Full wwPDB X-ray Structure Validation Report (i)

Nov 2, 2026 – 03:27 pm GMT

PDB ID	:	30EG / pdb_000030eg
Title	:	2-heptyl-4(1H)-quinolone synthase PqsBC in complex with D3
Authors	:	Borgert, S.R.; Blankenfeldt, W.
Deposited on	:	2026-04-21
Resolution	:	2.24 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:

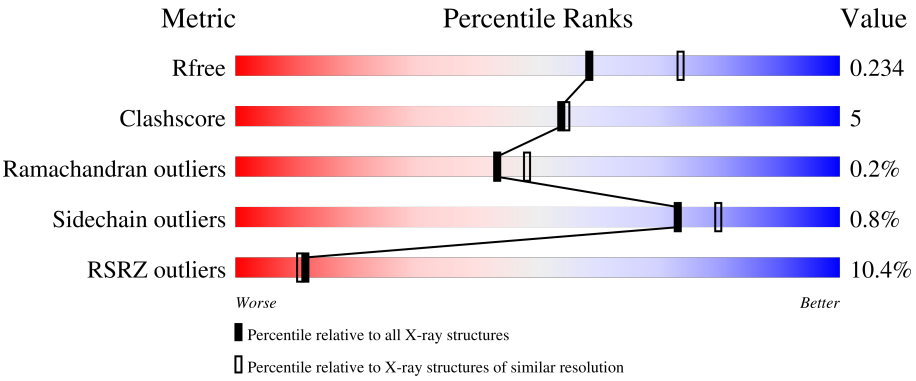
MolProbity	:	4-5-2 with Phenix2.2
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	2.2
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal 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 2.24 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.



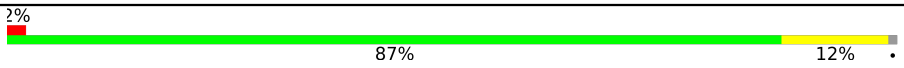
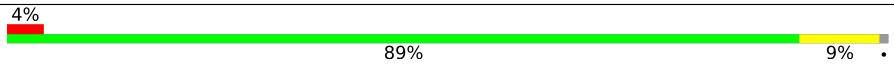
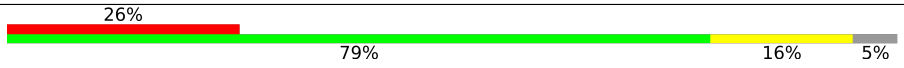
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3416 (2.26-2.22)
Clashscore	190562	3556 (2.26-2.22)
Ramachandran outliers	187476	3500 (2.26-2.22)
Sidechain outliers	187428	3501 (2.26-2.22)
RSRZ outliers	180081	3415 (2.26-2.22)

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	7% 88% 9% .
1	C	365	17% 81% 15% .
1	E	365	9% 85% 10% 5%
2	B	283	2% 90% 8% .
2	D	283	8% 88% 11% .

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Mol	Chain	Length	Quality of chain
2	F	283	
2	H	283	
3	G	365	

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 19752 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 2-heptyl-4(1H)-quinolone synthase subunit PqsC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	354	Total	C	N	O	S	0	1	0
			2728	1727	472	510	19			
1	C	350	Total	C	N	O	S	0	1	0
			2653	1683	450	501	19			
1	E	348	Total	C	N	O	S	0	0	0
			2666	1693	456	499	18			

There are 51 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
C	-16	MET	-	initiating methionine	UNP Q9I4X1
C	-15	GLY	-	expression tag	UNP Q9I4X1
C	-14	HIS	-	expression tag	UNP Q9I4X1
C	-13	HIS	-	expression tag	UNP Q9I4X1
C	-12	HIS	-	expression tag	UNP Q9I4X1
C	-11	HIS	-	expression tag	UNP Q9I4X1

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Chain	Residue	Modelled	Actual	Comment	Reference
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
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

- 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	0	0
			2117	1335	378	396	8			
2	D	279	Total	C	N	O	S	0	0	0
			2107	1330	375	394	8			
2	F	279	Total	C	N	O	S	0	0	0
			2117	1335	378	396	8			
2	H	279	Total	C	N	O	S	0	0	0
			2112	1333	376	395	8			

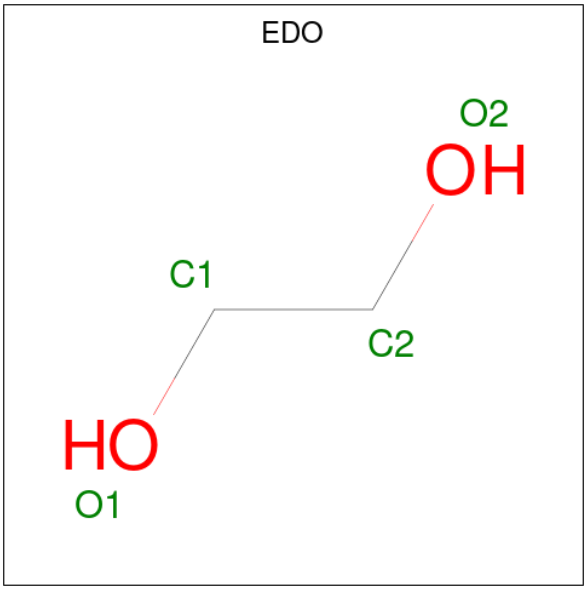
- Molecule 3 is a protein called 2-heptyl-4(1H)-quinolone synthase subunit PqsC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	G	348	Total	C	N	O	S	0	0	0
			2655	1686	455	495	19			

There are 17 discrepancies between the modelled and reference sequences:

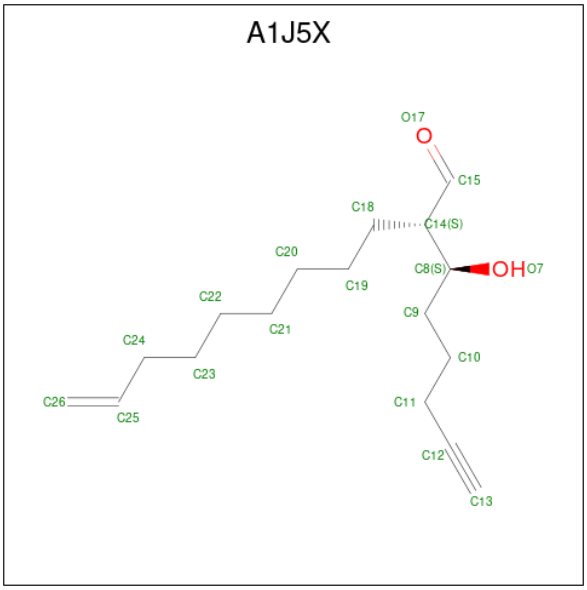
Chain	Residue	Modelled	Actual	Comment	Reference
G	-16	MET	-	initiating methionine	UNP Q9I4X1
G	-15	GLY	-	expression tag	UNP Q9I4X1
G	-14	HIS	-	expression tag	UNP Q9I4X1
G	-13	HIS	-	expression tag	UNP Q9I4X1
G	-12	HIS	-	expression tag	UNP Q9I4X1
G	-11	HIS	-	expression tag	UNP Q9I4X1
G	-10	HIS	-	expression tag	UNP Q9I4X1
G	-9	HIS	-	expression tag	UNP Q9I4X1
G	-8	ALA	-	expression tag	UNP Q9I4X1
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

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 4 2 2	0	0
4	C	1	Total C O 4 2 2	0	0

- Molecule 5 is (2 {S})-2-[(1 {S})-1-oxidanylhex-5-ynyl]undec-10-enal (CCD ID: A1J5X) (formula: $C_{17}H_{28}O_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 19 17 2	0	0
5	C	1	Total C O 19 17 2	0	0
5	E	1	Total C O 19 17 2	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	124	Total O 124 124	0	0
6	B	120	Total O 120 120	0	0
6	C	21	Total O 21 21	0	0
6	D	35	Total O 35 35	0	0

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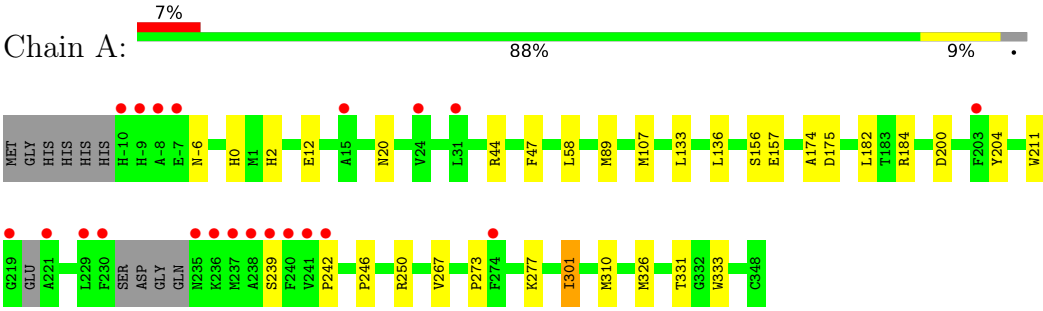
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	E	54	Total O 54 54	0	0
6	F	98	Total O 98 98	0	0
6	G	11	Total O 11 11	0	0
6	H	69	Total O 69 69	0	0

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.

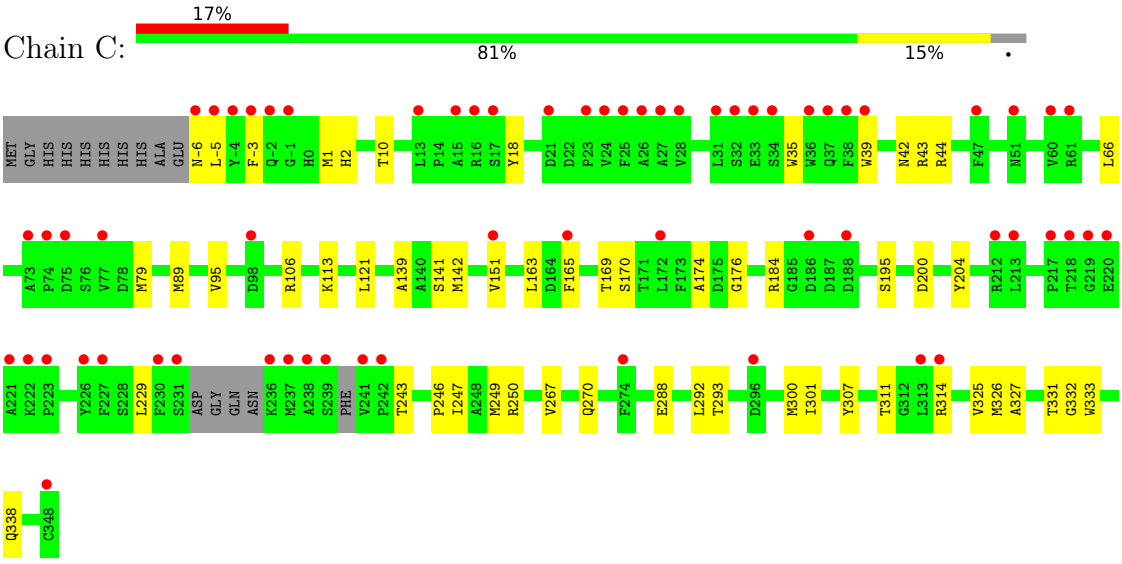
• Molecule 1: 2-heptyl-4(1H)-quinolone synthase subunit PqsC

Chain A:



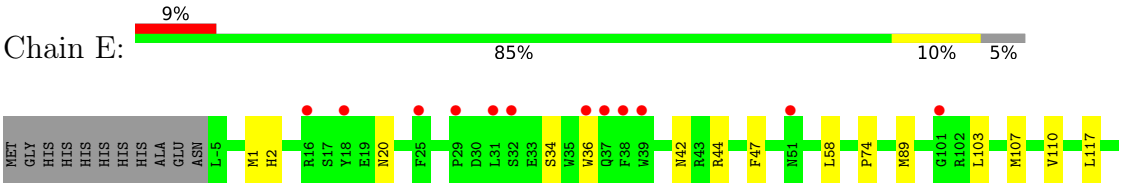
• Molecule 1: 2-heptyl-4(1H)-quinolone synthase subunit PqsC

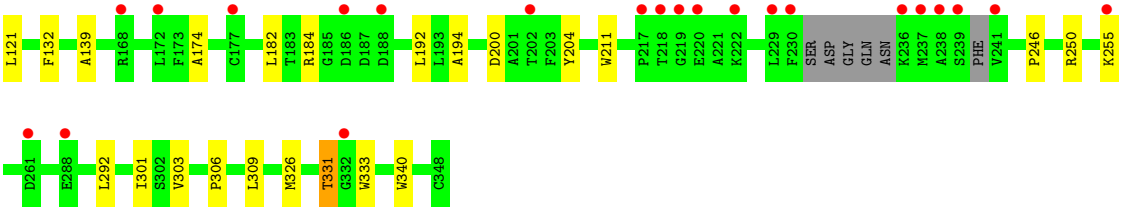
Chain C:



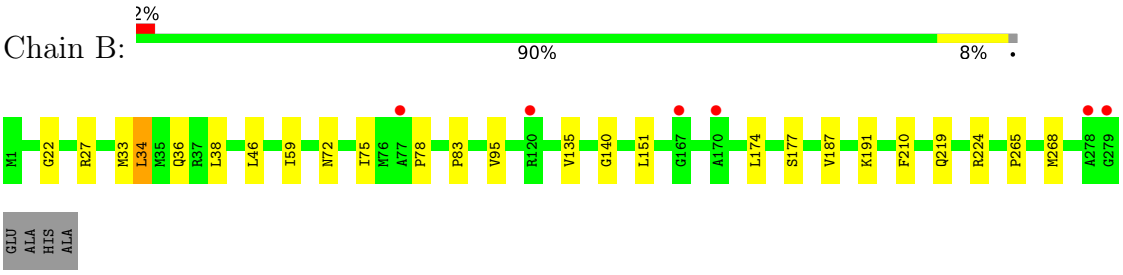
• Molecule 1: 2-heptyl-4(1H)-quinolone synthase subunit PqsC

Chain E:

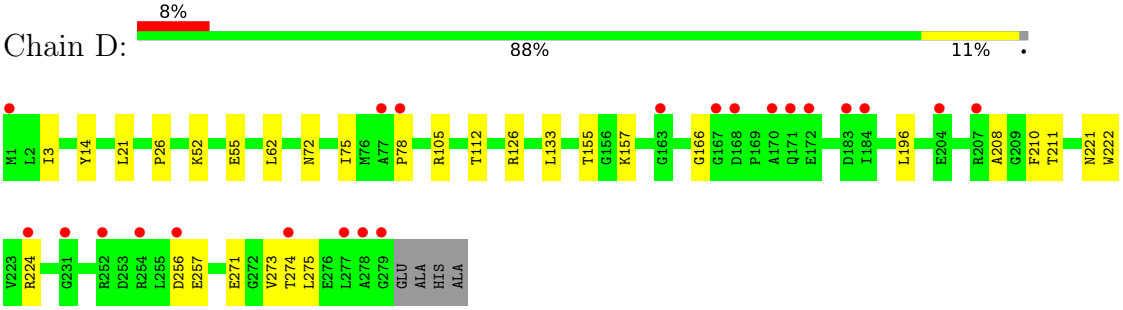




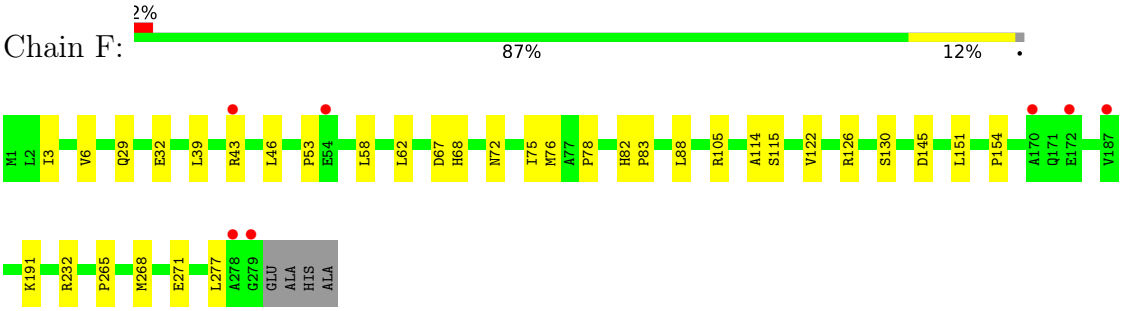
• Molecule 2: 2-heptyl-4(1H)-quinolone synthase subunit PqsB



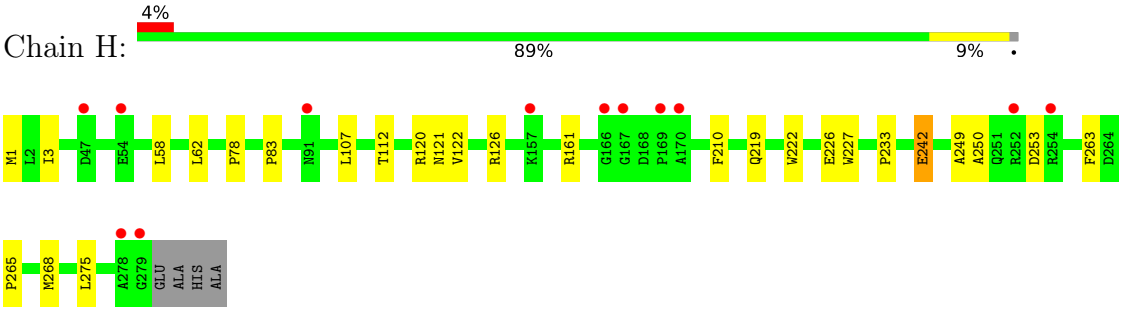
• Molecule 2: 2-heptyl-4(1H)-quinolone synthase subunit PqsB



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

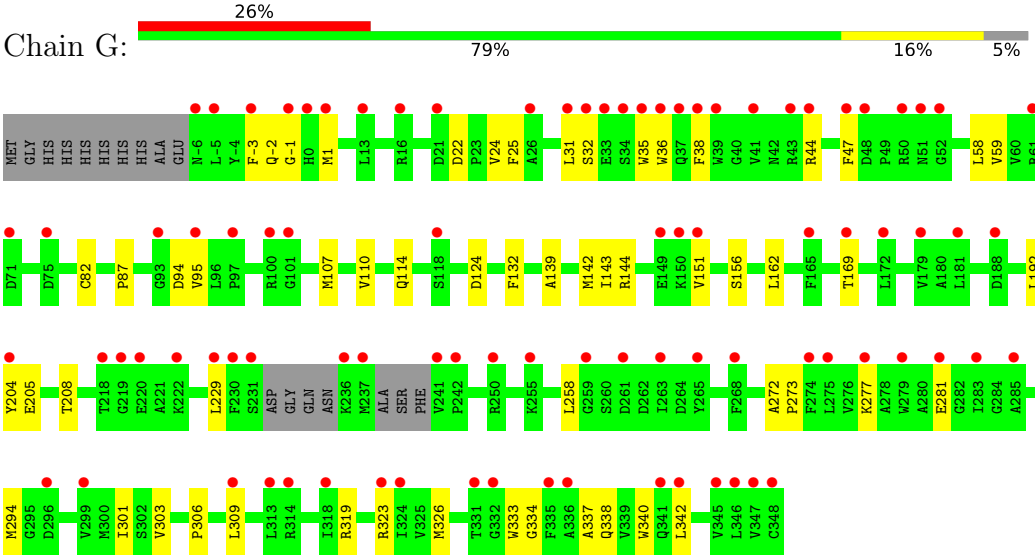


• Molecule 2: 2-heptyl-4(1H)-quinolone synthase subunit PqsB



• Molecule 3: 2-heptyl-4(1H)-quinolone synthase subunit PqsC

Chain G:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	158.83Å 226.07Å 135.05Å 90.00° 112.06° 90.00°	Depositor
Resolution (Å)	76.62 – 2.24 76.62 – 2.24	Depositor EDS
% Data completeness (in resolution range)	62.8 (76.62-2.24) 63.0 (76.62-2.24)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.43 (at 1.86Å)	Xtriage
Refinement program	PHENIX ("2.0_5936": ???)	Depositor
R, R_{free}	0.205 , 0.238 0.203 , 0.234	Depositor DCC
R_{free} test set	7428 reflections (2.01%)	wwPDB-VP
Wilson B-factor (Å^2)	23.7	Xtriage
Anisotropy	0.087	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 49.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	19752	wwPDB-VP
Average B, all atoms (Å^2)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.59% 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: A1J5X, CSX, EDO

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	0.27	0/2791	0.57	0/3788
1	C	0.25	0/2711	0.58	1/3686 (0.0%)
1	E	0.26	0/2725	0.56	0/3700
2	B	0.25	0/2158	0.57	0/2932
2	D	0.26	0/2148	0.59	0/2920
2	F	0.24	0/2158	0.52	0/2932
2	H	0.23	0/2153	0.52	0/2926
3	G	0.28	0/2706	0.60	0/3675
All	All	0.26	0/19550	0.57	1/26559 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	332	GLY	N-CA-C	-6.31	107.08	115.40

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	2728	0	2673	27	0
1	C	2653	0	2569	38	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	2666	0	2615	27	0
2	B	2117	0	2126	19	0
2	D	2107	0	2111	23	0
2	F	2117	0	2126	21	0
2	H	2112	0	2115	21	0
3	G	2655	0	2588	41	0
4	A	4	0	6	0	0
4	C	4	0	6	0	0
5	A	19	0	0	0	0
5	C	19	0	0	0	0
5	E	19	0	0	0	0
6	A	124	0	0	0	0
6	B	120	0	0	0	0
6	C	21	0	0	0	0
6	D	35	0	0	0	0
6	E	54	0	0	0	0
6	F	98	0	0	1	0
6	G	11	0	0	1	0
6	H	69	0	0	0	0
All	All	19752	0	18935	186	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 5.

All (186) 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:139:ALA:HB1	1:C:151:VAL:HG21	1.64	0.76
1:A:204:TYR:HE1	2:B:78:PRO:HB2	1.52	0.75
1:E:121:LEU:HD11	2:F:105:ARG:HD2	1.73	0.71
1:C:121:LEU:HD11	2:D:105:ARG:HD2	1.73	0.71
2:F:62:LEU:HB2	2:F:126:ARG:HG3	1.73	0.70
2:B:27:ARG:HH12	2:B:135:VAL:HG11	1.56	0.70
3:G:114:GLN:HB3	6:G:401:HOH:O	1.97	0.64
1:C:267:VAL:HB	1:C:326:MET:HG2	1.80	0.63
2:H:62:LEU:HB2	2:H:126:ARG:HG3	1.81	0.63
1:E:47:PHE:HE1	1:E:58:LEU:HD12	1.64	0.62
1:A:204:TYR:CE1	2:B:78:PRO:HB2	2.35	0.62
1:C:-5:LEU:HD13	2:H:249:ALA:HB1	1.81	0.62
2:H:226:GLU:HB3	2:H:242:GLU:OE2	1.99	0.62
1:A:12:GLU:HG2	1:A:58:LEU:HB3	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:79:MET:HB3	1:C:151:VAL:HG12	1.81	0.61
1:C:270:GLN:NE2	1:C:293:THR:H	1.99	0.60
3:G:169:THR:HG21	3:G:229:LEU:HD23	1.83	0.60
2:D:210:PHE:HD2	2:D:224:ARG:HD2	1.66	0.60
1:C:79:MET:HE1	1:C:142:MET:SD	2.40	0.60
2:D:273:VAL:HG13	2:D:275:LEU:HD11	1.83	0.59
3:G:87:PRO:HD2	3:G:162:LEU:HD21	1.85	0.59
2:D:155:THR:HB	2:D:157:LYS:HE2	1.84	0.59
2:F:72:ASN:HB3	2:F:75:ILE:HD12	1.85	0.59
1:A:200:ASP:O	1:A:333:TRP:HB2	2.02	0.58
3:G:132:PHE:CG	3:G:306:PRO:HG3	2.38	0.58
1:E:20:ASN:HD21	1:E:42:ASN:HA	1.66	0.57
1:E:44:ARG:HB2	1:E:174:ALA:HB2	1.84	0.57
1:C:200:ASP:O	1:C:333:TRP:HB2	2.04	0.57
2:F:105:ARG:HD3	2:F:271:GLU:OE2	2.05	0.57
2:B:33:MET:HE3	2:B:36:GLN:HG3	1.86	0.56
1:E:200:ASP:O	1:E:333:TRP:HB2	2.05	0.56
1:A:47:PHE:HE1	1:A:58:LEU:HD12	1.70	0.55
1:C:204:TYR:HE1	2:D:78:PRO:HB2	1.72	0.55
1:C:331:THR:C	1:C:333:TRP:H	2.14	0.55
1:C:1:MET:HE1	2:H:219:GLN:HG2	1.87	0.55
2:B:46:LEU:HD23	2:B:151:LEU:HD22	1.88	0.55
1:E:132:PHE:CG	1:E:306:PRO:HG3	2.42	0.55
2:F:130:SER:HB3	2:F:145:ASP:OD2	2.06	0.54
1:E:47:PHE:CE1	1:E:58:LEU:HD12	2.42	0.54
2:B:72:ASN:HB3	2:B:75:ILE:HD12	1.90	0.53
2:B:219:GLN:HG2	1:E:1:MET:HE1	1.88	0.53
1:E:246:PRO:O	1:E:250:ARG:HG3	2.08	0.53
2:B:265:PRO:HA	2:B:268:MET:HE2	1.90	0.53
1:E:20:ASN:ND2	1:E:42:ASN:HA	2.24	0.53
2:B:174:LEU:HD12	2:H:233:PRO:HB3	1.91	0.53
1:C:18:TYR:HB2	1:C:44:ARG:HD2	1.89	0.53
1:C:10:THR:HG22	1:C:66:LEU:HD12	1.90	0.52
1:C:249:MET:HE3	1:C:327:ALA:HB3	1.91	0.52
1:C:311:THR:HA	1:C:314:ARG:HG2	1.91	0.52
2:D:210:PHE:HB3	2:D:222:TRP:CZ3	2.44	0.52
1:C:44:ARG:HB2	1:C:174:ALA:HB2	1.92	0.52
2:D:105:ARG:HD3	2:D:271:GLU:OE2	2.09	0.52
2:D:221:ASN:HD22	3:G:-2:GLN:HA	1.74	0.52
1:C:-3:PHE:HE2	2:H:253:ASP:OD2	1.91	0.52
1:A:47:PHE:CE1	1:A:58:LEU:HD12	2.45	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:331:THR:C	1:E:333:TRP:H	2.17	0.51
2:D:221:ASN:HD22	3:G:-1:GLY:H	1.59	0.51
3:G:31:LEU:HD13	3:G:36:TRP:CZ2	2.46	0.51
3:G:277:LYS:O	3:G:281:GLU:HG3	2.11	0.51
1:A:267:VAL:HB	1:A:326:MET:HG2	1.93	0.50
1:C:204:TYR:CE1	2:D:78:PRO:HB2	2.47	0.50
3:G:47:PHE:CE1	3:G:58:LEU:HD12	2.46	0.50
1:E:309:LEU:HD22	1:E:326:MET:HE1	1.93	0.50
1:A:331:THR:C	1:A:333:TRP:H	2.19	0.50
2:D:72:ASN:HB3	2:D:75:ILE:HD12	1.93	0.50
2:H:58:LEU:HB3	2:H:122:VAL:HG22	1.92	0.50
3:G:143:ILE:HD11	3:G:151:VAL:HG23	1.94	0.50
1:E:20:ASN:HB3	1:E:36:TRP:CH2	2.47	0.49
2:F:265:PRO:HA	2:F:268:MET:HE2	1.93	0.49
3:G:94:ASP:HB2	3:G:114:GLN:HE22	1.77	0.49
3:G:25:PHE:HZ	3:G:44:ARG:HD2	1.77	0.49
2:B:210:PHE:CD2	2:B:224:ARG:HD2	2.47	0.49
2:F:3:ILE:HB	2:F:277:LEU:HD23	1.95	0.49
2:F:46:LEU:HD23	2:F:151:LEU:HD22	1.95	0.48
1:C:42:ASN:HB3	1:C:43:ARG:HH11	1.78	0.48
2:H:3:ILE:HD11	2:H:275:LEU:HD13	1.95	0.48
1:C:270:GLN:HE22	1:C:292:LEU:HA	1.78	0.48
3:G:31:LEU:HD13	3:G:36:TRP:HZ2	1.79	0.48
1:A:211:TRP:CG	2:B:191:LYS:HE2	2.48	0.48
2:D:256:ASP:OD2	2:D:257:GLU:OE2	2.32	0.47
2:F:67:ASP:OD1	2:F:68:HIS:HD2	1.96	0.47
1:A:107:MET:HE1	1:A:156:SER:HB3	1.96	0.47
1:C:246:PRO:O	1:C:250:ARG:HG3	2.14	0.47
2:B:27:ARG:HH22	2:B:135:VAL:HG12	1.80	0.47
2:D:3:ILE:HG13	2:D:275:LEU:HD23	1.95	0.47
1:E:103:LEU:HD22	2:F:76:MET:HE1	1.96	0.47
2:D:62:LEU:HB2	2:D:126:ARG:HD3	1.96	0.47
1:C:141:SER:HB3	2:D:112:THR:HG22	1.96	0.47
2:D:166:GLY:HA3	2:D:208:ALA:HB3	1.95	0.47
2:H:210:PHE:HB3	2:H:222:TRP:CZ3	2.50	0.47
1:A:89:MET:HE2	1:A:89:MET:HB3	1.80	0.47
2:D:210:PHE:CD2	2:D:224:ARG:HD2	2.48	0.47
2:F:53:PRO:HB3	2:F:88:LEU:HG	1.96	0.47
1:A:-6:ASN:HD21	2:F:232:ARG:NH2	2.13	0.46
1:C:89:MET:HE3	1:C:106:ARG:CZ	2.45	0.46
1:E:44:ARG:HH11	1:E:44:ARG:HG2	1.79	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:ARG:HH11	1:A:44:ARG:HG2	1.79	0.46
1:C:-3:PHE:CE2	2:H:253:ASP:OD2	2.68	0.46
3:G:200:ASP:O	3:G:333:TRP:HB2	2.15	0.46
3:G:192:LEU:HD11	3:G:338:GLN:HB3	1.97	0.46
2:F:115:SER:HA	6:F:302:HOH:O	2.14	0.46
3:G:319:ARG:HE	3:G:319:ARG:HB2	1.57	0.46
2:D:3:ILE:CG1	2:D:275:LEU:HD23	2.46	0.46
1:A:310:MET:HE2	1:A:310:MET:HB3	1.90	0.45
1:E:204:TYR:CE2	2:F:78:PRO:HD2	2.51	0.45
2:F:114:ALA:HB1	2:F:154:PRO:HG3	1.98	0.45
1:E:204:TYR:CE1	2:F:83:PRO:HG3	2.51	0.45
3:G:1:MET:HB3	3:G:144:ARG:HB3	1.99	0.45
1:C:163:LEU:HD22	1:C:170:SER:HA	1.99	0.45
3:G:107:MET:HE1	3:G:156:SER:HB3	1.99	0.45
3:G:38:PHE:O	3:G:273:PRO:HD2	2.17	0.45
1:A:204:TYR:CZ	2:B:83:PRO:HG3	2.51	0.45
1:C:176:GLY:HA3	1:C:300:MET:CE	2.47	0.45
1:C:307:TYR:O	1:C:311:THR:HG23	2.16	0.45
2:H:1:MET:HE3	2:H:107:LEU:HD22	1.98	0.45
3:G:326:MET:O	3:G:337:ALA:HA	2.17	0.45
2:B:34:LEU:HD22	2:B:38:LEU:HB2	1.99	0.44
3:G:142:MET:HE2	2:H:112:THR:HG21	1.99	0.44
3:G:169:THR:HG21	3:G:229:LEU:CD2	2.45	0.44
1:E:211:TRP:CG	2:F:191:LYS:HE2	2.53	0.44
2:F:58:LEU:HB3	2:F:122:VAL:HG22	1.99	0.44
3:G:258:LEU:HD21	3:G:323:ARG:HD2	1.99	0.44
3:G:272:ALA:HA	3:G:294:MET:HE1	2.00	0.44
2:D:21:LEU:HD13	2:D:196:LEU:HD22	1.98	0.44
1:C:195:SER:HB2	1:C:338:GLN:HG2	2.00	0.44
3:G:309:LEU:HD22	3:G:326:MET:HE1	2.00	0.44
1:E:74:PRO:HB3	1:E:117:LEU:HD12	2.00	0.43
2:D:52:LYS:HD2	2:D:55:GLU:CD	2.43	0.43
1:A:12:GLU:HG3	1:A:58:LEU:HD13	1.99	0.43
3:G:32:SER:O	3:G:35:TRP:HB3	2.18	0.43
3:G:204:TYR:CZ	2:H:83:PRO:HG3	2.54	0.43
3:G:303:VAL:C	3:G:306:PRO:HD2	2.43	0.43
2:F:29:GLN:N	2:F:32:GLU:OE2	2.44	0.43
3:G:192:LEU:HD13	3:G:340:TRP:CE2	2.54	0.43
1:A:239:SER:C	1:A:242:PRO:HD2	2.44	0.43
1:E:2:HIS:HB3	1:E:184:ARG:HG3	1.99	0.43
1:A:204:TYR:CE2	2:B:83:PRO:HG3	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:204:TYR:CE1	2:H:83:PRO:HG3	2.54	0.43
1:C:35:TRP:NE1	1:C:39:TRP:HE1	2.17	0.43
1:C:95:VAL:HG21	1:C:113:LYS:HB2	2.00	0.43
1:C:249:MET:HE2	1:C:325:VAL:CG1	2.49	0.43
1:E:194:ALA:HB1	1:E:255:LYS:HB3	2.01	0.43
3:G:205:GLU:H	3:G:205:GLU:CD	2.24	0.42
2:H:120:ARG:C	2:H:121:ASN:HD22	2.27	0.42
1:C:2:HIS:HB3	1:C:184:ARG:HG3	2.00	0.42
1:E:303:VAL:C	1:E:306:PRO:HD2	2.45	0.42
3:G:139:ALA:O	3:G:143:ILE:HG12	2.20	0.42
3:G:82:CYS:HB2	3:G:124:ASP:OD1	2.19	0.42
2:D:274:THR:C	2:D:275:LEU:HD12	2.45	0.42
3:G:95:VAL:HB	3:G:110:VAL:HG23	2.00	0.42
2:B:59:ILE:HB	2:B:95:VAL:HG22	2.01	0.42
2:F:39:LEU:HD21	2:F:43:ARG:CZ	2.50	0.42
1:A:20:ASN:HD22	1:A:20:ASN:N	2.17	0.42
1:C:165:PHE:CD2	1:C:170:SER:HB2	2.54	0.42
1:C:169:THR:HG22	1:C:229:LEU:HD12	2.02	0.42
3:G:59:VAL:HG21	3:G:156:SER:HB2	2.01	0.42
3:G:132:PHE:CD2	3:G:306:PRO:HG3	2.55	0.42
1:A:273:PRO:O	1:A:277:LYS:HG3	2.20	0.41
2:H:161:ARG:HE	2:H:161:ARG:HB2	1.65	0.41
2:B:22:GLY:HA2	2:B:140:GLY:O	2.21	0.41
1:E:192:LEU:HD13	1:E:340:TRP:CE2	2.55	0.41
3:G:-3:PHE:CD1	3:G:-3:PHE:N	2.88	0.41
3:G:208:THR:HA	2:H:78:PRO:HD3	2.02	0.41
1:E:139:ALA:HB3	1:E:182:LEU:HD12	2.02	0.41
1:E:292:LEU:HD23	1:E:292:LEU:HA	1.93	0.41
3:G:22:ASP:HB3	3:G:24:VAL:HG12	2.02	0.41
2:H:265:PRO:HA	2:H:268:MET:HE2	2.02	0.41
3:G:199:SER:HA	3:G:334:GLY:HA2	2.02	0.41
2:H:227:TRP:CD1	2:H:263:PHE:HD2	2.39	0.41
1:C:-3:PHE:CZ	2:H:250:ALA:HA	2.56	0.41
1:A:211:TRP:CD2	2:B:191:LYS:HE2	2.56	0.41
1:C:243:THR:O	1:C:247:ILE:HG13	2.21	0.41
2:B:27:ARG:HH22	2:B:135:VAL:CG1	2.34	0.41
1:A:157:GLU:HG3	1:A:301:ILE:HB	2.02	0.40
1:A:2:HIS:HB3	1:A:184:ARG:HG3	2.02	0.40
1:A:246:PRO:O	1:A:250:ARG:HG3	2.22	0.40
2:F:82:HIS:HB2	2:F:83:PRO:HD3	2.04	0.40
1:E:89:MET:HE1	1:E:107:MET:HB2	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:107:MET:HE3	1:E:107:MET:HB2	2.01	0.40
1:A:47:PHE:CE2	1:A:175:ASP:HB3	2.57	0.40
1:C:-5:LEU:HD13	2:H:249:ALA:CB	2.48	0.40
2:D:221:ASN:ND2	3:G:-2:GLN:HA	2.36	0.40
1:A:44:ARG:HB2	1:A:174:ALA:HB2	2.03	0.40
1:A:136:LEU:HD22	1:A:182:LEU:HD11	2.02	0.40
1:C:-6:ASN:HB3	1:C:-5:LEU:H	1.70	0.40
2:D:14:TYR:CE2	2:D:26:PRO:HG3	2.56	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	349/365 (96%)	336 (96%)	12 (3%)	1 (0%)	36	38
1	C	345/365 (94%)	330 (96%)	14 (4%)	1 (0%)	36	38
1	E	342/365 (94%)	332 (97%)	9 (3%)	1 (0%)	36	38
2	B	277/283 (98%)	272 (98%)	5 (2%)	0	100	100
2	D	277/283 (98%)	273 (99%)	4 (1%)	0	100	100
2	F	277/283 (98%)	273 (99%)	4 (1%)	0	100	100
2	H	277/283 (98%)	270 (98%)	7 (2%)	0	100	100
3	G	341/365 (93%)	331 (97%)	9 (3%)	1 (0%)	36	38
All	All	2485/2592 (96%)	2417 (97%)	64 (3%)	4 (0%)	43	48

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	301	ILE
1	C	301	ILE

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Mol	Chain	Res	Type
1	E	301	ILE
3	G	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	287/300 (96%)	283 (99%)	4 (1%)	59	69
1	C	276/300 (92%)	275 (100%)	1 (0%)	84	88
1	E	280/300 (93%)	277 (99%)	3 (1%)	65	74
2	B	221/223 (99%)	218 (99%)	3 (1%)	59	69
2	D	219/223 (98%)	217 (99%)	2 (1%)	70	78
2	F	221/223 (99%)	220 (100%)	1 (0%)	81	86
2	H	219/223 (98%)	218 (100%)	1 (0%)	81	86
3	G	275/299 (92%)	274 (100%)	1 (0%)	84	88
All	All	1998/2091 (96%)	1982 (99%)	16 (1%)	73	80

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	0[A]	HIS
1	A	0[B]	HIS
1	A	133	LEU
1	A	218	THR
2	B	34	LEU
2	B	177	SER
2	B	187	VAL
1	C	288	GLU
2	D	133	LEU
2	D	211	THR
1	E	34	SER
1	E	110	VAL
1	E	331	THR

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Mol	Chain	Res	Type
2	F	6	VAL
3	G	342	LEU
2	H	242	GLU

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	-6	ASN
1	A	20	ASN
1	A	37	GLN
1	A	198	HIS
1	A	291	GLN
2	B	91	ASN
1	C	270	GLN
2	D	4	GLN
2	D	121	ASN
2	D	221	ASN
1	E	-2	GLN
1	E	0	HIS
2	F	68	HIS
2	F	219	GLN
2	F	221	ASN
3	G	216	ASN
2	H	29	GLN
2	H	121	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue 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	$\#\|Z\|>2$
3	CSX	G	129	3	3,6,7	0.63	0	1,6,8	0.20	0

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	CSX	G	129	3	-	0/1/5/7	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates (i)

There are no oligosaccharides in this entry.

5.6 Ligand geometry (i)

5 ligands are 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	$\#\|Z\|>2$
5	A1J5X	A	402	1	18,18,18	0.19	0	14,19,19	0.33	0
4	EDO	A	401	-	3,3,3	0.25	0	2,2,2	0.42	0
5	A1J5X	E	401	1	18,18,18	0.35	0	14,19,19	0.44	0
4	EDO	C	401	-	3,3,3	0.26	0	2,2,2	0.30	0
5	A1J5X	C	402	1	18,18,18	0.34	0	14,19,19	0.28	0

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
5	A1J5X	A	402	1	-	5/19/20/20	-
4	EDO	A	401	-	-	0/1/1/1	-
5	A1J5X	E	401	1	-	9/19/20/20	-
4	EDO	C	401	-	-	1/1/1/1	-
5	A1J5X	C	402	1	-	9/19/20/20	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (24) torsion outliers are listed below:

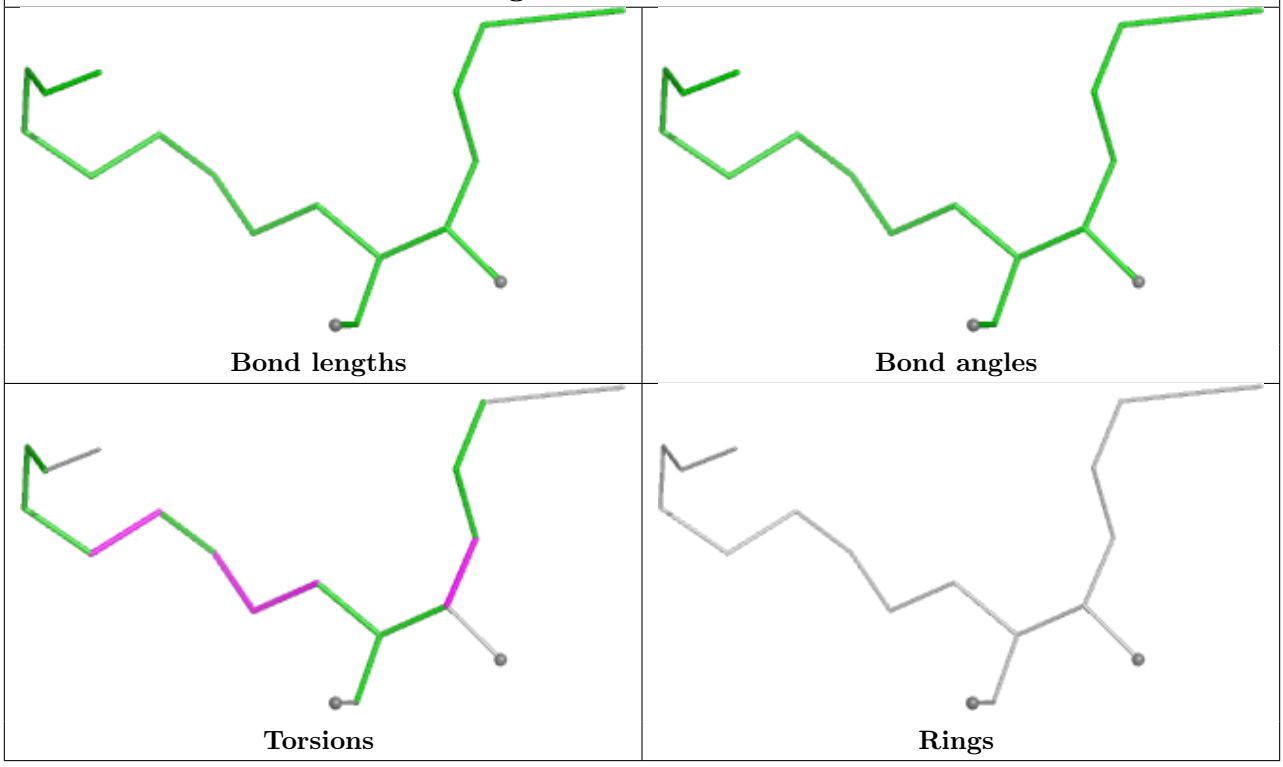
Mol	Chain	Res	Type	Atoms
5	A	402	A1J5X	C14-C8-C9-C10
5	C	402	A1J5X	C15-C14-C18-C19
5	C	402	A1J5X	C14-C8-C9-C10
5	E	401	A1J5X	C15-C14-C8-C9
5	E	401	A1J5X	C15-C14-C8-O7
5	E	401	A1J5X	C18-C14-C8-C9
5	E	401	A1J5X	C14-C8-C9-C10
5	E	401	A1J5X	C23-C24-C25-C26
5	E	401	A1J5X	O7-C8-C9-C10
5	C	402	A1J5X	C18-C19-C20-C21
5	A	402	A1J5X	O7-C8-C9-C10
5	C	402	A1J5X	O7-C8-C9-C10
5	C	402	A1J5X	C21-C22-C23-C24
5	C	402	A1J5X	C20-C21-C22-C23
5	A	402	A1J5X	C20-C21-C22-C23
5	C	402	A1J5X	C9-C10-C11-C12
5	A	402	A1J5X	C14-C18-C19-C20
5	A	402	A1J5X	C18-C19-C20-C21
5	E	401	A1J5X	C18-C19-C20-C21
4	C	401	EDO	O1-C1-C2-O2
5	C	402	A1J5X	C8-C14-C15-O17
5	E	401	A1J5X	C14-C18-C19-C20
5	E	401	A1J5X	C18-C14-C8-O7
5	C	402	A1J5X	C19-C20-C21-C22

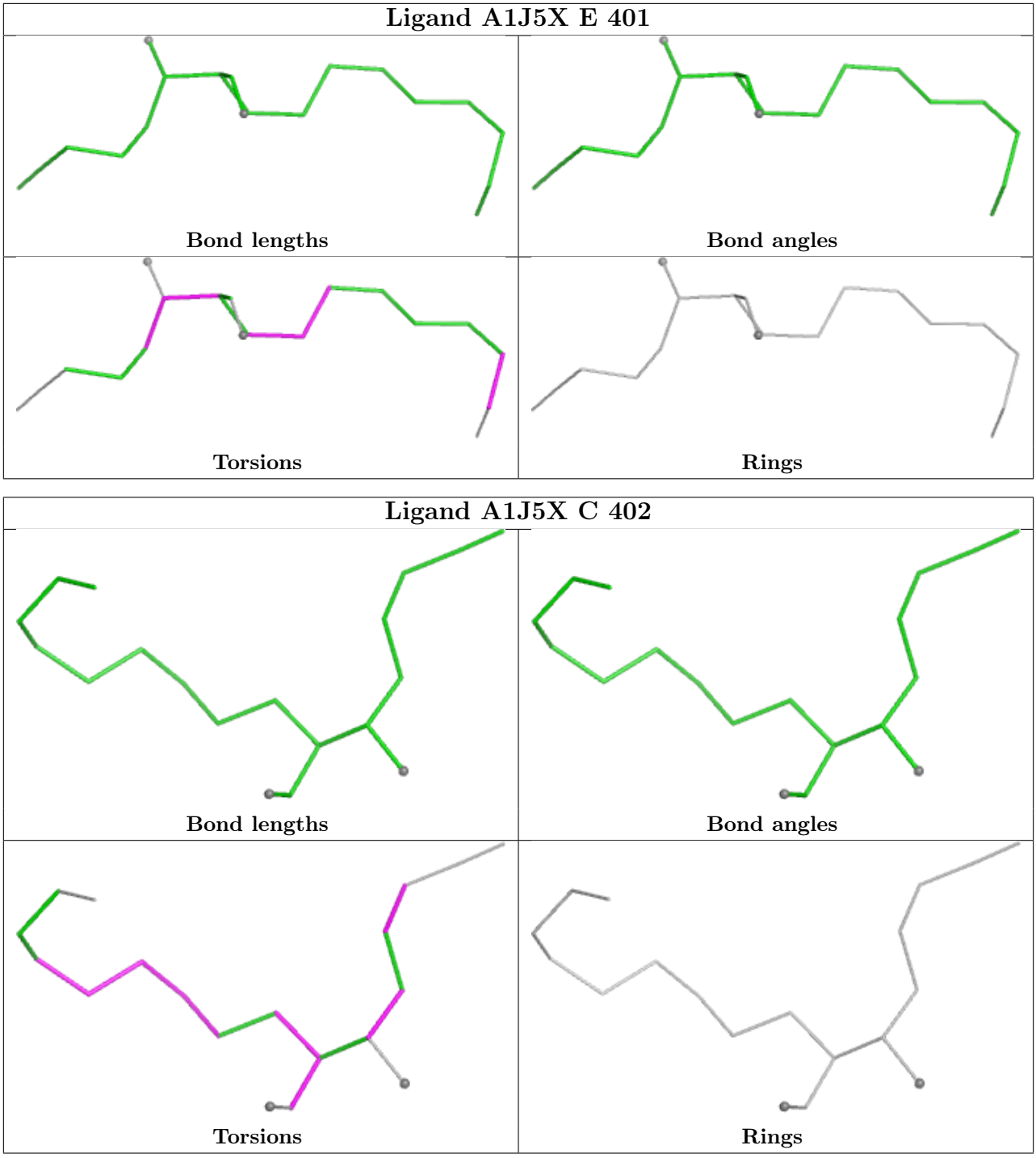
There are no ring outliers.

No monomer is involved in short contacts.

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 A1J5X A 402





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	354/365 (96%)	0.44	24 (6%) 23 21	14, 35, 76, 121	1 (0%)
1	C	350/365 (95%)	1.25	63 (18%) 3 3	34, 64, 99, 121	1 (0%)
1	E	348/365 (95%)	0.82	34 (9%) 13 12	27, 51, 79, 115	0
2	B	279/283 (98%)	-0.08	6 (2%) 62 64	11, 26, 54, 77	0
2	D	279/283 (98%)	0.72	22 (7%) 18 16	36, 51, 82, 114	0
2	F	279/283 (98%)	0.03	7 (2%) 58 59	14, 34, 60, 106	0
2	H	279/283 (98%)	0.59	12 (4%) 40 38	36, 49, 78, 107	0
3	G	347/365 (95%)	1.58	94 (27%) 1 1	54, 71, 98, 126	0
All	All	2515/2592 (97%)	0.71	262 (10%) 11 11	11, 49, 90, 126	2 (0%)

All (262) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	279	GLY	8.5
1	C	238	ALA	8.4
3	G	231	SER	7.9
3	G	218	THR	7.8
3	G	-5	LEU	7.8
2	D	77	ALA	7.4
2	H	279	GLY	7.4
1	C	231	SER	7.0
2	D	278	ALA	6.8
2	D	279	GLY	6.6
1	E	241	VAL	6.1
1	A	241	VAL	6.1
1	A	240	PHE	5.9
3	G	230	PHE	5.9
3	G	21	ASP	5.9
2	B	279	GLY	5.8

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Mol	Chain	Res	Type	RSRZ
1	A	-9	HIS	5.8
1	E	236	LYS	5.8
1	A	219	GLY	5.7
1	A	235	ASN	5.5
2	H	278	ALA	5.5
1	C	-5	LEU	5.4
2	B	77	ALA	5.3
1	A	236	LYS	5.3
1	E	218	THR	5.3
1	E	36	TRP	5.3
1	C	218	THR	5.3
3	G	275	LEU	5.2
2	H	170	ALA	5.2
1	E	238	ALA	5.1
1	C	220	GLU	5.1
3	G	241	VAL	5.1
1	E	230	PHE	5.0
1	A	238	ALA	5.0
3	G	242	PRO	5.0
3	G	274	PHE	4.9
1	C	239	SER	4.8
1	E	237	MET	4.8
2	F	278	ALA	4.8
3	G	38	PHE	4.7
3	G	237	MET	4.7
1	C	230	PHE	4.6
1	C	236	LYS	4.6
3	G	150	LYS	4.5
1	C	219	GLY	4.5
1	C	36	TRP	4.3
1	E	18	TYR	4.2
1	C	172	LEU	4.2
1	A	221	ALA	4.1
1	A	218	THR	4.1
1	C	51	ASN	4.1
1	A	242	PRO	4.1
2	F	54	GLU	4.1
1	C	-2	GLN	4.0
1	C	-3	PHE	4.0
3	G	37	GLN	4.0
1	C	242	PRO	4.0
3	G	51	ASN	4.0

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Mol	Chain	Res	Type	RSRZ
3	G	236	LYS	3.9
1	A	230	PHE	3.9
2	D	1	MET	3.9
3	G	332	GLY	3.9
3	G	31	LEU	3.8
3	G	290	TYR	3.7
2	D	172	GLU	3.7
2	D	183	ASP	3.7
3	G	32	SER	3.7
3	G	-3	PHE	3.6
1	C	-1	GLY	3.6
1	E	239	SER	3.6
1	A	239	SER	3.6
2	D	207	ARG	3.5
3	G	39	TRP	3.5
2	H	254	ARG	3.5
3	G	61	ARG	3.5
2	D	168	ASP	3.4
3	G	34	SER	3.4
1	A	237	MET	3.4
2	H	169	PRO	3.4
3	G	331	THR	3.4
2	H	54	GLU	3.4
2	D	252	ARG	3.4
3	G	261	ASP	3.3
1	C	32	SER	3.3
1	E	51	ASN	3.3
1	E	172	LEU	3.3
1	C	221	ALA	3.3
3	G	323	ARG	3.3
3	G	220	GLU	3.3
1	C	61	ARG	3.2
1	C	348	CYS	3.2
2	H	252	ARG	3.2
3	G	101	GLY	3.2
1	E	222	LYS	3.1
1	E	32	SER	3.1
3	G	348	CYS	3.1
1	E	261	ASP	3.1
1	E	229	LEU	3.1
3	G	346	LEU	3.1
1	C	188	ASP	3.1

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Mol	Chain	Res	Type	RSRZ
3	G	293	THR	3.0
1	C	37	GLN	3.0
1	E	288	GLU	3.0
1	C	-4	TYR	3.0
1	C	227	PHE	3.0
1	C	274	PHE	3.0
1	E	38	PHE	3.0
1	C	-6	ASN	3.0
3	G	43	ARG	2.9
2	D	204	GLU	2.9
3	G	71	ASP	2.9
1	C	31	LEU	2.9
2	B	278	ALA	2.9
1	C	241	VAL	2.9
1	E	37	GLN	2.9
3	G	341	GLN	2.9
3	G	255	LYS	2.9
2	H	166	GLY	2.9
1	C	27	ALA	2.9
3	G	336	ALA	2.9
3	G	222	LYS	2.8
1	C	23	PRO	2.8
1	C	313	LEU	2.8
1	E	31	LEU	2.8
1	C	38	PHE	2.8
3	G	33	GLU	2.8
1	E	202	THR	2.8
2	H	167	GLY	2.8
3	G	219	GLY	2.8
1	E	217	PRO	2.8
1	C	165	PHE	2.8
3	G	16	ARG	2.8
1	A	-10	HIS	2.8
1	C	186	ASP	2.8
1	A	217	PRO	2.8
1	E	332	GLY	2.7
1	C	74	PRO	2.7
2	D	224	ARG	2.7
2	D	277	LEU	2.7
3	G	41	VAL	2.7
1	C	39	TRP	2.7
3	G	36	TRP	2.7

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Mol	Chain	Res	Type	RSRZ
2	D	170	ALA	2.7
3	G	179	VAL	2.7
3	G	281	GLU	2.7
1	A	24	VAL	2.6
2	D	167	GLY	2.6
3	G	52	GLY	2.6
2	F	170	ALA	2.6
1	C	237	MET	2.6
3	G	229	LEU	2.6
1	C	73	ALA	2.6
1	C	25	PHE	2.6
3	G	165	PHE	2.6
3	G	335	PHE	2.6
1	C	28	VAL	2.6
3	G	283	ILE	2.6
1	E	220	GLU	2.6
2	D	254	ARG	2.6
1	E	219	GLY	2.6
2	D	78	PRO	2.6
2	D	184	ILE	2.6
1	E	186	ASP	2.5
2	F	43	ARG	2.5
3	G	314	ARG	2.5
3	G	172	LEU	2.5
2	H	157	LYS	2.5
3	G	347	VAL	2.5
3	G	259	GLY	2.5
2	F	187	VAL	2.4
3	G	47	PHE	2.4
3	G	268	PHE	2.4
2	H	47	ASP	2.4
2	B	120	ARG	2.4
2	D	274	THR	2.4
1	A	31	LEU	2.4
2	D	231	GLY	2.4
3	G	95	VAL	2.4
3	G	299	VAL	2.4
1	C	226	TYR	2.4
1	C	33	GLU	2.4
1	C	296	ASP	2.4
3	G	48	ASP	2.4
3	G	70	SER	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	47	PHE	2.4
1	A	229	LEU	2.4
3	G	342	LEU	2.4
1	C	21	ASP	2.3
1	C	13	LEU	2.3
3	G	292	LEU	2.3
1	E	39	TRP	2.3
1	E	16	ARG	2.3
1	C	77	VAL	2.3
3	G	13	LEU	2.3
3	G	181	LEU	2.3
3	G	-6	ASN	2.3
1	C	217	PRO	2.3
1	E	29	PRO	2.3
1	A	-8	ALA	2.3
3	G	118	SER	2.3
3	G	250	ARG	2.3
3	G	318	ILE	2.3
1	C	24	VAL	2.3
1	E	255	LYS	2.3
2	D	256	ASP	2.3
1	C	212	ARG	2.3
3	G	68	ARG	2.3
3	G	265	TYR	2.3
3	G	169	THR	2.3
1	C	222	LYS	2.3
3	G	151	VAL	2.3
3	G	313	LEU	2.2
1	A	216	ASN	2.2
1	A	15	ALA	2.2
3	G	35	TRP	2.2
3	G	279	TRP	2.2
3	G	44	ARG	2.2
2	D	163	GLY	2.2
2	D	171	GLN	2.2
1	C	15	ALA	2.2
3	G	26	ALA	2.2
3	G	285	ALA	2.2
1	C	60	VAL	2.2
1	C	151	VAL	2.2
1	C	314	ARG	2.2
3	G	100	ARG	2.2

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Mol	Chain	Res	Type	RSRZ
3	G	296	ASP	2.2
1	E	168	ARG	2.2
1	A	274	PHE	2.2
3	G	1	MET	2.2
3	G	93	GLY	2.2
3	G	50	ARG	2.1
2	B	170	ALA	2.1
3	G	204	TYR	2.1
3	G	277	LYS	2.1
1	C	75	ASP	2.1
3	G	0	HIS	2.1
1	E	25	PHE	2.1
2	B	167	GLY	2.1
1	A	-7	GLU	2.1
3	G	149	GLU	2.1
1	E	188	ASP	2.1
2	H	91	ASN	2.1
3	G	75	ASP	2.1
1	C	223	PRO	2.1
3	G	345	VAL	2.1
1	A	203	PHE	2.1
1	C	34	SER	2.1
1	C	26	ALA	2.1
2	F	172	GLU	2.1
1	C	98	ASP	2.1
1	C	17	SER	2.0
1	E	177	CYS	2.0
1	C	213	LEU	2.0
3	G	309	LEU	2.0
3	G	263	ILE	2.0
3	G	324	ILE	2.0
1	C	16	ARG	2.0
1	E	101	GLY	2.0
3	G	-1	GLY	2.0
3	G	188	ASP	2.0
3	G	97	PRO	2.0

6.2 Non-standard residues in protein, DNA, RNA chains (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	CSX	G	129	7/8	0.92	0.11	55,56,71,72	0

6.3 Carbohydrates (i)

There are no oligosaccharides in this entry.

6.4 Ligands (i)

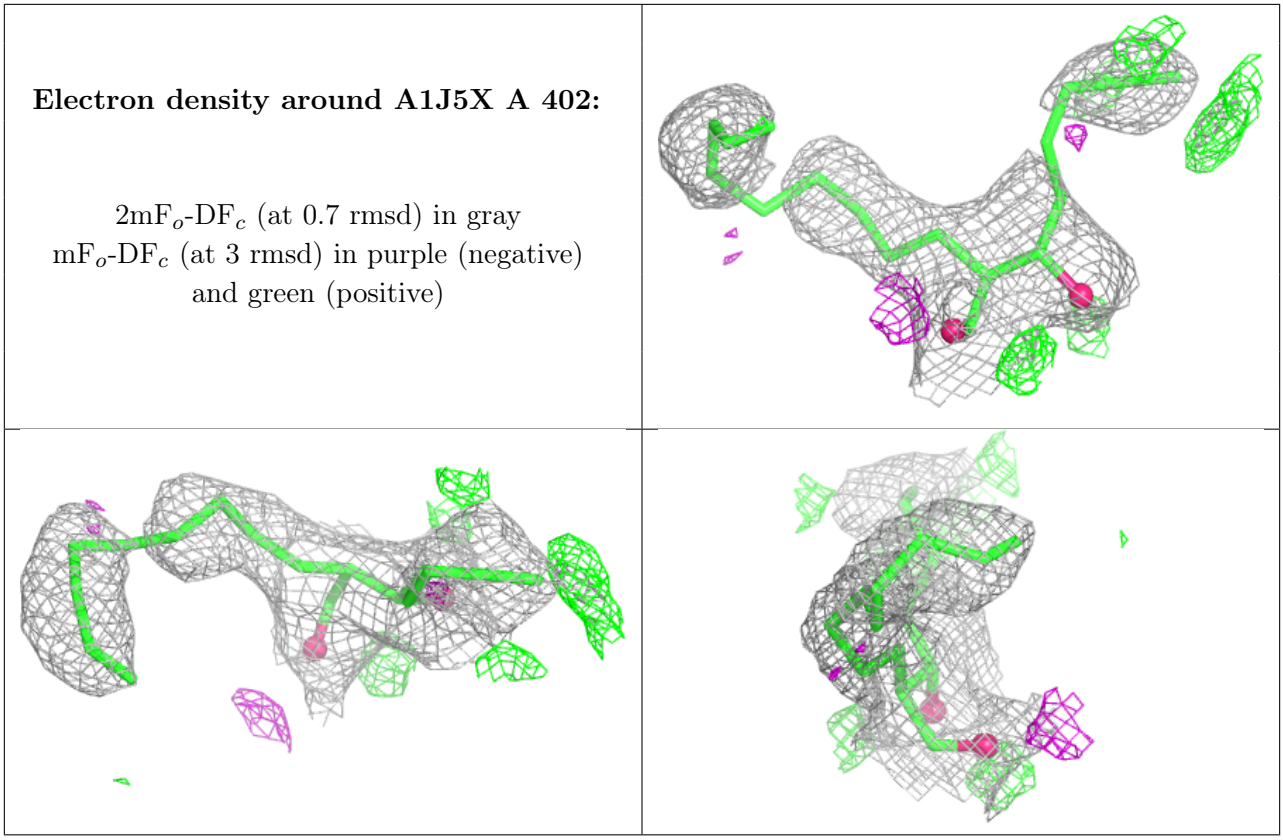
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' 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
4	EDO	A	401	4/4	0.85	0.19	41,43,50,52	0
5	A1J5X	A	402	19/19	0.88	0.19	40,50,72,77	0
5	A1J5X	C	402	19/19	0.89	0.17	50,59,65,71	0
4	EDO	C	401	4/4	0.91	0.15	33,45,50,55	0
5	A1J5X	E	401	19/19	0.91	0.15	38,48,66,66	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.

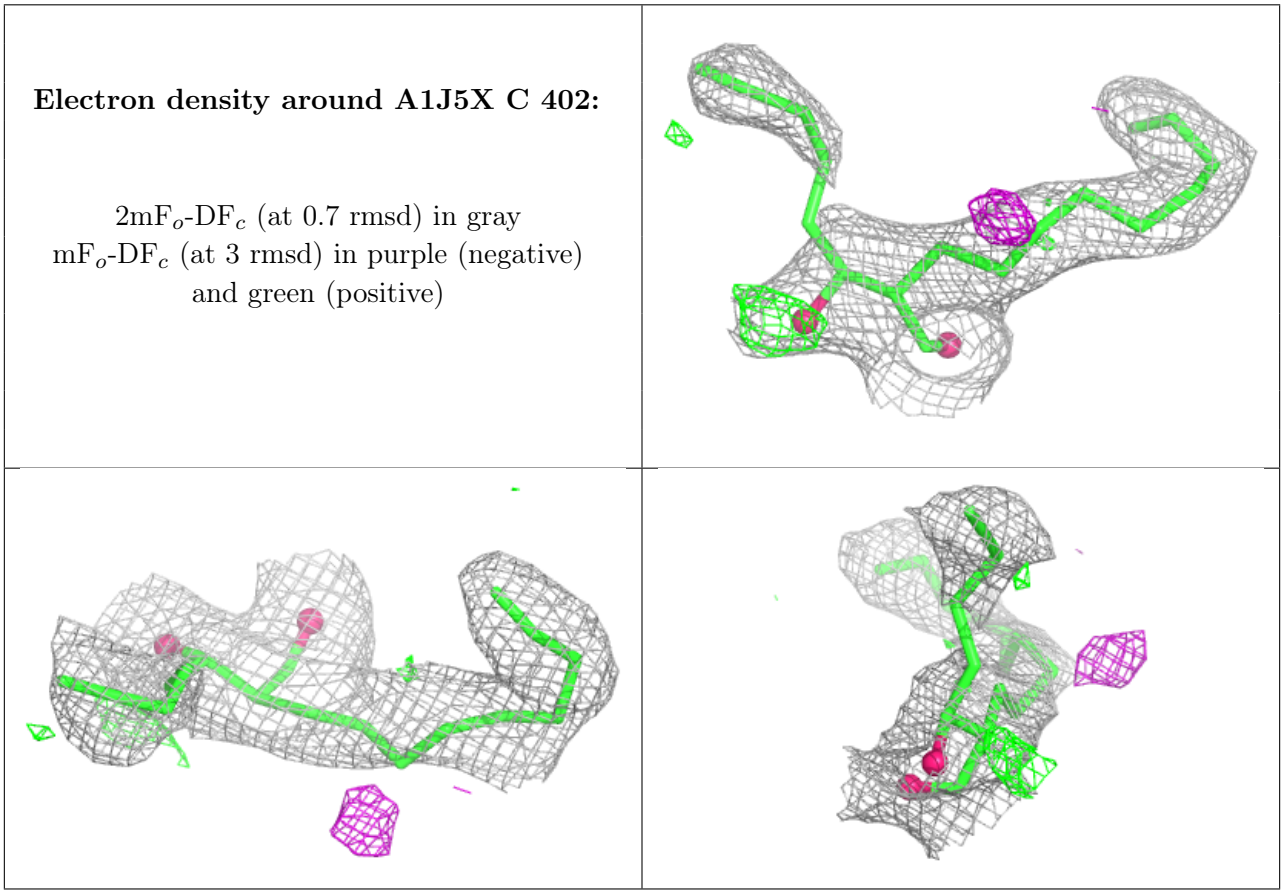
Electron density around A1J5X A 402:

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



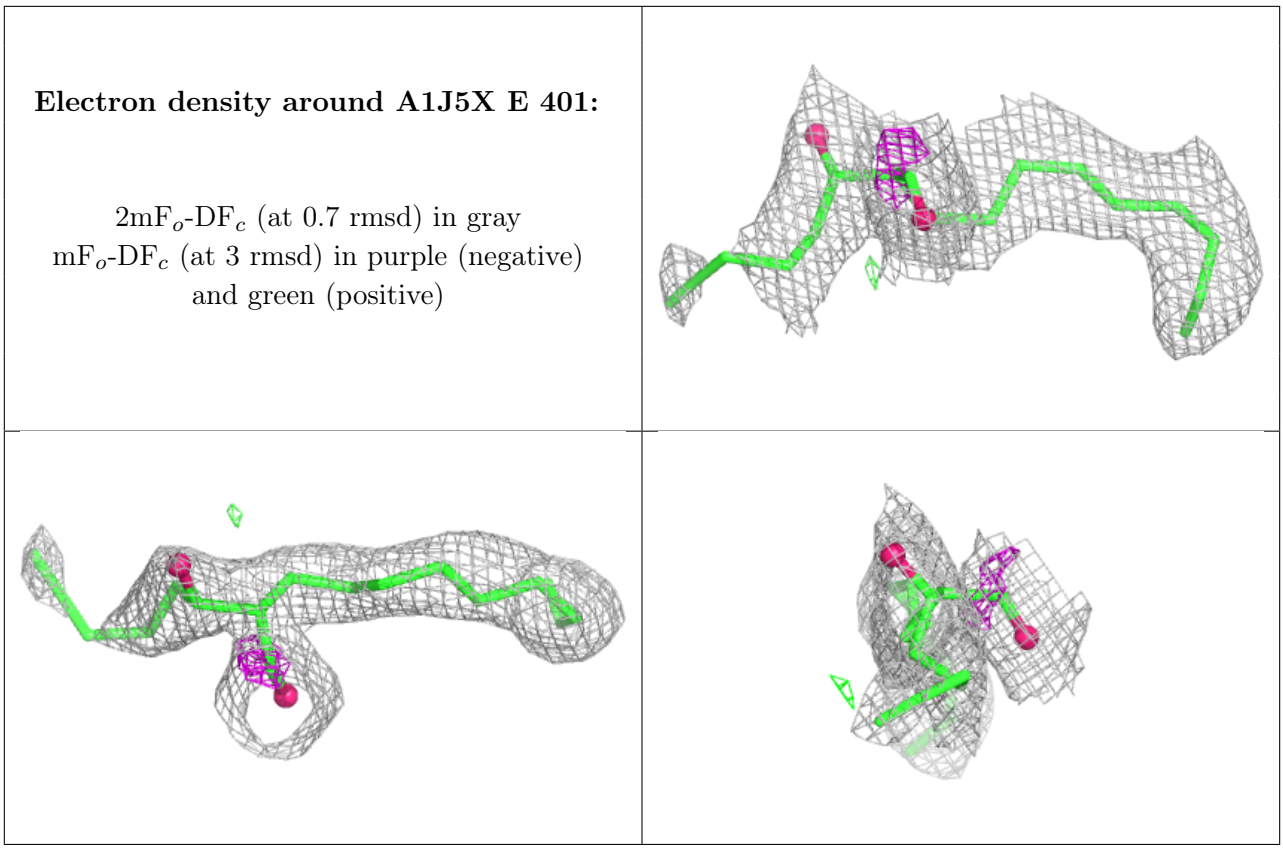
Electron density around A1J5X C 402:

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



Electron density around A1J5X 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)



6.5 Other polymers

There are no such residues in this entry.