE	1:E:226:TYR:HB2	1.73	0.53
2:F:120:ARG:NH2	2:F:153:ARG:HD3	2.24	0.53
2:D:122[B]:VAL:HG23	2:D:152:CYS:HB2	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:46:LEU:HD21	2:H:56[A]:ILE:HD11	1.92	0.52
2:F:72:ASN:HB3	2:F:75:ILE:HD12	1.91	0.52
1:A:332:GLY:N	1:A:333:TRP:HA	2.24	0.52
1:A:204:TYR:CE1	2:B:78:PRO:HD2	2.45	0.52
2:D:54:GLU:CD	2:D:54:GLU:H	2.18	0.52
1:G:42:ASN:O	1:G:43:ARG:HG3	2.09	0.52
1:C:47:PHE:CE1	1:C:58:LEU:HD12	2.45	0.52
1:G:136:LEU:HD22	1:G:182:LEU:HD21	1.92	0.52
1:C:211:TRP:CG	2:D:191:LYS:HE2	2.44	0.52
2:D:210:PHE:CD2	2:D:224[A]:ARG:HD2	2.43	0.52
1:E:103:LEU:HD22	2:F:76:MET:HE1	1.92	0.51
1:E:212:ARG:HH21	1:E:226:TYR:HB3	1.75	0.51
2:D:207:ARG:NE	2:D:224[A]:ARG:HH12	2.08	0.51
1:C:223:PRO:HB3	2:D:187:VAL:HG12	1.93	0.51
2:H:71[A]:GLU:CD	2:H:71[A]:GLU:H	2.19	0.51
1:A:173:PHE:HE2	1:A:227:PHE:CE2	2.29	0.51
1:C:249:MET:HE2	1:C:325:VAL:HG11	1.92	0.51
1:A:195:SER:HB2	1:A:338:GLN:HG2	1.93	0.50
2:F:232:ARG:HG3	2:F:232:ARG:HH11	1.76	0.50
1:A:331:THR:HA	1:A:333:TRP:CE3	2.46	0.50
2:H:46:LEU:HD23	2:H:151:LEU:HD22	1.93	0.50
1:E:40:GLY:H	1:E:273:PRO:HD3	1.76	0.50
1:G:204:TYR:CE1	2:H:78:PRO:HD2	2.46	0.50
2:H:46:LEU:HD11	2:H:56[B]:ILE:HD11	1.94	0.50
1:G:192:LEU:HD23	1:G:340:TRP:CE2	2.47	0.50
1:A:244:ASN:HB3	1:A:335:PHE:CG	2.46	0.50
2:H:230:GLN:HG3	5:H:352:HOH:O	2.12	0.50
1:C:204:TYR:CE1	2:D:83:PRO:HD3	2.47	0.49
1:A:210:ARG:HD2	1:A:212:ARG:HD3	1.95	0.49
2:B:99:THR:HA	5:B:376:HOH:O	2.11	0.49
1:A:128:GLU:HG2	1:A:129(A):ALA:N	2.28	0.49
1:C:237:MET:C	1:C:239:SER:H	2.21	0.49
1:A:44:ARG:HB2	1:A:174:ALA:HB2	1.95	0.49
2:D:39:LEU:HD11	2:D:43:ARG:HE	1.77	0.49
1:E:175:ASP:O	1:E:300:MET:HE3	2.13	0.49
1:G:239:SER:C	1:G:242:PRO:HD2	2.38	0.48
1:E:3:LYS:HE3	1:E:188:ASP:OD2	2.12	0.48
2:H:210:PHE:CD2	2:H:224[A]:ARG:HD2	2.48	0.48
1:G:204:TYR:CE1	2:H:83:PRO:HD3	2.48	0.48
1:E:13:LEU:CD2	1:E:300:MET:HE1	2.41	0.48
1:C:89:MET:HE1	1:C:107:MET:HB2	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:236:LYS:HD3	1:G:237:MET:H	1.79	0.48
1:A:244:ASN:HB3	1:A:335:PHE:CD1	2.48	0.48
1:A:127:MET:HE1	1:A:134:LEU:HD13	1.95	0.48
1:C:204:TYR:HE1	2:D:78[A]:PRO:HB2	1.79	0.48
1:C:35:TRP:HD1	1:C:36:TRP:CD1	2.32	0.48
1:A:206:VAL:O	1:A:229:LEU:HA	2.14	0.47
1:A:247:ILE:HD13	1:A:250:ARG:HE	1.79	0.47
2:B:126[A]:ARG:HE	2:B:240[A]:PRO:HG3	1.79	0.47
1:C:237:MET:HB3	1:C:240:PHE:HB3	1.96	0.47
1:E:-2:GLN:HA	2:H:221:ASN:HD22	1.79	0.47
1:A:157:GLU:OE2	1:A:301:ILE:HB	2.14	0.47
2:F:25:ARG:NH2	2:F:142:ALA:H	2.13	0.47
1:G:275:LEU:HA	1:G:275:LEU:HD13	1.70	0.47
1:G:31:LEU:HD23	1:G:31:LEU:HA	1.72	0.47
1:A:246:PRO:O	1:A:250:ARG:HG3	2.15	0.47
1:A:0:HIS:HB3	5:A:596:HOH:O	2.15	0.46
1:G:31:LEU:HB2	5:G:504:HOH:O	2.15	0.46
1:G:42:ASN:HB2	1:G:43:ARG:NH1	2.30	0.46
1:A:129(A):ALA:HB3	1:A:330:ALA:HA	1.97	0.46
1:A:204:TYR:HE1	2:B:78:PRO:HB2	1.80	0.46
1:A:331:THR:HG22	3:A:401:A1J61:C16	2.45	0.46
1:C:16:ARG:HG2	1:C:18:TYR:CE1	2.50	0.46
2:D:197:PRO:O	2:D:199:GLN:HG3	2.14	0.46
1:E:12:GLU:C	1:E:13:LEU:HD23	2.39	0.46
1:E:18:TYR:HB2	1:E:44:ARG:HD2	1.96	0.46
2:H:79:LYS:O	2:H:83:PRO:HD2	2.16	0.46
2:F:105:ARG:HD3	2:F:271:GLU:OE2	2.15	0.46
1:E:-3:PHE:CE1	2:H:224[B]:ARG:HD2	2.51	0.46
2:D:71[A]:GLU:HG2	5:D:377:HOH:O	2.15	0.46
2:F:67:ASP:HB2	2:F:73:ARG:O	2.16	0.46
2:B:204:GLU:HG3	5:B:451:HOH:O	2.16	0.46
2:D:20:PRO:HG2	2:D:198:ALA:HB1	1.98	0.46
2:H:197:PRO:O	2:H:199:GLN:HG3	2.16	0.46
1:A:331:THR:HA	1:A:333:TRP:CD2	2.50	0.46
1:G:130:ALA:H	1:G:330:ALA:HB2	1.81	0.45
1:G:190:CYS:HA	1:G:341:GLN:O	2.16	0.45
1:G:203:PHE:CE2	1:G:237:MET:HA	2.51	0.45
1:A:182:LEU:N	1:A:182:LEU:HD12	2.31	0.45
1:A:214:PRO:HB2	1:A:217:PRO:HA	1.99	0.45
1:C:237:MET:HB2	1:C:238:ALA:H	1.51	0.45
2:D:35:MET:HE2	2:D:80:ILE:HD11	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:207:ARG:HG2	2:D:224[A]:ARG:NH1	2.32	0.45
1:C:194:ALA:HB3	1:C:256:ALA:HB2	1.99	0.45
2:F:84:LEU:O	2:F:88:LEU:HB2	2.17	0.45
1:A:157:GLU:HA	1:A:157:GLU:OE1	2.17	0.45
3:A:401:A1J61:C19	5:A:535:HOH:O	2.64	0.45
2:F:98:LEU:HD23	2:F:98:LEU:HA	1.79	0.45
2:B:67:ASP:HB2	2:B:73:ARG:O	2.16	0.45
1:G:198:HIS:CE1	1:G:335:PHE:HB2	2.51	0.45
2:F:217:PHE:CB	2:F:220:LEU:HD12	2.48	0.44
2:D:277:LEU:HD23	2:D:277:LEU:HA	1.85	0.44
1:G:265:TYR:OH	1:G:291:GLN:HG3	2.18	0.44
1:A:198:HIS:CG	1:A:251:ARG:HH12	2.36	0.44
1:A:204:TYR:CE1	2:B:78:PRO:HB2	2.53	0.44
2:B:175:PRO:HA	2:B:268:MET:HE3	2.00	0.44
1:G:20:ASN:HB3	1:G:36:TRP:CZ2	2.53	0.44
1:G:272:ALA:C	1:G:294:MET:HE1	2.43	0.44
1:A:128:GLU:HB3	1:A:330:ALA:HB1	1.99	0.44
1:A:69[B]:SER:OG	1:A:346:LEU:HD21	2.18	0.44
1:A:157:GLU:HG2	1:A:303:VAL:HG23	1.99	0.44
1:C:280:ALA:HB1	1:C:285:ALA:HB3	2.00	0.44
1:E:157:GLU:OE1	1:E:157:GLU:HA	2.16	0.44
1:E:255:LYS:NZ	5:E:501:HOH:O	2.48	0.44
1:G:20:ASN:HB3	1:G:36:TRP:CE2	2.52	0.44
1:G:79:MET:HE2	1:G:81:ILE:HG12	2.00	0.44
2:H:83:PRO:O	2:H:87:VAL:HG23	2.17	0.44
2:F:22:GLY:HA2	2:F:140:GLY:O	2.17	0.44
1:G:332:GLY:N	1:G:333:TRP:HA	2.33	0.44
2:H:159:ALA:O	2:H:273:VAL:HA	2.17	0.44
1:A:194:ALA:HB3	1:A:256:ALA:HB2	1.99	0.43
1:A:82:CYS:HB3	1:A:107:MET:HE3	1.99	0.43
2:B:226[B]:GLU:HG2	2:B:232:ARG:NE	2.34	0.43
1:C:31:LEU:HD23	1:C:31:LEU:HA	1.73	0.43
1:C:331:THR:C	1:C:333:TRP:HA	2.43	0.43
2:F:126:ARG:NE	2:F:240[A]:PRO:HG3	2.33	0.43
2:B:77:ALA:HA	2:B:78:PRO:C	2.43	0.43
2:B:83:PRO:O	2:B:87:VAL:HG23	2.19	0.43
2:B:177:SER:HB3	5:B:454:HOH:O	2.18	0.43
1:G:89:MET:HE2	1:G:89:MET:HB3	1.83	0.43
1:A:274:PHE:CZ	1:A:275:LEU:HG	2.53	0.43
1:A:318:ILE:HD13	1:A:318:ILE:HG21	1.85	0.43
1:C:243:THR:C	1:C:246:PRO:HD2	2.44	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:103:LEU:HD21	2:F:150:LEU:HD11	2.01	0.43
2:H:210:PHE:HD2	2:H:224[A]:ARG:HD2	1.84	0.43
1:E:60:VAL:HG23	1:E:115:LEU:HD21	2.01	0.43
2:H:67:ASP:HB2	2:H:73:ARG:O	2.19	0.43
1:A:299:VAL:HG11	5:A:535:HOH:O	2.18	0.43
2:B:258:LEU:HD12	2:B:258:LEU:C	2.44	0.43
1:E:297:THR:HB	1:E:300:MET:SD	2.59	0.43
2:F:166:GLY:HA3	2:F:208:ALA:HB3	1.99	0.43
2:D:62:LEU:HB2	2:D:126:ARG:HG2	2.00	0.43
1:G:55:GLU:HB3	1:G:158:TYR:HB2	2.01	0.43
1:G:107:MET:HE1	1:G:156:SER:HB3	2.00	0.43
1:G:263:ILE:HD13	1:G:263:ILE:HA	1.85	0.43
1:E:229:LEU:HD13	3:E:401:A1J61:F7	2.09	0.42
1:G:35:TRP:HD1	1:G:36:TRP:CD1	2.37	0.42
2:D:79:LYS:HD3	2:D:97:ASP:CG	2.45	0.42
1:E:47:PHE:HE2	1:E:158:TYR:HA	1.84	0.42
1:E:190:CYS:HA	1:E:341:GLN:O	2.19	0.42
1:E:224:ARG:HB2	1:E:226:TYR:CE2	2.54	0.42
2:F:239:GLY:HA3	2:F:240[A]:PRO:HD2	1.87	0.42
1:E:204:TYR:CE1	2:F:78:PRO:HD2	2.54	0.42
2:F:277:LEU:HD23	2:F:277:LEU:HA	1.87	0.42
2:D:39:LEU:HB3	2:D:40:PRO:HD3	2.01	0.42
2:H:58:LEU:HB3	2:H:122[A]:VAL:HG22	2.01	0.42
2:D:227:TRP:CE2	2:D:240[B]:PRO:HD3	2.54	0.42
1:C:22:ASP:CG	1:C:23:PRO:HD2	2.44	0.42
1:C:128:GLU:HB3	1:C:330:ALA:HB1	2.02	0.42
2:D:207:ARG:CD	2:D:224[A]:ARG:HH12	2.33	0.42
1:G:115:LEU:HB2	1:G:117:LEU:HD13	2.01	0.42
2:F:175:PRO:HA	2:F:268:MET:HE3	2.01	0.42
1:G:334:GLY:HA3	5:G:512:HOH:O	2.19	0.42
1:C:157:GLU:OE1	1:C:157:GLU:HA	2.19	0.42
2:H:274:THR:O	2:H:275:LEU:HD23	2.20	0.42
1:C:20:ASN:HB2	1:C:36:TRP:CZ2	2.55	0.42
2:D:8:VAL:HG12	2:D:248:ARG:HH12	1.85	0.42
1:A:223:PRO:HB3	2:B:187:VAL:HG12	2.02	0.41
1:C:59:VAL:HG21	1:C:156:SER:HB2	2.02	0.41
1:E:172:LEU:HD12	3:E:401:A1J61:C24	2.50	0.41
1:E:176:GLY:HA3	1:E:300:MET:CE	2.50	0.41
2:F:126:ARG:HG2	2:F:126:ARG:HH11	1.85	0.41
1:C:99:LEU:HD12	1:C:99:LEU:HA	1.82	0.41
2:H:179:PRO:HD2	2:H:193:HIS:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:197:PRO:O	2:B:199:GLN:HG3	2.20	0.41
1:A:203:PHE:N	1:A:203:PHE:CD2	2.88	0.41
2:F:260:LEU:O	2:F:272:GLY:HA2	2.21	0.41
2:B:210:PHE:CD2	2:B:224:ARG:HD2	2.55	0.41
1:C:310:MET:HB3	1:C:310:MET:HE2	1.74	0.41
2:D:261:ILE:HG21	2:D:261:ILE:HD13	1.77	0.41
1:G:2:HIS:CE1	1:G:184:ARG:CZ	3.04	0.41
1:G:39:TRP:CE3	1:G:272:ALA:HB2	2.56	0.41
1:G:267:VAL:O	1:G:326:MET:HA	2.21	0.41
1:E:-1:GLY:H	2:H:221:ASN:HD22	1.69	0.41
1:A:109:ASN:HB3	5:A:627:HOH:O	2.20	0.40
1:A:267:VAL:HB	1:A:326:MET:HG2	2.02	0.40
2:B:98:LEU:HA	2:B:98:LEU:HD23	1.87	0.40
1:G:181:LEU:HD22	1:G:183:THR:HG23	2.04	0.40
1:A:297:THR:HB	1:A:300:MET:HE2	2.03	0.40
2:D:70:ILE:HD11	2:D:75:ILE:HG21	2.04	0.40
2:F:70:ILE:HD11	2:F:75:ILE:HG21	2.04	0.40
1:G:32:SER:OG	1:G:34:SER:HB3	2.21	0.40
1:C:130:ALA:H	1:C:330:ALA:HB2	1.86	0.40
1:C:237:MET:HB3	1:C:240:PHE:CB	2.52	0.40
1:E:89:MET:O	1:E:101:GLY:HA2	2.21	0.40
1:A:129(A):ALA:HB3	1:A:329:ALA:O	2.21	0.40
1:C:275:LEU:HD23	1:C:275:LEU:HA	1.91	0.40
2:F:101:SER:HB2	5:F:336:HOH:O	2.20	0.40
1:G:141:SER:HB3	2:H:112:THR:HG22	2.03	0.40
1:A:66:LEU:HD23	1:A:66:LEU:C	2.46	0.40
1:G:157:GLU:OE1	1:G:157:GLU:HA	2.21	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	343/365 (94%)	332 (97%)	10 (3%)	1 (0%)	36	28
1	C	346/365 (95%)	333 (96%)	11 (3%)	2 (1%)	21	12
1	E	349/365 (96%)	336 (96%)	12 (3%)	1 (0%)	36	28
1	G	351/365 (96%)	337 (96%)	13 (4%)	1 (0%)	36	28
2	B	287/283 (101%)	284 (99%)	3 (1%)	0	100	100
2	D	287/283 (101%)	282 (98%)	5 (2%)	0	100	100
2	F	287/283 (101%)	282 (98%)	5 (2%)	0	100	100
2	H	286/283 (101%)	282 (99%)	4 (1%)	0	100	100
All	All	2536/2592 (98%)	2468 (97%)	63 (2%)	5 (0%)	43	36

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	238	ALA
1	E	301	ILE
1	G	301	ILE
1	A	301	ILE
1	C	301	ILE

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	285/299 (95%)	279 (98%)	6 (2%)	47	41
1	C	288/299 (96%)	284 (99%)	4 (1%)	59	56
1	E	288/299 (96%)	286 (99%)	2 (1%)	76	76
1	G	283/299 (95%)	276 (98%)	7 (2%)	42	34
2	B	231/223 (104%)	228 (99%)	3 (1%)	61	58
2	D	232/223 (104%)	231 (100%)	1 (0%)	84	85
2	F	231/223 (104%)	226 (98%)	5 (2%)	45	40
2	H	229/223 (103%)	227 (99%)	2 (1%)	70	70

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	2067/2088 (99%)	2037 (98%)	30 (2%)	57 54

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-6	ASN
1	A	136	LEU
1	A	195	SER
1	A	247	ILE
1	A	261	ASP
1	A	331	THR
2	B	37	ARG
2	B	122	VAL
2	B	196	LEU
1	C	32	SER
1	C	88	ILE
1	C	159	ILE
1	C	346	LEU
2	D	54	GLU
1	E	32	SER
1	E	192	LEU
2	F	31	ASP
2	F	37	ARG
2	F	174	LEU
2	F	182	THR
2	F	196	LEU
1	G	-5	LEU
1	G	32	SER
1	G	117	LEU
1	G	166	THR
1	G	181	LEU
1	G	231	SER
1	G	275	LEU
2	H	187	VAL
2	H	191	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	135	ASN
2	B	251	GLN

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Mol	Chain	Res	Type
1	C	0	HIS
1	C	20	ASN
1	C	109	ASN
1	C	126	GLN
1	E	0	HIS
1	E	216	ASN
2	F	186	GLN
2	F	219	GLN
2	H	4	GLN
2	H	29	GLN
2	H	221	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 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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	A1J61	A	401	1	21,22,22	0.56	0	25,30,30	0.87	1 (4%)
3	A1J61	C	401	1	21,22,22	0.47	0	25,30,30	1.19	2 (8%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
3	A1J61	E	401	1	21,22,22	0.27	0	25,30,30	0.91	1 (4%)
3	A1J61	G	401	1	21,22,22	0.63	0	25,30,30	0.94	2 (8%)

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	A1J61	A	401	1	-	5/15/17/17	0/2/2/2
3	A1J61	C	401	1	-	2/15/17/17	0/2/2/2
3	A1J61	E	401	1	-	8/15/17/17	0/2/2/2
3	A1J61	G	401	1	-	7/15/17/17	0/2/2/2

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	C	401	A1J61	S13-C14-N15	4.01	126.31	122.57
3	E	401	A1J61	C16-N15-C19	3.48	122.13	117.12
3	G	401	A1J61	C16-N15-C19	3.12	121.62	117.12
3	A	401	A1J61	S13-C14-N15	2.67	125.06	122.57
3	G	401	A1J61	S13-C14-N15	2.28	124.69	122.57
3	C	401	A1J61	C16-N15-C19	2.25	120.36	117.12

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	401	A1J61	C20-C19-N15-C14
3	A	401	A1J61	O22-C19-N15-C14
3	C	401	A1J61	C20-C19-N15-C14
3	E	401	A1J61	N23-C14-N15-C19
3	E	401	A1J61	S13-C14-N15-C16
3	E	401	A1J61	S13-C14-N15-C19
3	E	401	A1J61	C20-C19-N15-C14
3	E	401	A1J61	O22-C19-N15-C14
3	G	401	A1J61	N15-C16-C17-C18
3	G	401	A1J61	N23-C14-N15-C16
3	G	401	A1J61	N23-C14-N15-C19

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Mol	Chain	Res	Type	Atoms
3	G	401	A1J61	S13-C14-N15-C16
3	G	401	A1J61	S13-C14-N15-C19
3	A	401	A1J61	N15-C16-C17-C18
3	A	401	A1J61	O22-C19-N15-C16
3	E	401	A1J61	O22-C19-N15-C16
3	E	401	A1J61	N23-C14-N15-C16
3	G	401	A1J61	C20-C19-N15-C14
3	A	401	A1J61	C20-C19-N15-C16
3	E	401	A1J61	C20-C19-N15-C16
3	G	401	A1J61	O22-C19-N15-C16
3	C	401	A1J61	O22-C19-N15-C16

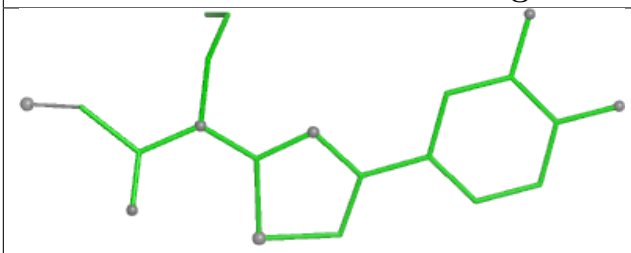
There are no ring outliers.

3 monomers are involved in 5 short contacts:

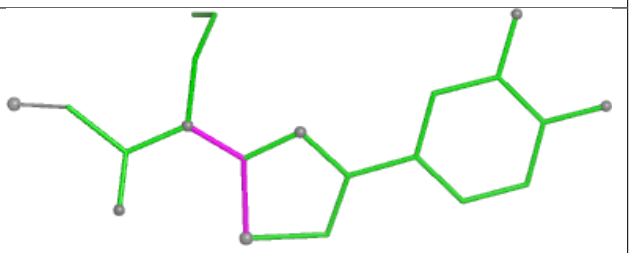
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	401	A1J61	2	0
3	C	401	A1J61	1	0
3	E	401	A1J61	2	0

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 then 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.

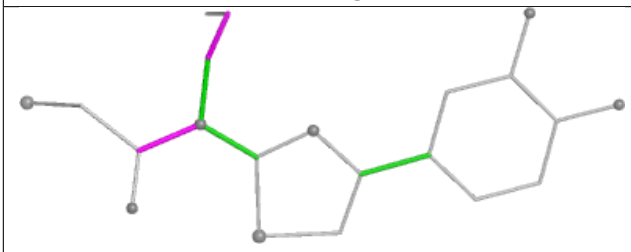
Ligand A1J61 A 401



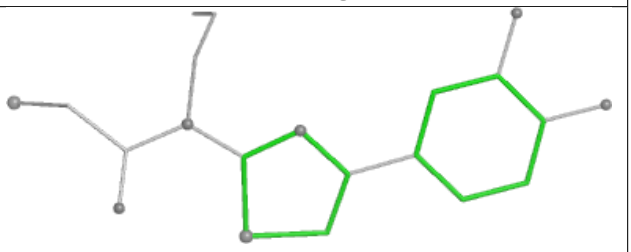
Bond lengths



Bond angles

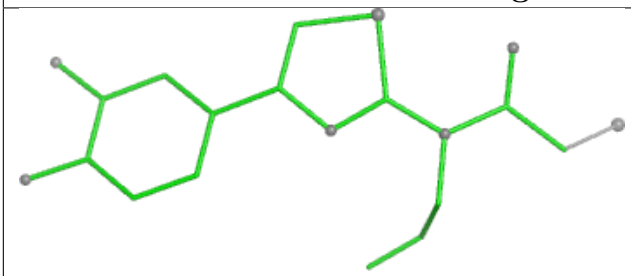


Torsions

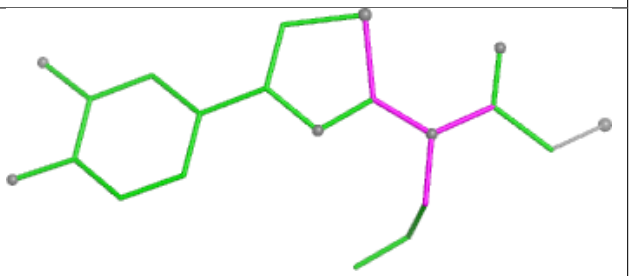


Rings

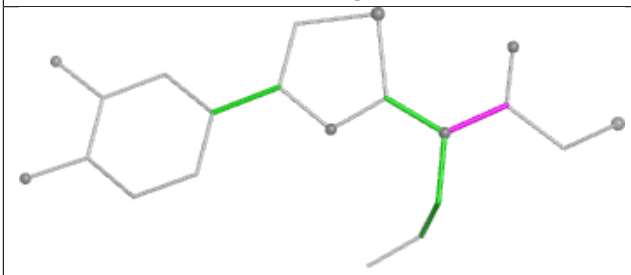
Ligand A1J61 C 401



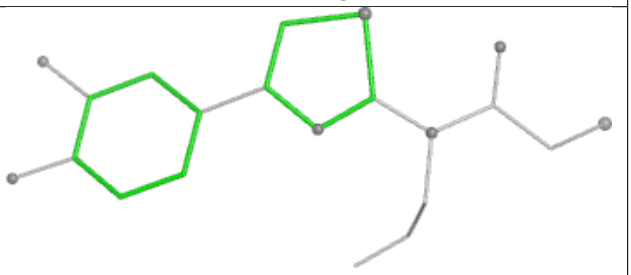
Bond lengths



Bond angles

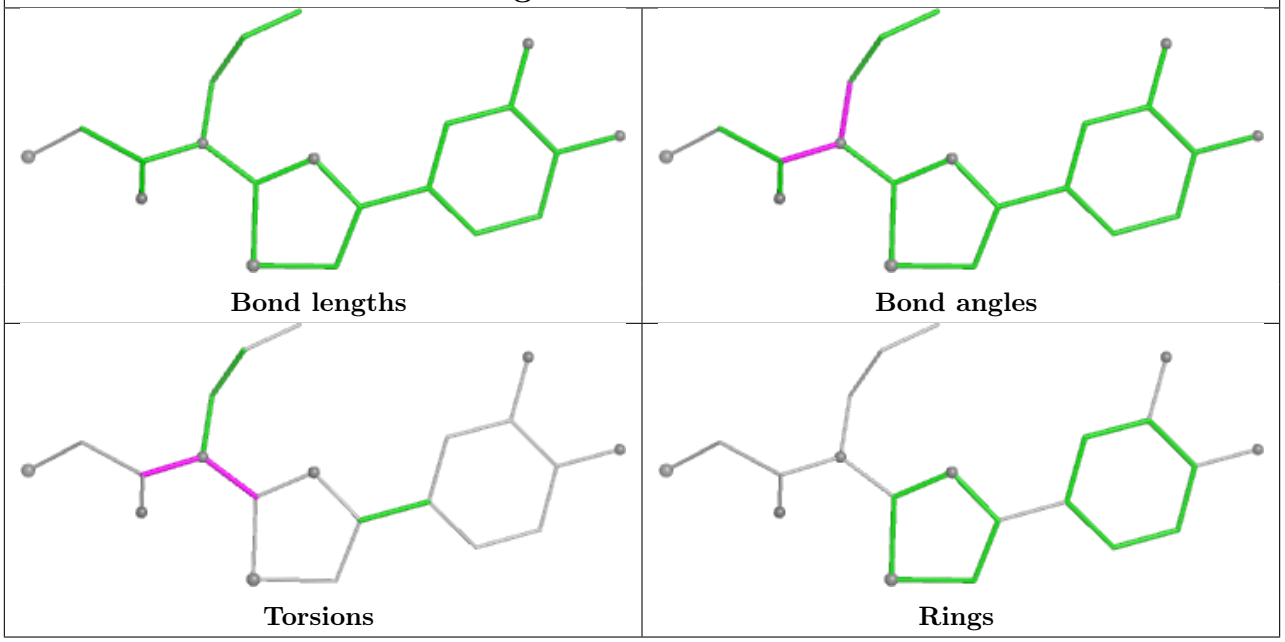


Torsions

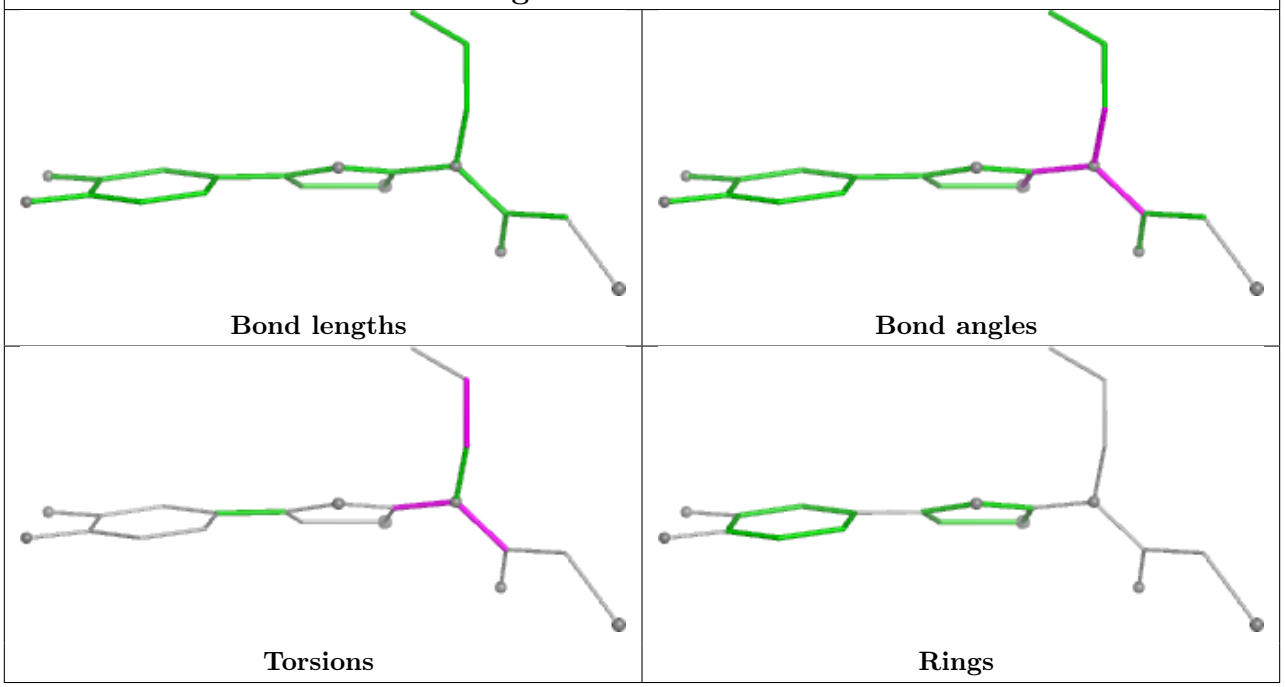


Rings

Ligand A1J61 E 401



Ligand A1J61 G 401



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} 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	345/365 (94%)	0.31	31 (8%) 15 17	16, 40, 75, 100	4 (1%)
1	C	349/365 (95%)	0.35	24 (6%) 23 26	24, 49, 81, 100	4 (1%)
1	E	350/365 (95%)	0.70	35 (10%) 12 14	24, 60, 97, 110	6 (1%)
1	G	350/365 (95%)	1.22	57 (16%) 4 5	36, 69, 90, 107	5 (1%)
2	B	279/283 (98%)	-0.52	1 (0%) 88 91	11, 26, 53, 70	9 (3%)
2	D	278/283 (98%)	-0.32	2 (0%) 84 88	13, 34, 59, 71	11 (3%)
2	F	279/283 (98%)	0.29	11 (3%) 43 50	24, 50, 78, 105	10 (3%)
2	H	279/283 (98%)	0.05	8 (2%) 53 60	22, 45, 68, 98	9 (3%)
All	All	2509/2592 (96%)	0.30	169 (6%) 24 27	11, 48, 84, 110	58 (2%)

All (169) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	36	TRP	8.1
1	C	36	TRP	8.0
2	H	279	GLY	7.7
1	A	219	GLY	7.3
1	E	219	GLY	6.9
1	A	230	PHE	6.7
1	E	238	ALA	6.6
1	A	204	TYR	6.5
2	F	279	GLY	5.9
1	C	238	ALA	5.8
1	A	242	PRO	5.5
1	E	36	TRP	5.4
1	A	241	VAL	5.4
1	G	-5	LEU	5.3
1	G	218	THR	5.3
1	G	236	LYS	5.2

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Mol	Chain	Res	Type	RSRZ
1	G	238	ALA	5.1
1	A	247	ILE	5.0
1	G	231	SER	4.9
1	G	204	TYR	4.9
1	E	240	PHE	4.9
2	B	279	GLY	4.8
1	E	231	SER	4.7
1	C	231	SER	4.7
1	C	236	LYS	4.7
1	A	335	PHE	4.5
1	G	277[A]	LYS	4.5
1	C	218	THR	4.5
1	E	221	ALA	4.4
1	G	240	PHE	4.4
1	A	331	THR	4.3
1	E	236	LYS	4.2
1	A	243	THR	4.2
1	A	203	PHE	4.1
1	C	221	ALA	4.0
1	A	196	ALA	4.0
1	E	239	SER	4.0
1	G	274	PHE	3.9
1	G	37	GLN	3.9
2	F	278	ALA	3.9
2	D	187	VAL	3.9
1	C	240	PHE	3.9
1	A	221	ALA	3.8
1	A	128	GLU	3.8
1	G	33	GLU	3.8
1	E	218	THR	3.7
1	C	230	PHE	3.7
1	G	239	SER	3.7
1	E	241	VAL	3.7
1	C	219	GLY	3.7
1	A	206	VAL	3.6
2	D	278	ALA	3.5
1	E	51	ASN	3.5
1	E	230	PHE	3.4
1	A	218	THR	3.4
1	C	237	MET	3.4
1	G	203	PHE	3.3
2	F	182	THR	3.3

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Mol	Chain	Res	Type	RSRZ
2	H	54[A]	GLU	3.3
1	E	204	TYR	3.3
1	A	201	ALA	3.3
1	G	129(A)	ALA	3.3
1	A	227	PHE	3.3
1	G	220	GLU	3.2
1	A	205	GLU	3.2
1	G	243	THR	3.2
1	E	242	PRO	3.2
1	C	217	PRO	3.1
1	E	213	LEU	3.1
1	E	-3	PHE	3.1
1	E	217	PRO	3.1
1	G	255	LYS	3.1
1	C	-5	LEU	3.0
1	E	31	LEU	2.9
1	G	219	GLY	2.9
1	A	197	GLU	2.9
1	A	333	TRP	2.9
1	G	340	TRP	2.9
1	A	244	ASN	2.9
1	G	241	VAL	2.9
1	C	204	TYR	2.9
1	G	-3	PHE	2.9
1	E	-6	ASN	2.9
1	A	37	GLN	2.8
1	A	248	ALA	2.7
1	C	165	PHE	2.7
1	G	31	LEU	2.7
2	F	187	VAL	2.7
1	G	35	TRP	2.6
1	E	210	ARG	2.6
1	G	314	ARG	2.6
1	G	39	TRP	2.6
1	G	230	PHE	2.6
1	A	217	PRO	2.6
1	G	313	LEU	2.6
2	H	31	ASP	2.5
1	G	24	VAL	2.5
1	G	74	PRO	2.5
1	C	18	TYR	2.4
1	G	229	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	E	222	LYS	2.4
1	C	24	VAL	2.4
1	A	259	GLY	2.4
1	G	257	GLY	2.4
1	E	-5	LEU	2.4
1	G	32	SER	2.4
1	G	148	ALA	2.4
1	C	32	SER	2.4
1	G	34	SER	2.4
2	H	71[A]	GLU	2.4
1	E	23	PRO	2.4
1	A	229	LEU	2.3
1	G	253	LEU	2.3
2	F	172	GLU	2.3
2	F	170	ALA	2.3
1	G	75	ASP	2.3
1	E	237	MET	2.3
1	E	226	TYR	2.3
1	E	24	VAL	2.3
1	E	172	LEU	2.3
1	C	50	ARG	2.2
2	F	233	PRO	2.2
1	C	222	LYS	2.2
1	G	265	TYR	2.2
1	G	283	ILE	2.2
1	E	37	GLN	2.2
1	C	26	ALA	2.2
1	G	266	PHE	2.2
1	G	263	ILE	2.2
1	G	68	ARG	2.2
1	G	250	ARG	2.2
1	G	259	GLY	2.2
1	C	20	ASN	2.1
1	E	165	PHE	2.1
1	E	60	VAL	2.1
1	A	332	GLY	2.1
1	G	242	PRO	2.1
2	H	92	ARG	2.1
1	A	330	ALA	2.1
1	E	26	ALA	2.1
2	H	278	ALA	2.1
2	H	119	TYR	2.1

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Mol	Chain	Res	Type	RSRZ
1	G	188	ASP	2.1
1	A	34	SER	2.1
1	G	4	VAL	2.1
1	G	345	VAL	2.1
1	C	37	GLN	2.1
1	C	203	PHE	2.1
1	G	318	ILE	2.1
1	A	250	ARG	2.1
1	G	189	SER	2.1
1	G	284	GLY	2.1
1	G	279	TRP	2.1
1	G	190	CYS	2.1
1	E	27	ALA	2.1
1	G	280	ALA	2.1
1	E	47	PHE	2.0
2	F	138	GLU	2.0
1	G	40	GLY	2.0
2	F	230	GLN	2.0
1	E	28	VAL	2.0
1	G	348	CYS	2.0
2	F	277	LEU	2.0
1	A	329	ALA	2.0
1	E	16	ARG	2.0
2	F	120	ARG	2.0
2	H	170	ALA	2.0
1	C	29	PRO	2.0
1	G	16	ARG	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,

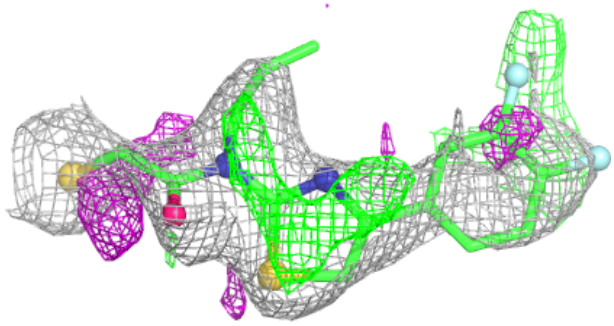
median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

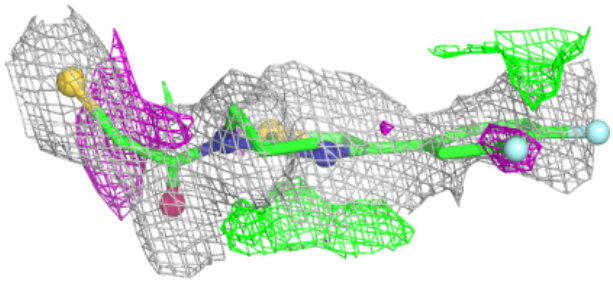
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	A1J61	E	401	21/21	0.83	0.22	65,85,102,113	0
3	A1J61	G	401	21/21	0.84	0.22	66,84,90,93	0
3	A1J61	C	401	21/21	0.90	0.17	54,69,90,96	0
4	CL	C	402	1/1	0.90	0.20	67,67,67,67	0
4	CL	A	402	1/1	0.91	0.16	66,66,66,66	0
3	A1J61	A	401	21/21	0.91	0.15	45,60,71,78	0

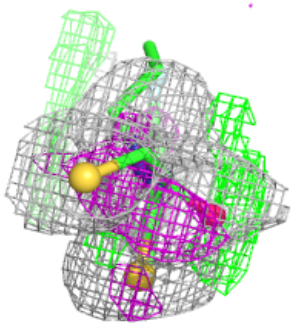
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around A1J61 E 401:

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

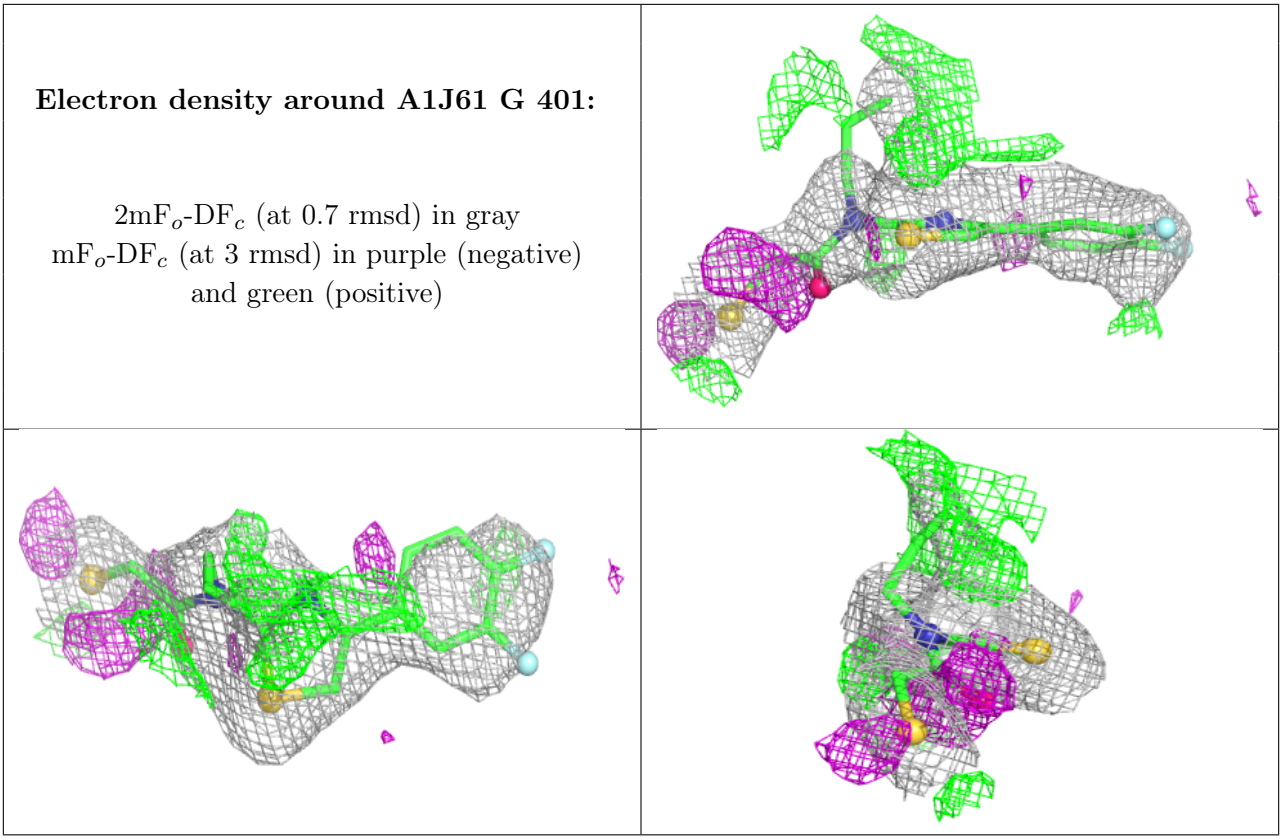






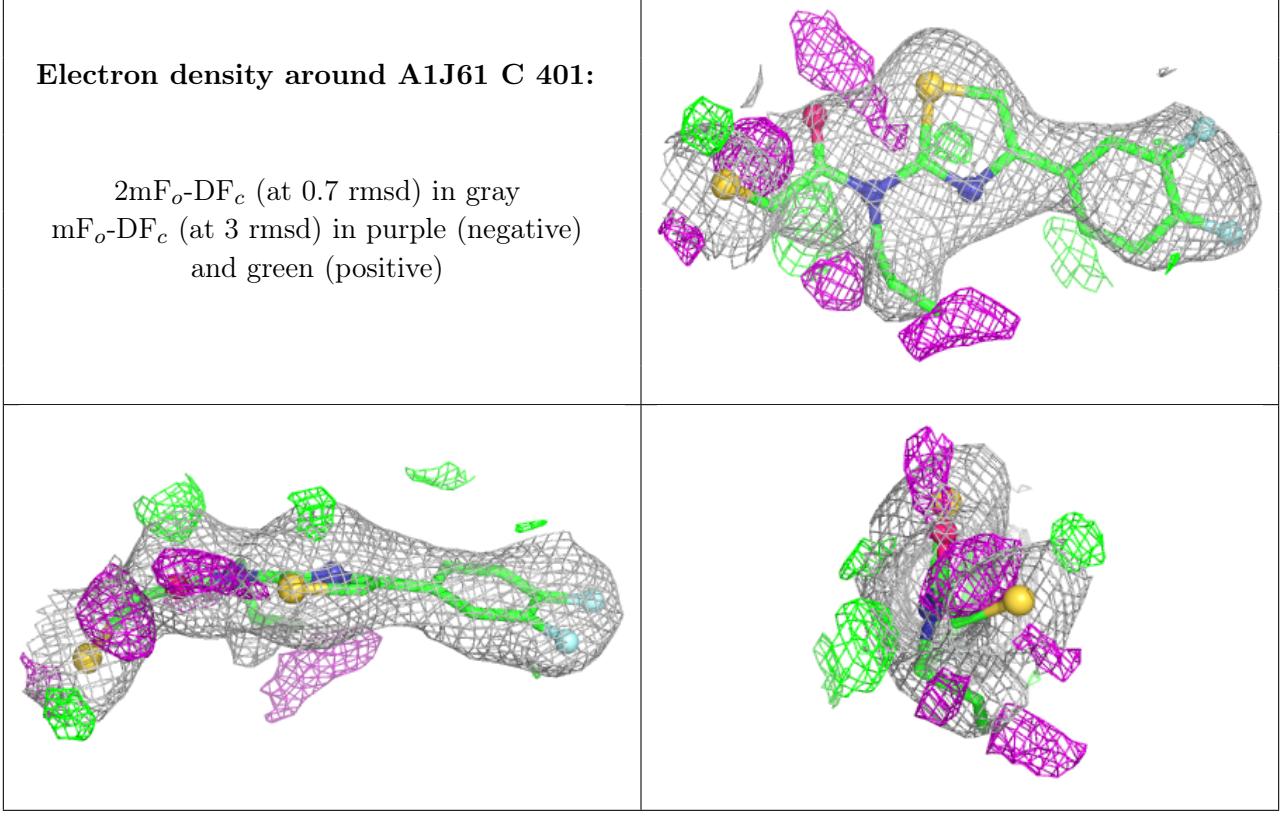
Electron density around A1J61 G 401:

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



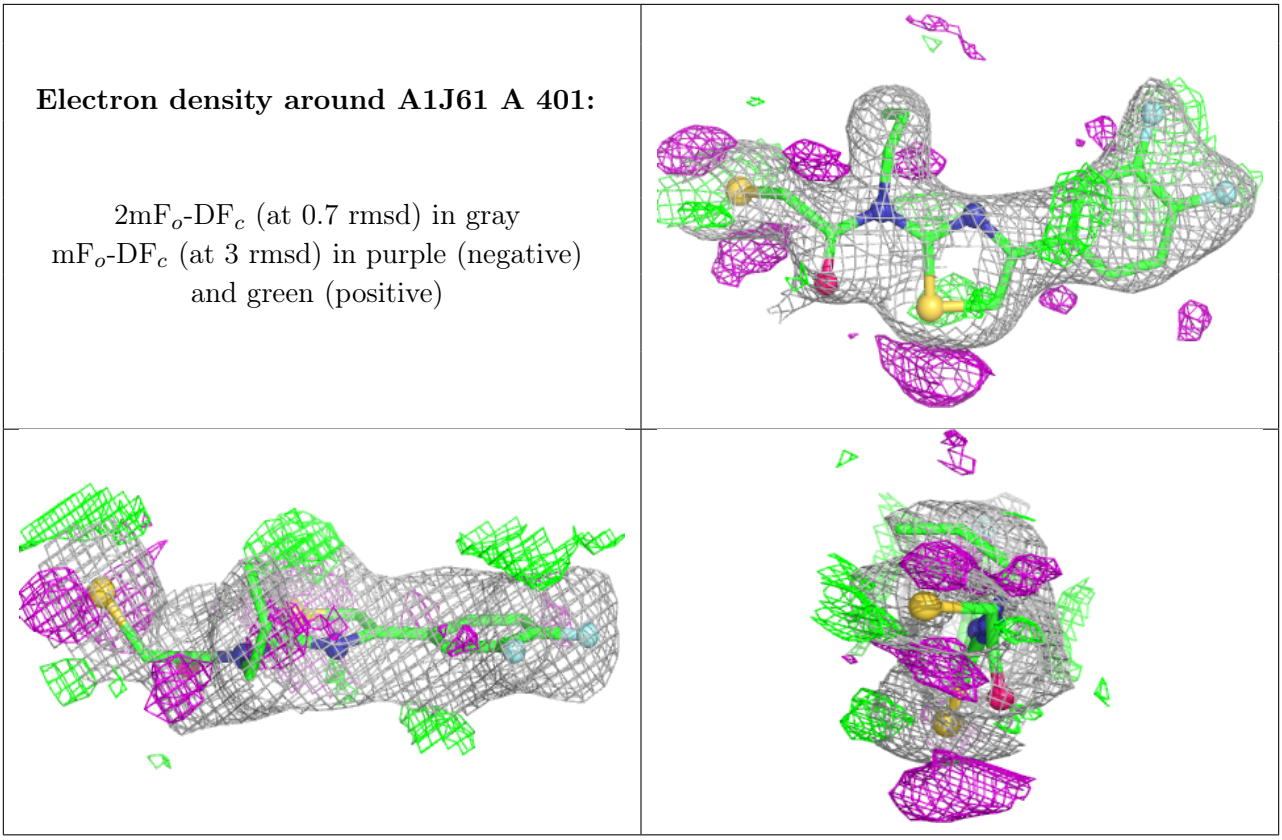
Electron density around A1J61 C 401:

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



Electron density around A1J61 A 401:

$2mF_o$-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.