



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2024 – 07:50 PM EST

PDB ID : 3WCM  
Title : The complex structure of HsSQS with ligand, ER119884  
Authors : Shang, N.; Li, Q.; Ko, T.P.; Chan, H.C.; Huang, C.H.; Ren, F.; Zheng, Y.;  
Zhu, Z.; Chen, C.C.; Guo, R.T.  
Deposited on : 2013-05-28  
Resolution : 2.06 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](mailto: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>.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.36  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.36

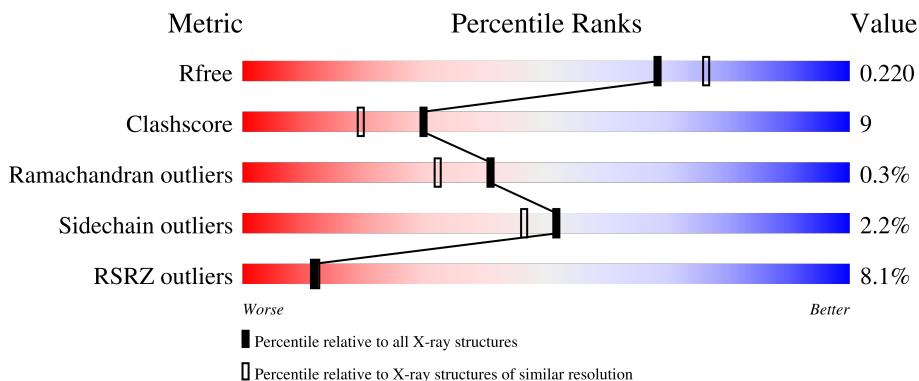
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.06 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.



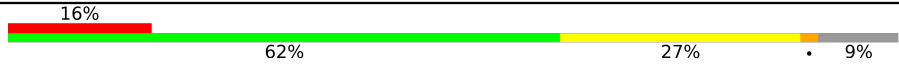
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	2684 (2.08-2.04)
Clashscore	141614	2801 (2.08-2.04)
Ramachandran outliers	138981	2768 (2.08-2.04)
Sidechain outliers	138945	2768 (2.08-2.04)
RSRZ outliers	127900	2646 (2.08-2.04)

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  $\geq 3$ , 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	360	<div style="display: flex; align-items: center;"> <div style="width: 4%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 77%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 14%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 9%; height: 10px; background-color: grey;"></div> </div>
1	B	360	<div style="display: flex; align-items: center;"> <div style="width: 7%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 74%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 17%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 8%; height: 10px; background-color: grey;"></div> </div>
1	C	360	<div style="display: flex; align-items: center;"> <div style="width: 6%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 74%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 17%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 8%; height: 10px; background-color: grey;"></div> </div>
1	D	360	<div style="display: flex; align-items: center;"> <div style="width: 7%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 74%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 16%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 10%; height: 10px; background-color: grey;"></div> </div>
1	E	360	<div style="display: flex; align-items: center;"> <div style="width: 5%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 75%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 16%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 8%; height: 10px; background-color: grey;"></div> </div>

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Mol	Chain	Length	Quality of chain
1	F	360	 <p>A horizontal bar chart representing the quality of the chain. The bar is divided into four segments: a red segment (16%), a green segment (62%), a yellow segment (27%), and a grey segment (9%). The percentages are labeled below the corresponding segments.</p>

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 17280 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 Squalene synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	329	Total 2659	C 1695	N 449	O 497	S 18	0	0	0
1	B	332	Total 2678	C 1705	N 452	O 503	S 18	0	0	0
1	C	332	Total 2683	C 1708	N 456	O 501	S 18	0	0	0
1	D	325	Total 2632	C 1674	N 446	O 494	S 18	0	0	0
1	E	330	Total 2664	C 1696	N 451	O 499	S 18	0	0	0
1	F	328	Total 2648	C 1686	N 448	O 496	S 18	0	0	0

There are 144 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	11	MET	-	expression tag	UNP P37268
A	12	GLY	-	expression tag	UNP P37268
A	13	SER	-	expression tag	UNP P37268
A	14	SER	-	expression tag	UNP P37268
A	15	HIS	-	expression tag	UNP P37268
A	16	HIS	-	expression tag	UNP P37268
A	17	HIS	-	expression tag	UNP P37268
A	18	HIS	-	expression tag	UNP P37268
A	19	HIS	-	expression tag	UNP P37268
A	20	HIS	-	expression tag	UNP P37268
A	21	SER	-	expression tag	UNP P37268
A	22	SER	-	expression tag	UNP P37268
A	23	GLY	-	expression tag	UNP P37268
A	24	LEU	-	expression tag	UNP P37268
A	25	VAL	-	expression tag	UNP P37268
A	26	PRO	-	expression tag	UNP P37268
A	27	ARG	-	expression tag	UNP P37268

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Chain	Residue	Modelled	Actual	Comment	Reference
A	28	GLY	-	expression tag	UNP P37268
A	29	SER	-	expression tag	UNP P37268
A	30	HIS	-	expression tag	UNP P37268
A	248	LEU	LYS	engineered mutation	UNP P37268
A	315	LEU	LYS	engineered mutation	UNP P37268
A	318	LEU	LYS	engineered mutation	UNP P37268
A	353	ASN	ASP	SEE REMARK 999	UNP P37268
B	11	MET	-	expression tag	UNP P37268
B	12	GLY	-	expression tag	UNP P37268
B	13	SER	-	expression tag	UNP P37268
B	14	SER	-	expression tag	UNP P37268
B	15	HIS	-	expression tag	UNP P37268
B	16	HIS	-	expression tag	UNP P37268
B	17	HIS	-	expression tag	UNP P37268
B	18	HIS	-	expression tag	UNP P37268
B	19	HIS	-	expression tag	UNP P37268
B	20	HIS	-	expression tag	UNP P37268
B	21	SER	-	expression tag	UNP P37268
B	22	SER	-	expression tag	UNP P37268
B	23	GLY	-	expression tag	UNP P37268
B	24	LEU	-	expression tag	UNP P37268
B	25	VAL	-	expression tag	UNP P37268
B	26	PRO	-	expression tag	UNP P37268
B	27	ARG	-	expression tag	UNP P37268
B	28	GLY	-	expression tag	UNP P37268
B	29	SER	-	expression tag	UNP P37268
B	30	HIS	-	expression tag	UNP P37268
B	248	LEU	LYS	engineered mutation	UNP P37268
B	315	LEU	LYS	engineered mutation	UNP P37268
B	318	LEU	LYS	engineered mutation	UNP P37268
B	353	ASN	ASP	SEE REMARK 999	UNP P37268
C	11	MET	-	expression tag	UNP P37268
C	12	GLY	-	expression tag	UNP P37268
C	13	SER	-	expression tag	UNP P37268
C	14	SER	-	expression tag	UNP P37268
C	15	HIS	-	expression tag	UNP P37268
C	16	HIS	-	expression tag	UNP P37268
C	17	HIS	-	expression tag	UNP P37268
C	18	HIS	-	expression tag	UNP P37268
C	19	HIS	-	expression tag	UNP P37268
C	20	HIS	-	expression tag	UNP P37268
C	21	SER	-	expression tag	UNP P37268

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Chain	Residue	Modelled	Actual	Comment	Reference
C	22	SER	-	expression tag	UNP P37268
C	23	GLY	-	expression tag	UNP P37268
C	24	LEU	-	expression tag	UNP P37268
C	25	VAL	-	expression tag	UNP P37268
C	26	PRO	-	expression tag	UNP P37268
C	27	ARG	-	expression tag	UNP P37268
C	28	GLY	-	expression tag	UNP P37268
C	29	SER	-	expression tag	UNP P37268
C	30	HIS	-	expression tag	UNP P37268
C	248	LEU	LYS	engineered mutation	UNP P37268
C	315	LEU	LYS	engineered mutation	UNP P37268
C	318	LEU	LYS	engineered mutation	UNP P37268
C	353	ASN	ASP	SEE REMARK 999	UNP P37268
D	11	MET	-	expression tag	UNP P37268
D	12	GLY	-	expression tag	UNP P37268
D	13	SER	-	expression tag	UNP P37268
D	14	SER	-	expression tag	UNP P37268
D	15	HIS	-	expression tag	UNP P37268
D	16	HIS	-	expression tag	UNP P37268
D	17	HIS	-	expression tag	UNP P37268
D	18	HIS	-	expression tag	UNP P37268
D	19	HIS	-	expression tag	UNP P37268
D	20	HIS	-	expression tag	UNP P37268
D	21	SER	-	expression tag	UNP P37268
D	22	SER	-	expression tag	UNP P37268
D	23	GLY	-	expression tag	UNP P37268
D	24	LEU	-	expression tag	UNP P37268
D	25	VAL	-	expression tag	UNP P37268
D	26	PRO	-	expression tag	UNP P37268
D	27	ARG	-	expression tag	UNP P37268
D	28	GLY	-	expression tag	UNP P37268
D	29	SER	-	expression tag	UNP P37268
D	30	HIS	-	expression tag	UNP P37268
D	248	LEU	LYS	engineered mutation	UNP P37268
D	315	LEU	LYS	engineered mutation	UNP P37268
D	318	LEU	LYS	engineered mutation	UNP P37268
D	353	ASN	ASP	SEE REMARK 999	UNP P37268
E	11	MET	-	expression tag	UNP P37268
E	12	GLY	-	expression tag	UNP P37268
E	13	SER	-	expression tag	UNP P37268
E	14	SER	-	expression tag	UNP P37268
E	15	HIS	-	expression tag	UNP P37268

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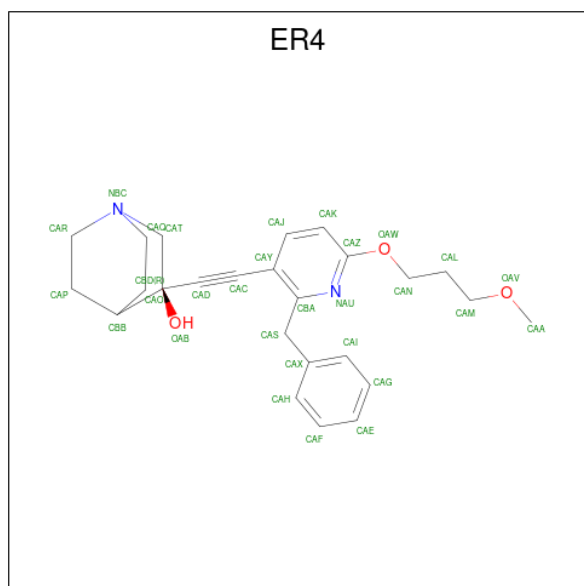
Chain	Residue	Modelled	Actual	Comment	Reference
E	16	HIS	-	expression tag	UNP P37268
E	17	HIS	-	expression tag	UNP P37268
E	18	HIS	-	expression tag	UNP P37268
E	19	HIS	-	expression tag	UNP P37268
E	20	HIS	-	expression tag	UNP P37268
E	21	SER	-	expression tag	UNP P37268
E	22	SER	-	expression tag	UNP P37268
E	23	GLY	-	expression tag	UNP P37268
E	24	LEU	-	expression tag	UNP P37268
E	25	VAL	-	expression tag	UNP P37268
E	26	PRO	-	expression tag	UNP P37268
E	27	ARG	-	expression tag	UNP P37268
E	28	GLY	-	expression tag	UNP P37268
E	29	SER	-	expression tag	UNP P37268
E	30	HIS	-	expression tag	UNP P37268
E	248	LEU	LYS	engineered mutation	UNP P37268
E	315	LEU	LYS	engineered mutation	UNP P37268
E	318	LEU	LYS	engineered mutation	UNP P37268
E	353	ASN	ASP	SEE REMARK 999	UNP P37268
F	11	MET	-	expression tag	UNP P37268
F	12	GLY	-	expression tag	UNP P37268
F	13	SER	-	expression tag	UNP P37268
F	14	SER	-	expression tag	UNP P37268
F	15	HIS	-	expression tag	UNP P37268
F	16	HIS	-	expression tag	UNP P37268
F	17	HIS	-	expression tag	UNP P37268
F	18	HIS	-	expression tag	UNP P37268
F	19	HIS	-	expression tag	UNP P37268
F	20	HIS	-	expression tag	UNP P37268
F	21	SER	-	expression tag	UNP P37268
F	22	SER	-	expression tag	UNP P37268
F	23	GLY	-	expression tag	UNP P37268
F	24	LEU	-	expression tag	UNP P37268
F	25	VAL	-	expression tag	UNP P37268
F	26	PRO	-	expression tag	UNP P37268
F	27	ARG	-	expression tag	UNP P37268
F	28	GLY	-	expression tag	UNP P37268
F	29	SER	-	expression tag	UNP P37268
F	30	HIS	-	expression tag	UNP P37268
F	248	LEU	LYS	engineered mutation	UNP P37268
F	315	LEU	LYS	engineered mutation	UNP P37268
F	318	LEU	LYS	engineered mutation	UNP P37268

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Chain	Residue	Modelled	Actual	Comment	Reference
F	353	ASN	ASP	SEE REMARK 999	UNP P37268

- Molecule 2 is (3R)-3-[[2-benzyl-6-(3-methoxypropoxy)pyridin-3-yl]ethynyl]-1-azabicyclo[2.2.2]octan-3-ol (three-letter code: ER4) (formula: C<sub>25</sub>H<sub>30</sub>N<sub>2</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	Total	C	N	O	0	0
			30	25	2	3		
2	B	1	Total	C	N	O	0	0
			30	25	2	3		
2	C	1	Total	C	N	O	0	0
			30	25	2	3		
2	D	1	Total	C	N	O	0	0
			30	25	2	3		
2	E	1	Total	C	N	O	0	0
			30	25	2	3		
2	F	1	Total	C	N	O	0	0
			30	25	2	3		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	273	Total	O	0	0
			273	273		
3	B	234	Total	O	0	0
			234	234		

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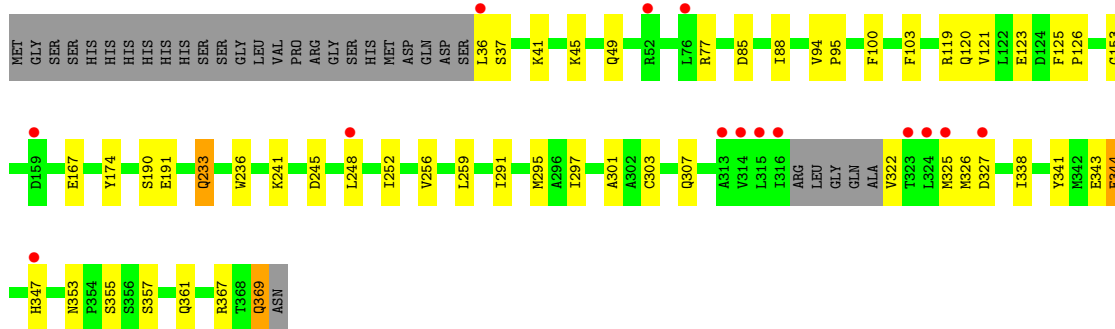
<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
3	C	236	Total 236	O 236	0	0
3	D	168	Total 168	O 168	0	0
3	E	141	Total 141	O 141	0	0
3	F	84	Total 84	O 84	0	0

### 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.

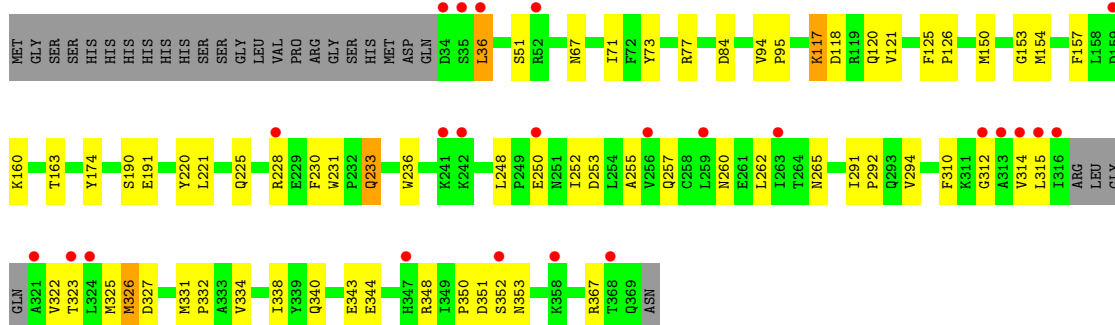
- Molecule 1: Squalene synthase

Chain A: 




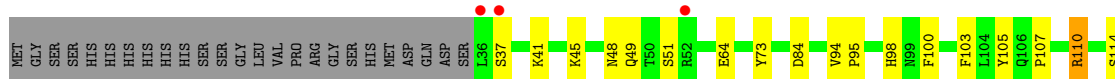
- Molecule 1: Squalene synthase

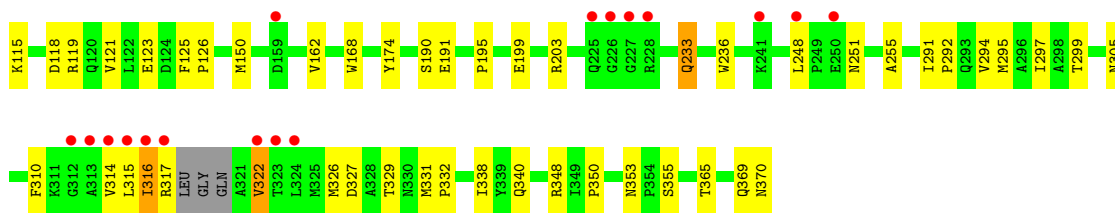
Chain B: 



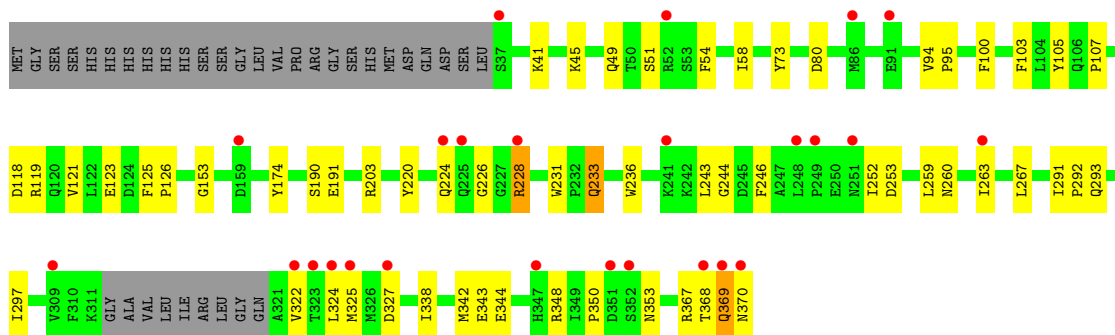
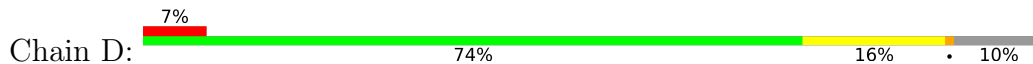
- Molecule 1: Squalene synthase

Chain C: 

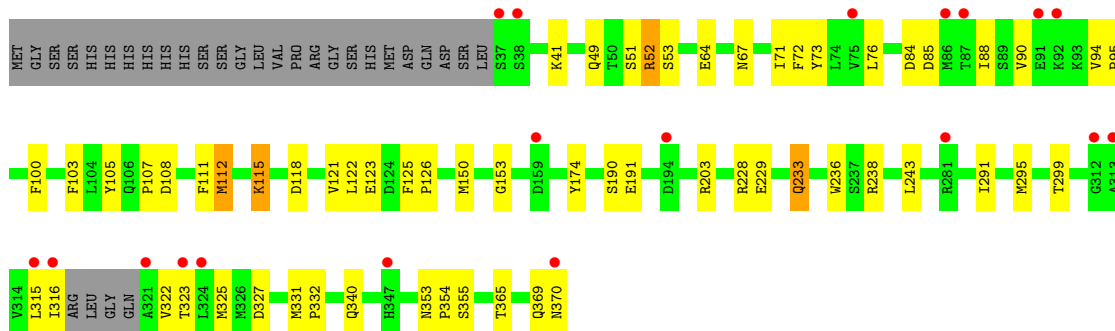
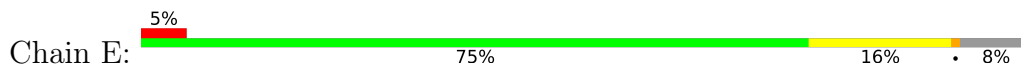




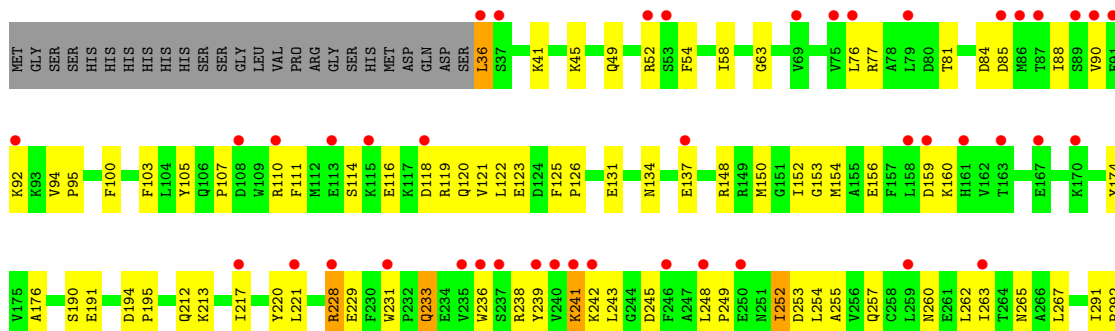
• Molecule 1: Squalene synthase

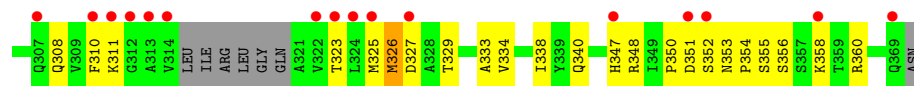


• Molecule 1: Squalene synthase



• Molecule 1: Squalene synthase





## 4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	85.62Å 153.45Å 92.12Å 90.00° 90.86° 90.00°	Depositor
Resolution (Å)	25.00 – 2.06 24.94 – 2.06	Depositor EDS
% Data completeness (in resolution range)	(Not available) (25.00-2.06) 95.9 (24.94-2.06)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.20 (at 2.06Å)	Xtrriage
Refinement program	CNS 1.2	Depositor
R, $R_{free}$	0.186 , 0.222 0.185 , 0.220	Depositor DCC
$R_{free}$ test set	7126 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	25.2	Xtrriage
Anisotropy	0.669	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 59.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.024 for h,-k,-l	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	17280	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	37.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.81% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ER4

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	0.35	0/2713	0.53	0/3672
1	B	0.33	0/2732	0.53	0/3698
1	C	0.33	0/2737	0.52	0/3704
1	D	0.32	0/2686	0.52	0/3635
1	E	0.29	0/2718	0.49	0/3679
1	F	0.28	0/2702	0.48	0/3657
All	All	0.32	0/16288	0.51	0/22045

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 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	2659	0	2635	42	0
1	B	2678	0	2649	52	0
1	C	2683	0	2659	61	0
1	D	2632	0	2596	44	0
1	E	2664	0	2635	47	0
1	F	2648	0	2618	82	0
2	A	30	0	30	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	30	0	30	1	0
2	C	30	0	30	0	0
2	D	30	0	30	1	0
2	E	30	0	30	0	0
2	F	30	0	30	1	0
3	A	273	0	0	3	0
3	B	234	0	0	1	0
3	C	236	0	0	5	0
3	D	168	0	0	1	0
3	E	141	0	0	3	0
3	F	84	0	0	3	0
All	All	17280	0	15972	302	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 9.

All (302) 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:D:322:VAL:HG13	1:E:291:ILE:HG21	1.45	0.98
1:D:260:ASN:HD22	1:D:353:ASN:ND2	1.69	0.89
1:F:228:ARG:HH11	1:F:228:ARG:H	1.14	0.89
1:D:369:GLN:O	1:D:370:ASN:HB2	1.75	0.87
1:F:131:GLU:HA	1:F:134:ASN:HD22	1.41	0.85
1:F:260:ASN:HD22	1:F:353:ASN:ND2	1.74	0.84
1:B:326:MET:HE3	1:C:291:ILE:HD11	1.60	0.81
1:C:95:PRO:HA	3:C:572:HOH:O	1.80	0.81
1:F:233:GLN:HA	1:F:236:TRP:NE1	1.96	0.81
1:D:368:THR:HG21	1:E:41:LYS:HA	1.66	0.78
1:E:150:MET:HG3	1:E:174:TYR:O	1.85	0.75
1:C:45:LYS:O	1:C:49:GLN:HG3	1.85	0.74
1:C:348:ARG:HB3	1:C:348:ARG:HH11	1.51	0.74
1:F:45:LYS:HG2	1:F:49:GLN:HE21	1.53	0.73
1:F:94:VAL:HB	1:F:95:PRO:HD3	1.70	0.73
1:D:94:VAL:HB	1:D:95:PRO:HD3	1.70	0.72
1:B:233:GLN:HA	1:B:236:TRP:NE1	2.05	0.71
1:A:291:ILE:HG21	1:C:322:VAL:HG22	1.71	0.70
1:C:195:PRO:O	1:C:199:GLU:HG3	1.92	0.70
1:B:77:ARG:HD3	2:B:401:ER4:H7	1.74	0.69
1:F:228:ARG:HH11	1:F:228:ARG:N	1.88	0.69
1:A:357:SER:O	1:A:361:GLN:HG3	1.92	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:353:ASN:ND2	1:A:355:SER:H	1.91	0.69
1:F:228:ARG:HB2	1:F:228:ARG:NH1	2.08	0.68
1:E:322:VAL:CG2	1:F:291:ILE:HG21	2.23	0.68
1:E:365:THR:O	1:E:369:GLN:HG3	1.94	0.67
1:F:118:ASP:O	1:F:121:VAL:HG22	1.96	0.66
1:D:224:GLN:HE21	1:D:244:GLY:HA2	1.61	0.66
1:E:322:VAL:HG12	1:E:340:GLN:NE2	2.11	0.66
1:B:322:VAL:HG13	1:C:291:ILE:HG21	1.77	0.66
1:C:348:ARG:HB3	1:C:348:ARG:NH1	2.12	0.65
1:D:233:GLN:HA	1:D:236:TRP:NE1	2.11	0.65
1:E:90:VAL:O	1:E:94:VAL:HG23	1.97	0.65
1:E:353:ASN:ND2	1:E:355:SER:H	1.94	0.65
1:C:322:VAL:HG12	1:C:340:GLN:CD	2.17	0.65
1:B:260:ASN:HD22	1:B:353:ASN:ND2	1.96	0.63
1:F:150:MET:HG3	1:F:174:TYR:O	1.97	0.63
1:C:100:PHE:HA	1:C:103:PHE:CD2	2.34	0.63
1:D:327:ASP:HB2	1:F:325:MET:O	1.98	0.63
1:C:315:LEU:O	1:C:315:LEU:HD12	1.99	0.63
1:A:326:MET:HA	1:B:327:ASP:HB2	1.81	0.62
1:D:260:ASN:HD22	1:D:353:ASN:HD22	1.46	0.62
1:F:242:LYS:HB2	1:F:245:ASP:OD2	1.98	0.62
1:E:52:ARG:HD2	1:E:53:SER:N	2.15	0.62
1:F:260:ASN:HD22	1:F:353:ASN:HD22	1.48	0.62
1:A:245:ASP:HA	1:A:248:LEU:HD13	1.82	0.61
1:F:131:GLU:HA	1:F:134:ASN:ND2	2.14	0.61
1:D:45:LYS:O	1:D:49:GLN:HG3	2.00	0.61
1:F:228:ARG:H	1:F:228:ARG:NH1	1.94	0.61
1:E:233:GLN:HA	1:E:236:TRP:NE1	2.15	0.61
1:B:221:LEU:HD23	1:B:312:GLY:HA2	1.83	0.60
1:E:295:MET:HA	1:E:295:MET:CE	2.32	0.60
1:E:323:THR:HB	1:E:340:GLN:OE1	2.01	0.60
1:E:100:PHE:HA	1:E:103:PHE:CD2	2.36	0.60
1:E:322:VAL:HG23	1:F:291:ILE:HG21	1.83	0.59
1:C:233:GLN:HA	1:C:236:TRP:NE1	2.17	0.59
3:A:1113:HOH:O	1:C:370:ASN:HB2	2.03	0.58
1:B:253:ASP:O	1:B:257:GLN:HG3	2.03	0.58
1:C:98:HIS:HB2	3:C:572:HOH:O	2.04	0.58
1:F:90:VAL:O	1:F:94:VAL:HG23	2.03	0.58
1:B:323:THR:CG2	1:B:340:GLN:HG3	2.34	0.58
1:A:343:GLU:OE1	1:A:367:ARG:NH2	2.37	0.57
1:E:52:ARG:CD	1:E:52:ARG:H	2.17	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:118:ASP:O	1:C:121:VAL:HG22	2.03	0.57
1:F:100:PHE:HA	1:F:103:PHE:CD2	2.38	0.57
1:D:119:ARG:O	1:D:123:GLU:HG3	2.04	0.57
1:D:291:ILE:HB	1:D:292:PRO:HD3	1.86	0.57
1:E:315:LEU:HD12	1:E:315:LEU:O	2.05	0.57
1:F:36:LEU:HB2	1:F:41:LYS:HE3	1.86	0.56
1:C:305:ASN:HB3	1:C:348:ARG:HH12	1.70	0.56
1:D:100:PHE:HA	1:D:103:PHE:CD2	2.41	0.56
1:C:150:MET:HG3	1:C:174:TYR:O	2.06	0.56
1:F:85:ASP:HB3	1:F:88:ILE:HD13	1.88	0.55
1:F:253:ASP:O	1:F:257:GLN:HG3	2.07	0.55
1:C:348:ARG:O	1:C:350:PRO:HD3	2.07	0.55
1:E:52:ARG:HD2	1:E:53:SER:H	1.69	0.55
1:E:112:MET:HE1	1:E:123:GLU:HB3	1.89	0.55
1:E:331:MET:HB3	1:E:332:PRO:HD3	1.89	0.55
1:A:322:VAL:O	1:A:326:MET:HG2	2.07	0.55
1:F:220:TYR:HB2	1:F:231:TRP:CZ2	2.42	0.55
1:F:36:LEU:N	1:F:36:LEU:HD22	2.21	0.55
1:C:94:VAL:HB	1:C:95:PRO:HD3	1.89	0.55
1:D:105:TYR:O	1:D:107:PRO:HD3	2.06	0.55
1:D:41:LYS:HE2	3:F:516:HOH:O	2.08	0.54
1:E:118:ASP:O	1:E:121:VAL:HG22	2.07	0.54
1:F:323:THR:HB	1:F:340:GLN:OE1	2.06	0.54
1:F:326:MET:HE2	1:F:333:ALA:HA	1.89	0.54
1:E:323:THR:H	1:E:340:GLN:HE22	1.55	0.54
1:D:369:GLN:O	1:D:370:ASN:CB	2.52	0.54
1:E:105:TYR:O	1:E:107:PRO:HD3	2.07	0.54
1:D:125:PHE:N	1:D:126:PRO:CD	2.71	0.54
1:D:324:LEU:HD12	1:D:325:MET:HG2	1.90	0.54
1:E:125:PHE:N	1:E:126:PRO:CD	2.70	0.54
1:C:322:VAL:HG12	1:C:340:GLN:OE1	2.06	0.54
1:A:327:ASP:HB3	1:C:326:MET:HA	1.90	0.53
1:E:67:ASN:O	1:E:71:ILE:HG12	2.09	0.53
1:A:325:MET:HE2	1:B:294:VAL:HG11	1.89	0.53
1:C:314:VAL:CG1	1:C:315:LEU:N	2.72	0.53
1:E:370:ASN:HB3	3:F:522:HOH:O	2.08	0.53
1:F:88:ILE:HG23	1:F:92:LYS:HB3	1.91	0.52
1:F:88:ILE:N	1:F:88:ILE:HD12	2.24	0.52
1:A:325:MET:O	1:B:327:ASP:HB2	2.09	0.52
1:E:325:MET:O	1:F:327:ASP:HB3	2.10	0.52
1:F:116:GLU:O	1:F:119:ARG:HG2	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:36:LEU:HG	1:A:37:SER:N	2.24	0.52
1:C:348:ARG:HH11	1:C:348:ARG:CB	2.22	0.52
1:E:51:SER:HB2	1:E:73:TYR:CZ	2.44	0.52
1:A:325:MET:CE	1:B:294:VAL:HG11	2.39	0.52
1:C:37:SER:O	1:C:41:LYS:HG3	2.10	0.52
1:C:64:GLU:HG3	3:C:556:HOH:O	2.10	0.52
1:D:343:GLU:OE2	1:D:367:ARG:NE	2.43	0.52
1:D:348:ARG:O	1:D:350:PRO:HD3	2.10	0.52
1:D:51:SER:HB2	1:D:73:TYR:CZ	2.44	0.52
1:E:228:ARG:HG3	1:E:228:ARG:HH11	1.75	0.52
1:A:233:GLN:HA	1:A:236:TRP:NE1	2.25	0.51
1:C:314:VAL:HG12	1:C:315:LEU:N	2.25	0.51
1:F:263:ILE:O	1:F:267:LEU:HG	2.10	0.51
1:A:252:ILE:HG13	1:A:307:GLN:HG2	1.92	0.51
1:C:105:TYR:O	1:C:107:PRO:HD3	2.10	0.51
1:D:293:GLN:O	1:D:297:ILE:HG13	2.11	0.51
1:D:220:TYR:HB2	1:D:231:TRP:CZ2	2.44	0.51
1:D:350:PRO:HG2	1:D:353:ASN:HB2	1.93	0.51
1:C:365:THR:O	1:C:369:GLN:HG3	2.09	0.51
1:F:326:MET:CE	1:F:333:ALA:HA	2.41	0.51
1:F:291:ILE:HB	1:F:292:PRO:HD3	1.92	0.51
1:B:325:MET:O	1:C:327:ASP:HB2	2.10	0.50
1:C:248:LEU:HB2	1:C:251:ASN:HD22	1.75	0.50
1:F:118:ASP:HA	1:F:120:GLN:OE1	2.12	0.50
1:A:100:PHE:HA	1:A:103:PHE:CD2	2.47	0.50
1:A:291:ILE:CG2	1:C:322:VAL:HG22	2.41	0.50
1:F:81:THR:HG23	1:F:116:GLU:HG3	1.92	0.50
1:B:325:MET:CE	1:C:295:MET:HG3	2.42	0.50
1:B:150:MET:O	1:B:154:MET:HG3	2.10	0.50
1:A:36:LEU:HG	1:A:37:SER:H	1.76	0.50
1:C:114:SER:C	1:C:115:LYS:HD2	2.33	0.50
1:D:233:GLN:HG3	1:D:236:TRP:CZ2	2.46	0.50
1:F:52:ARG:HG3	3:F:583:HOH:O	2.11	0.50
1:F:255:ALA:HB1	1:F:310:PHE:CE2	2.47	0.50
1:B:125:PHE:N	1:B:126:PRO:CD	2.75	0.49
1:F:308:GLN:OE1	1:F:311:LYS:HD2	2.11	0.49
1:F:92:LYS:O	1:F:95:PRO:HD2	2.12	0.49
1:F:233:GLN:HA	1:F:236:TRP:CD1	2.47	0.49
1:B:118:ASP:O	1:B:121:VAL:HG22	2.12	0.49
1:E:115:LYS:HZ2	1:E:115:LYS:HB3	1.77	0.49
1:E:299:THR:HA	1:E:316:ILE:HD12	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:241:LYS:HD3	3:A:1188:HOH:O	2.12	0.49
1:A:94:VAL:HB	1:A:95:PRO:HD3	1.95	0.49
1:C:291:ILE:HB	1:C:292:PRO:HD3	1.94	0.49
1:C:327:ASP:OD2	1:C:329:THR:HG23	2.13	0.49
1:A:301:ALA:HB1	1:A:344:GLU:HG2	1.95	0.49
1:C:299:THR:HA	1:C:316:ILE:HD12	1.95	0.49
1:C:125:PHE:N	1:C:126:PRO:CD	2.76	0.48
1:F:239:TYR:O	1:F:254:LEU:HD13	2.13	0.48
1:E:64:GLU:HG3	3:E:506:HOH:O	2.13	0.48
1:F:252:ILE:HG23	1:F:253:ASP:N	2.28	0.48
1:B:344:GLU:O	1:B:348:ARG:HD3	2.12	0.48
1:E:295:MET:HA	1:E:295:MET:HE2	1.95	0.48
1:B:252:ILE:HG23	1:B:253:ASP:N	2.28	0.48
1:C:233:GLN:HA	1:C:236:TRP:CD1	2.49	0.48
1:F:153:GLY:O	1:F:156:GLU:HB3	2.14	0.48
1:F:262:LEU:O	1:F:265:ASN:HB3	2.13	0.48
1:F:190:SER:O	1:F:191:GLU:HB2	2.14	0.48
1:F:150:MET:O	1:F:154:MET:HG3	2.14	0.48
1:A:252:ILE:CG1	1:A:307:GLN:HG2	2.44	0.47
1:C:353:ASN:ND2	1:C:355:SER:H	2.12	0.47
3:E:580:HOH:O	1:F:329:THR:HG21	2.13	0.47
1:B:77:ARG:HD2	3:B:627:HOH:O	2.14	0.47
1:D:153:GLY:HA3	1:D:174:TYR:CG	2.48	0.47
1:A:45:LYS:O	1:A:49:GLN:HG3	2.14	0.47
1:B:325:MET:HE2	1:C:295:MET:HG3	1.97	0.47
1:D:220:TYR:OH	1:D:246:PHE:HB2	2.14	0.47
1:A:119:ARG:O	1:A:123:GLU:HG3	2.15	0.47
1:E:233:GLN:HA	1:E:236:TRP:CD1	2.50	0.47
1:A:125:PHE:N	1:A:126:PRO:CD	2.78	0.47
1:B:325:MET:HE3	1:C:294:VAL:CG1	2.45	0.47
1:E:190:SER:O	1:E:191:GLU:HB2	2.14	0.47
1:F:236:TRP:CG	1:F:243:LEU:HD13	2.50	0.47
1:A:252:ILE:O	1:A:256:VAL:HG23	2.15	0.46
1:A:326:MET:HA	1:B:327:ASP:CB	2.45	0.46
1:B:331:MET:HB3	1:B:332:PRO:HD3	1.96	0.46
1:F:334:VAL:O	1:F:338:ILE:HG13	2.15	0.46
1:A:85:ASP:HB3	1:A:88:ILE:HD12	1.97	0.46
1:C:110:ARG:HB3	1:C:126:PRO:HD3	1.98	0.46
1:C:315:LEU:O	1:C:317:ARG:N	2.48	0.46
1:F:228:ARG:HB2	1:F:228:ARG:CZ	2.44	0.46
1:A:190:SER:O	1:A:191:GLU:HB2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:343:GLU:OE1	1:B:367:ARG:NH2	2.49	0.46
1:E:115:LYS:HB3	1:E:115:LYS:NZ	2.30	0.46
1:A:245:ASP:HA	1:A:248:LEU:CD1	2.46	0.46
1:A:369:GLN:HE22	1:B:36:LEU:CD2	2.29	0.46
1:B:220:TYR:HB2	1:B:231:TRP:CZ2	2.50	0.46
1:A:36:LEU:CG	1:A:37:SER:H	2.27	0.46
1:B:348:ARG:O	1:B:350:PRO:HD3	2.15	0.46
1:B:291:ILE:HB	1:B:292:PRO:HD3	1.97	0.46
1:C:331:MET:HB3	1:C:332:PRO:HD3	1.98	0.46
1:D:190:SER:O	1:D:191:GLU:HB2	2.15	0.46
1:D:233:GLN:HA	1:D:236:TRP:CE2	2.50	0.46
1:B:163:THR:HA	1:B:233:GLN:HB3	1.98	0.46
1:B:326:MET:HA	1:C:327:ASP:OD2	2.15	0.46
1:F:255:ALA:HB1	1:F:310:PHE:CZ	2.50	0.46
1:C:190:SER:O	1:C:191:GLU:HB2	2.16	0.46
1:D:118:ASP:O	1:D:121:VAL:HG22	2.16	0.46
1:F:213:LYS:O	1:F:217:ILE:HG13	2.16	0.46
1:A:153:GLY:HA3	1:A:174:TYR:CG	2.51	0.46
1:A:295:MET:SD	1:C:322:VAL:HG23	2.56	0.46
1:B:190:SER:O	1:B:191:GLU:HB2	2.15	0.46
1:E:94:VAL:HB	1:E:95:PRO:HD3	1.97	0.46
1:E:354:PRO:HG2	3:E:515:HOH:O	2.14	0.46
1:B:120:GLN:HA	1:B:120:GLN:NE2	2.30	0.45
1:B:221:LEU:O	1:B:225:GLN:HG3	2.16	0.45
1:D:224:GLN:C	1:D:226:GLY:H	2.20	0.45
1:F:125:PHE:N	1:F:126:PRO:CD	2.78	0.45
1:B:157:PHE:CE2	1:B:160:LYS:HE3	2.51	0.45
1:C:94:VAL:HG11	3:C:704:HOH:O	2.16	0.45
1:B:255:ALA:HB1	1:B:310:PHE:CZ	2.52	0.45
1:B:314:VAL:HG12	1:B:315:LEU:N	2.31	0.45
1:B:94:VAL:HB	1:B:95:PRO:HD3	1.98	0.45
1:E:85:ASP:HB3	1:E:88:ILE:HD12	1.99	0.45
1:B:334:VAL:O	1:B:338:ILE:HG13	2.17	0.45
1:F:252:ILE:O	1:F:255:ALA:HB3	2.17	0.45
1:B:51:SER:HB2	1:B:73:TYR:CZ	2.52	0.45
1:B:117:LYS:HG3	1:B:118:ASP:N	2.31	0.45
1:D:224:GLN:OE1	1:D:224:GLN:HA	2.17	0.44
1:F:156:GLU:HG2	1:F:160:LYS:HE2	1.98	0.44
1:E:72:PHE:O	1:E:76:LEU:HG	2.17	0.44
1:E:150:MET:HB2	1:E:150:MET:HE3	1.78	0.44
1:D:228:ARG:HG2	3:D:668:HOH:O	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:350:PRO:C	1:F:352:SER:H	2.19	0.44
1:A:77:ARG:NH1	2:A:901:ER4:H9	2.32	0.44
1:E:153:GLY:HA3	1:E:174:TYR:CG	2.52	0.44
1:F:351:ASP:HB3	1:F:360:ARG:HH22	1.83	0.44
1:C:297:ILE:CD1	1:C:338:ILE:HG12	2.48	0.44
1:C:299:THR:HA	1:C:316:ILE:CD1	2.47	0.44
1:D:338:ILE:O	1:D:342:MET:HG2	2.17	0.44
1:D:343:GLU:OE1	1:D:367:ARG:NH2	2.51	0.44
1:E:353:ASN:HD22	1:E:355:SER:H	1.61	0.44
1:F:111:PHE:HB3	1:F:122:LEU:HB3	1.98	0.44
1:F:148:ARG:O	1:F:152:ILE:HG13	2.18	0.44
1:B:228:ARG:HD2	1:B:230:PHE:CE2	2.52	0.44
1:C:255:ALA:HB1	1:C:310:PHE:CZ	2.52	0.44
1:A:120:GLN:HG2	1:A:121:VAL:N	2.33	0.43
1:F:114:SER:O	1:F:119:ARG:HD3	2.18	0.43
1:E:233:GLN:HA	1:E:236:TRP:CE2	2.53	0.43
1:C:305:ASN:HB3	1:C:348:ARG:NH1	2.32	0.43
1:F:229:GLU:HB3	1:F:243:LEU:HD23	2.00	0.43
1:F:236:TRP:CE2	1:F:243:LEU:HB2	2.53	0.43
1:C:94:VAL:CG1	3:C:704:HOH:O	2.67	0.43
1:D:263:ILE:O	1:D:267:LEU:HG	2.19	0.43
1:F:241:LYS:HB2	1:F:241:LYS:NZ	2.33	0.43
1:C:119:ARG:O	1:C:123:GLU:HG3	2.18	0.43
1:F:76:LEU:HD12	2:F:401:ER4:H28	2.00	0.43
1:F:248:LEU:HA	1:F:249:PRO:HD3	1.92	0.43
1:B:343:GLU:OE1	1:C:48:ASN:HB3	2.19	0.43
1:D:260:ASN:HB2	1:D:353:ASN:HD21	1.83	0.43
1:F:236:TRP:C	1:F:238:ARG:H	2.21	0.43
1:D:344:GLU:CD	1:D:348:ARG:HH12	2.23	0.42
1:E:322:VAL:HG22	1:F:291:ILE:HG21	2.00	0.42
1:C:51:SER:HB2	1:C:73:TYR:CZ	2.54	0.42
1:F:221:LEU:HD23	1:F:311:LYS:O	2.19	0.42
1:F:233:GLN:HE21	1:F:233:GLN:HB3	1.60	0.42
1:A:120:GLN:HB2	3:A:1125:HOH:O	2.19	0.42
1:B:350:PRO:HB2	1:B:353:ASN:H	1.84	0.42
1:D:260:ASN:HD22	1:D:353:ASN:HD21	1.62	0.42
1:E:322:VAL:HG12	1:E:340:GLN:CD	2.39	0.42
1:F:54:PHE:O	1:F:58:ILE:HG13	2.20	0.42
1:F:348:ARG:O	1:F:350:PRO:HD3	2.19	0.42
1:B:350:PRO:C	1:B:352:SER:N	2.73	0.42
1:F:355:SER:O	1:F:358:LYS:N	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:259:LEU:O	1:D:263:ILE:HG13	2.20	0.42
1:F:105:TYR:O	1:F:107:PRO:HD3	2.20	0.42
1:B:67:ASN:O	1:B:71:ILE:HG12	2.20	0.42
1:C:350:PRO:HG2	1:C:353:ASN:HB2	2.01	0.42
1:F:353:ASN:HA	1:F:354:PRO:HD3	1.85	0.42
1:A:297:ILE:CD1	1:A:338:ILE:HG12	2.50	0.41
1:B:120:GLN:HA	1:B:120:GLN:HE21	1.85	0.41
1:C:316:ILE:HD13	1:C:316:ILE:HA	1.93	0.41
1:F:260:ASN:ND2	1:F:353:ASN:ND2	2.55	0.41
1:B:262:LEU:O	1:B:265:ASN:HB3	2.21	0.41
1:D:236:TRP:CD1	1:D:243:LEU:HD13	2.55	0.41
1:E:111:PHE:HB3	1:E:122:LEU:HB3	2.02	0.41
1:F:119:ARG:HD2	1:F:123:GLU:OE2	2.21	0.41
1:F:159:ASP:OD2	1:F:160:LYS:N	2.54	0.41
1:F:228:ARG:HH11	1:F:228:ARG:HB2	1.85	0.41
1:A:341:TYR:O	1:A:344:GLU:HB3	2.21	0.41
1:B:248:LEU:HB3	1:B:250:GLU:OE1	2.21	0.41
1:E:229:GLU:HB3	1:E:243:LEU:HD23	2.02	0.41
1:F:355:SER:O	1:F:356:SER:C	2.59	0.41
1:B:233:GLN:HB3	1:B:233:GLN:HE21	1.60	0.41
1:B:343:GLU:HA	1:B:367:ARG:NH2	2.36	0.41
1:D:54:PHE:O	1:D:58:ILE:HG13	2.20	0.41
1:D:80:ASP:OD1	2:D:401:ER4:H8	2.21	0.41
1:A:233:GLN:HE21	1:A:233:GLN:HB3	1.65	0.40
1:A:259:LEU:HD21	1:A:303:CYS:O	2.21	0.40
1:C:100:PHE:HA	1:C:103:PHE:CE2	2.55	0.40
1:C:162:VAL:HG11	1:C:168:TRP:HA	2.02	0.40
1:F:176:ALA:CB	1:F:212:GLN:HB2	2.51	0.40
1:B:153:GLY:HA3	1:B:174:TYR:CG	2.56	0.40
1:F:120:GLN:HG2	1:F:121:VAL:N	2.37	0.40
1:A:41:LYS:NZ	1:C:365:THR:HA	2.37	0.40
1:A:233:GLN:HA	1:A:236:TRP:CD1	2.56	0.40
1:D:252:ILE:HG23	1:D:253:ASP:N	2.36	0.40
1:F:194:ASP:OD1	1:F:195:PRO:HD2	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	325/360 (90%)	317 (98%)	8 (2%)	0	100	100
1	B	328/360 (91%)	315 (96%)	12 (4%)	1 (0%)	41	32
1	C	328/360 (91%)	318 (97%)	9 (3%)	1 (0%)	41	32
1	D	321/360 (89%)	304 (95%)	16 (5%)	1 (0%)	41	32
1	E	326/360 (91%)	315 (97%)	11 (3%)	0	100	100
1	F	324/360 (90%)	306 (94%)	16 (5%)	2 (1%)	25	15
All	All	1952/2160 (90%)	1875 (96%)	72 (4%)	5 (0%)	41	32

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	36	LEU
1	D	369	GLN
1	C	316	ILE
1	F	252	ILE
1	F	63	GLY

### 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	294/320 (92%)	289 (98%)	5 (2%)	60	57
1	B	296/320 (92%)	291 (98%)	5 (2%)	60	57
1	C	296/320 (92%)	291 (98%)	5 (2%)	60	57

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	291/320 (91%)	288 (99%)	3 (1%)	76	75
1	E	294/320 (92%)	284 (97%)	10 (3%)	37	30
1	F	292/320 (91%)	282 (97%)	10 (3%)	37	30
All	All	1763/1920 (92%)	1725 (98%)	38 (2%)	52	46

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	167	GLU
1	A	233	GLN
1	A	344	GLU
1	A	347	HIS
1	A	369	GLN
1	B	84	ASP
1	B	117	LYS
1	B	233	GLN
1	B	326	MET
1	B	351	ASP
1	C	84	ASP
1	C	110	ARG
1	C	203	ARG
1	C	233	GLN
1	C	322	VAL
1	D	203	ARG
1	D	228	ARG
1	D	233	GLN
1	E	49	GLN
1	E	52	ARG
1	E	84	ASP
1	E	108	ASP
1	E	112	MET
1	E	115	LYS
1	E	203	ARG
1	E	233	GLN
1	E	238	ARG
1	E	327	ASP
1	F	36	LEU
1	F	77	ARG
1	F	84	ASP
1	F	110	ARG
1	F	137	GLU

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Mol	Chain	Res	Type
1	F	228	ARG
1	F	233	GLN
1	F	241	LYS
1	F	326	MET
1	F	347	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	225	GLN
1	A	233	GLN
1	A	257	GLN
1	A	308	GLN
1	A	353	ASN
1	A	361	GLN
1	A	369	GLN
1	B	120	GLN
1	B	225	GLN
1	B	233	GLN
1	B	308	GLN
1	C	225	GLN
1	C	233	GLN
1	C	251	ASN
1	C	257	GLN
1	D	225	GLN
1	D	233	GLN
1	D	257	GLN
1	D	353	ASN
1	D	370	ASN
1	E	225	GLN
1	E	233	GLN
1	E	257	GLN
1	E	293	GLN
1	E	353	ASN
1	E	370	ASN
1	F	134	ASN
1	F	225	GLN
1	F	233	GLN
1	F	257	GLN
1	F	293	GLN
1	F	353	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](#)

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

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

6 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
2	ER4	A	901	-	31,33,33	2.38	4 (12%)	36,45,45	1.09	2 (5%)
2	ER4	C	401	-	31,33,33	2.47	4 (12%)	36,45,45	1.08	2 (5%)
2	ER4	F	401	-	31,33,33	2.53	4 (12%)	36,45,45	1.06	2 (5%)
2	ER4	B	401	-	31,33,33	2.36	3 (9%)	36,45,45	1.07	2 (5%)
2	ER4	E	401	-	31,33,33	2.52	4 (12%)	36,45,45	1.06	2 (5%)
2	ER4	D	401	-	31,33,33	2.43	4 (12%)	36,45,45	1.09	2 (5%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ER4	A	901	-	-	4/14/37/37	0/5/4/4
2	ER4	C	401	-	-	5/14/37/37	0/5/4/4

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ER4	F	401	-	-	5/14/37/37	0/5/4/4
2	ER4	B	401	-	-	5/14/37/37	0/5/4/4
2	ER4	E	401	-	-	5/14/37/37	0/5/4/4
2	ER4	D	401	-	-	4/14/37/37	0/5/4/4

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	401	ER4	CAS-CBA	-11.75	1.42	1.51
2	F	401	ER4	CAS-CBA	-11.61	1.42	1.51
2	C	401	ER4	CAS-CBA	-11.32	1.43	1.51
2	D	401	ER4	CAS-CBA	-11.18	1.43	1.51
2	A	901	ER4	CAS-CBA	-10.98	1.43	1.51
2	B	401	ER4	CAS-CBA	-10.89	1.43	1.51
2	F	401	ER4	CAS-CAX	-3.22	1.41	1.52
2	E	401	ER4	CAS-CAX	-3.14	1.42	1.52
2	D	401	ER4	CAZ-NAU	3.10	1.38	1.33
2	A	901	ER4	CAS-CAX	-3.07	1.42	1.52
2	C	401	ER4	CAS-CAX	-2.99	1.42	1.52
2	F	401	ER4	CAZ-NAU	2.97	1.38	1.33
2	D	401	ER4	CAS-CAX	-2.96	1.42	1.52
2	C	401	ER4	CAZ-NAU	2.95	1.38	1.33
2	A	901	ER4	CAZ-NAU	2.89	1.38	1.33
2	B	401	ER4	CAS-CAX	-2.81	1.43	1.52
2	B	401	ER4	CAZ-NAU	2.67	1.37	1.33
2	E	401	ER4	CAZ-NAU	2.63	1.37	1.33
2	F	401	ER4	CAH-CAX	2.25	1.43	1.38
2	D	401	ER4	CAH-CAX	2.22	1.43	1.38
2	C	401	ER4	CAH-CAX	2.21	1.43	1.38
2	E	401	ER4	CAH-CAX	2.16	1.43	1.38
2	A	901	ER4	CAH-CAX	2.13	1.43	1.38

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	ER4	CAZ-NAU-CBA	3.72	122.97	117.64
2	E	401	ER4	CAZ-NAU-CBA	3.70	122.94	117.64
2	C	401	ER4	CAZ-NAU-CBA	3.65	122.87	117.64
2	D	401	ER4	CAZ-NAU-CBA	3.63	122.84	117.64
2	A	901	ER4	CAZ-NAU-CBA	3.62	122.82	117.64
2	F	401	ER4	CAZ-NAU-CBA	3.49	122.63	117.64

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	ER4	CAJ-CAY-CAC	-2.92	114.48	120.11
2	C	401	ER4	CAJ-CAY-CAC	-2.91	114.49	120.11
2	A	901	ER4	CAJ-CAY-CAC	-2.90	114.51	120.11
2	D	401	ER4	CAJ-CAY-CAC	-2.83	114.64	120.11
2	E	401	ER4	CAJ-CAY-CAC	-2.65	115.00	120.11
2	F	401	ER4	CAJ-CAY-CAC	-2.62	115.05	120.11

There are no chirality outliers.

All (28) torsion outliers are listed below:

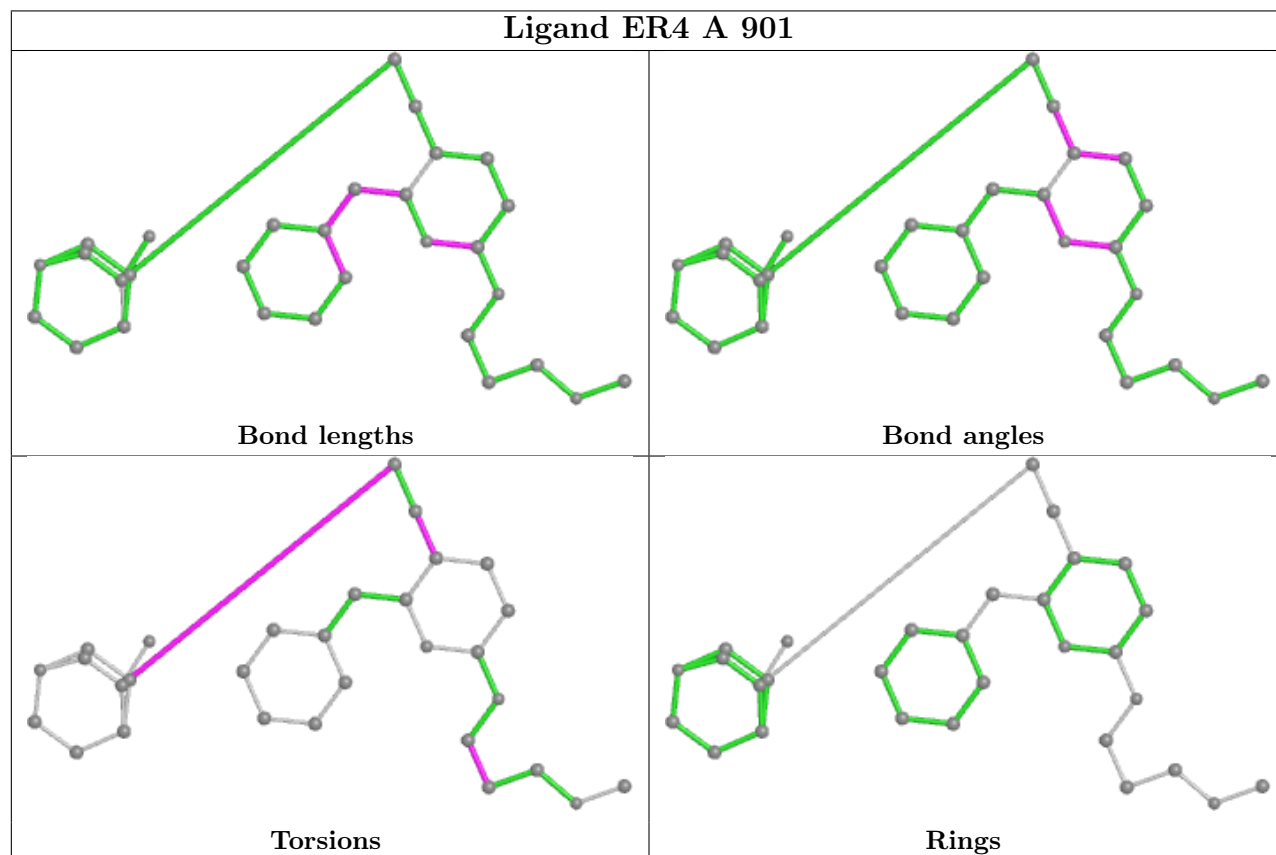
Mol	Chain	Res	Type	Atoms
2	F	401	ER4	CAC-CAD-CBD-CAT
2	B	401	ER4	CAD-CAC-CAY-CBA
2	B	401	ER4	CAY-CAC-CAD-CBD
2	D	401	ER4	CAY-CAC-CAD-CBD
2	E	401	ER4	CAD-CAC-CAY-CBA
2	A	901	ER4	CAD-CAC-CAY-CBA
2	F	401	ER4	CAD-CAC-CAY-CBA
2	B	401	ER4	CAD-CAC-CAY-CAJ
2	D	401	ER4	CAD-CAC-CAY-CBA
2	C	401	ER4	CAY-CAC-CAD-CBD
2	F	401	ER4	CAM-CAL-CAN-OAW
2	A	901	ER4	CAD-CAC-CAY-CAJ
2	C	401	ER4	CAD-CAC-CAY-CAJ
2	E	401	ER4	CAD-CAC-CAY-CAJ
2	F	401	ER4	CAD-CAC-CAY-CAJ
2	D	401	ER4	CAD-CAC-CAY-CAJ
2	A	901	ER4	CAM-CAL-CAN-OAW
2	E	401	ER4	CAN-CAL-CAM-OAV
2	C	401	ER4	CAD-CAC-CAY-CBA
2	D	401	ER4	CAM-CAL-CAN-OAW
2	B	401	ER4	CAN-CAL-CAM-OAV
2	C	401	ER4	CAM-CAL-CAN-OAW
2	B	401	ER4	CAM-CAL-CAN-OAW
2	E	401	ER4	CAM-CAL-CAN-OAW
2	A	901	ER4	CAC-CAD-CBD-CAT
2	C	401	ER4	CAC-CAD-CBD-CAT
2	E	401	ER4	CAC-CAD-CBD-CAT
2	F	401	ER4	CAL-CAM-OAV-CAA

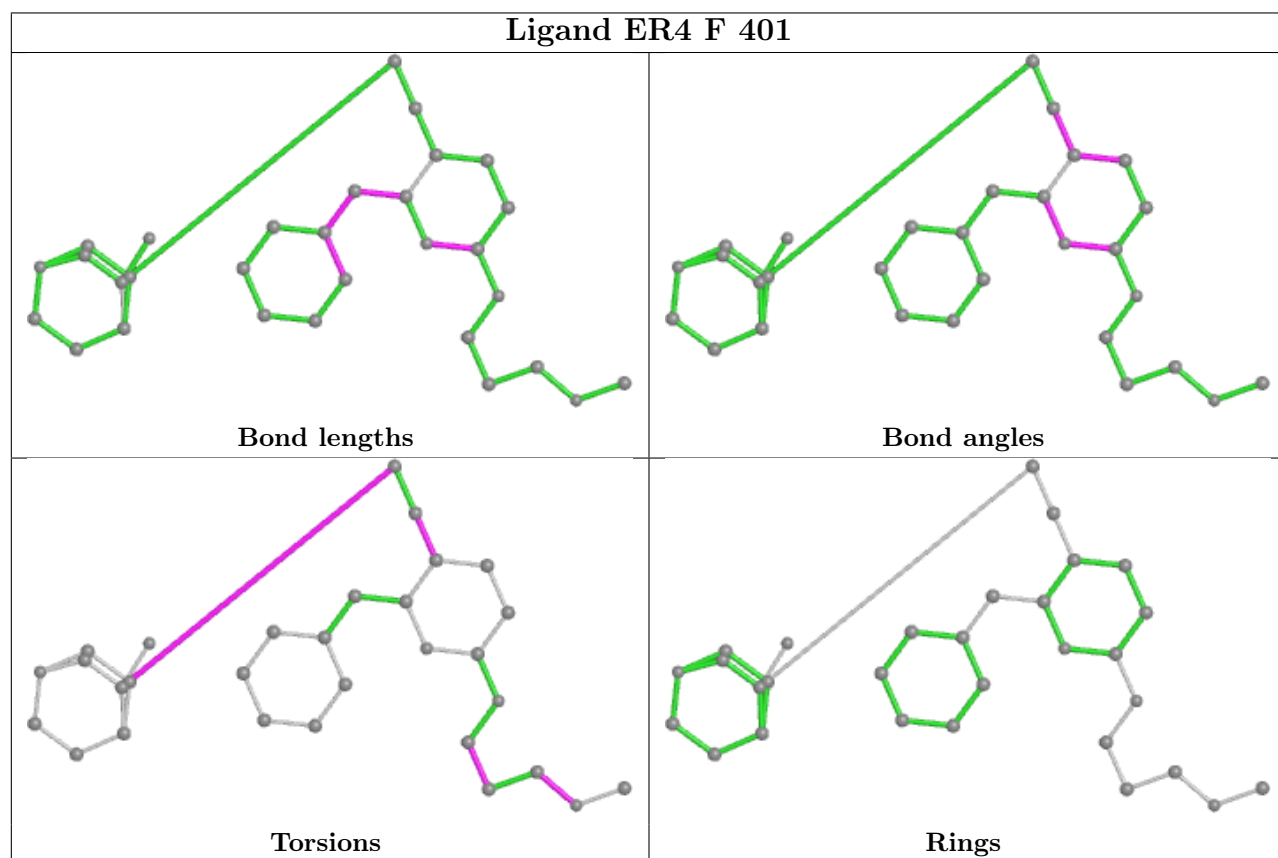
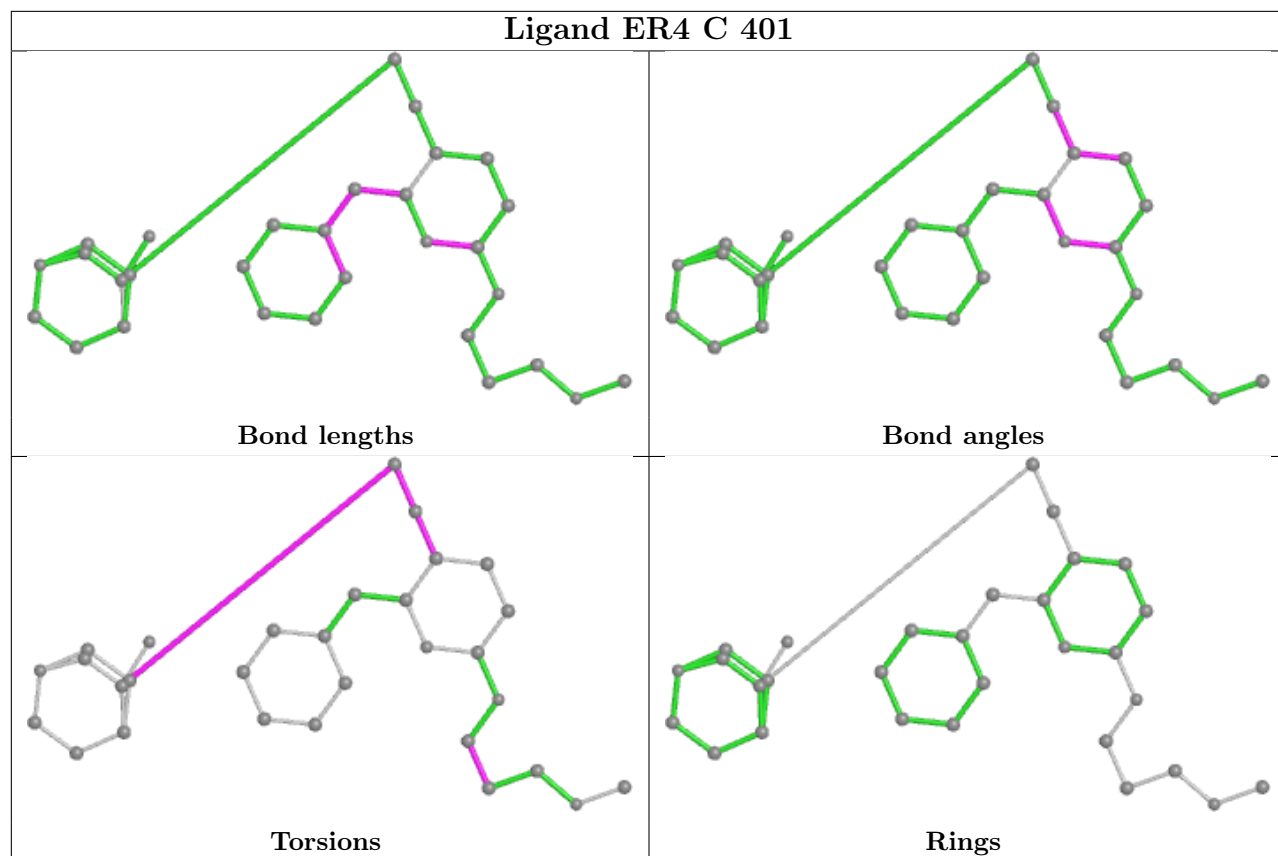
There are no ring outliers.

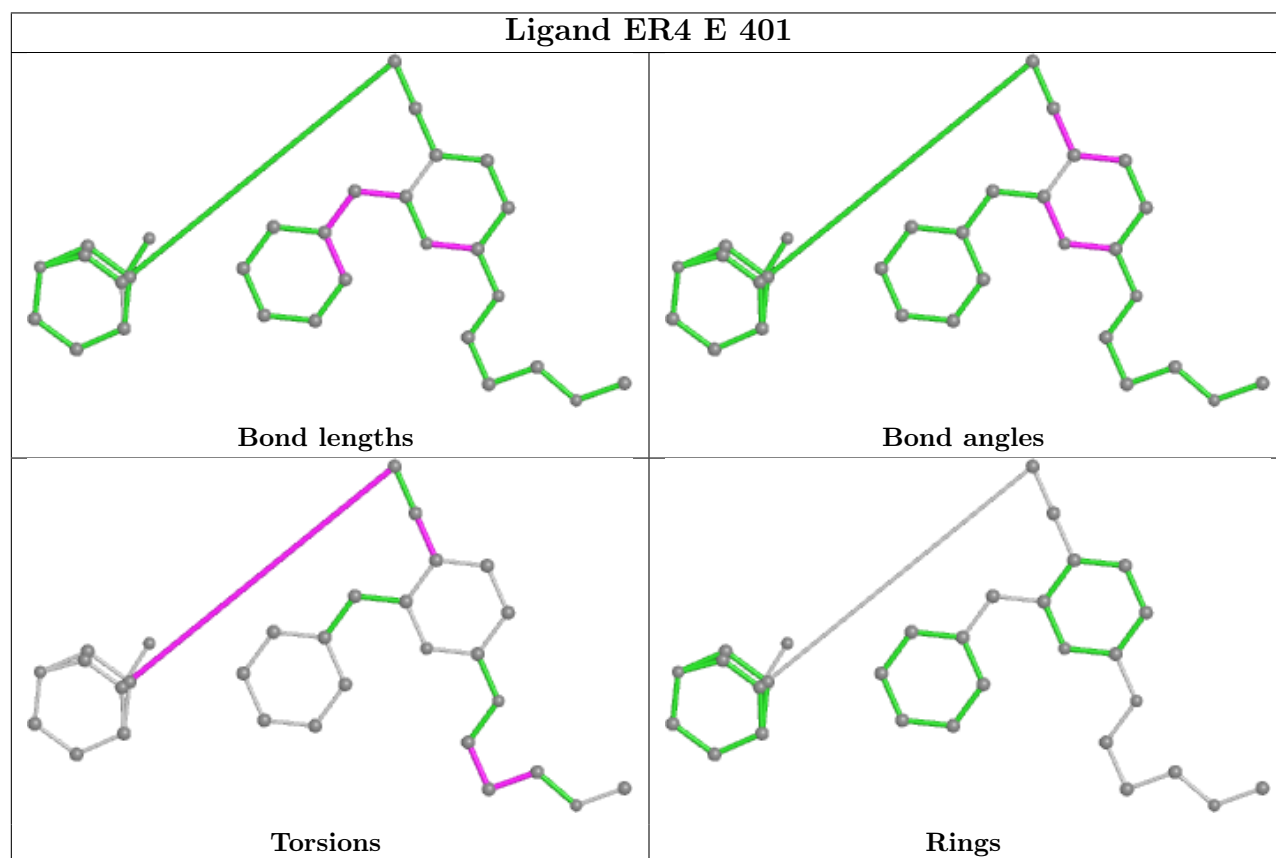
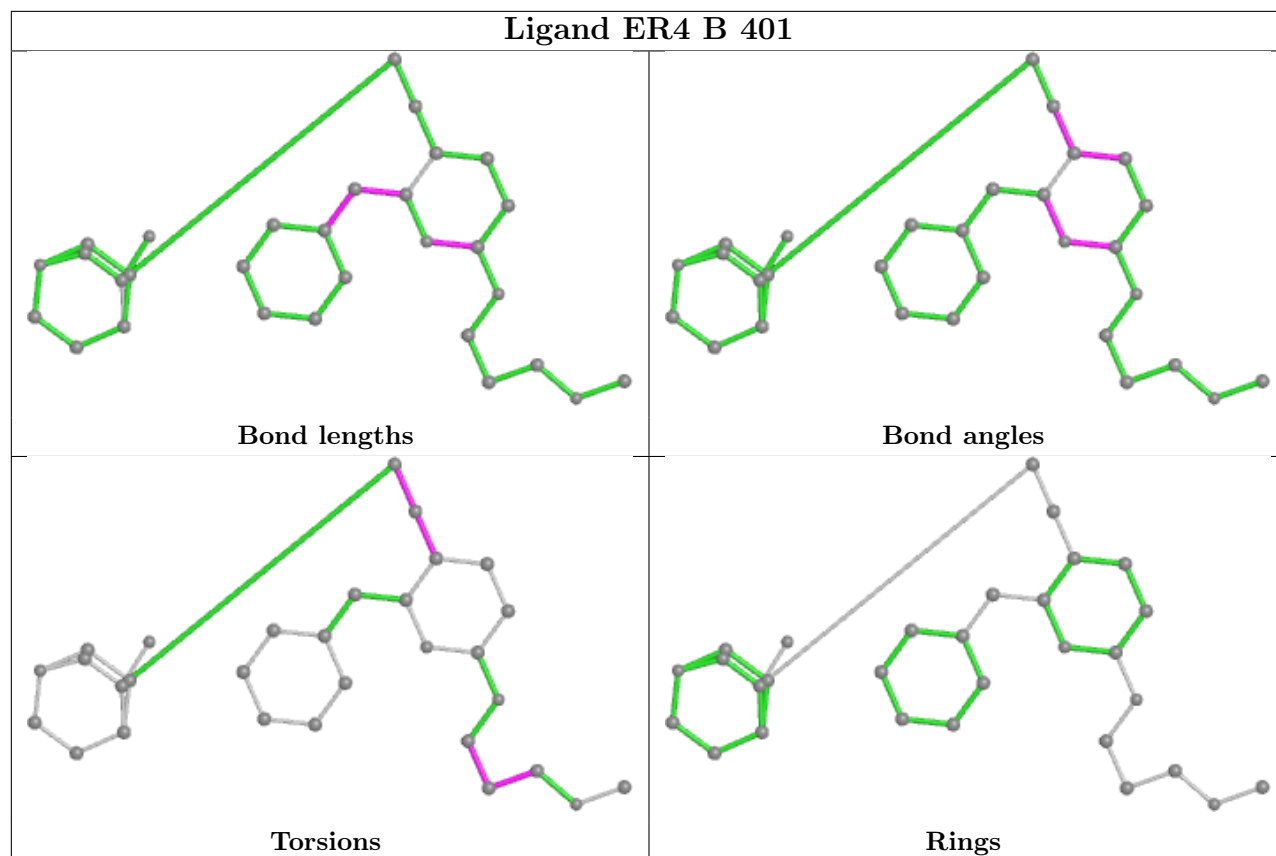
4 monomers are involved in 4 short contacts:

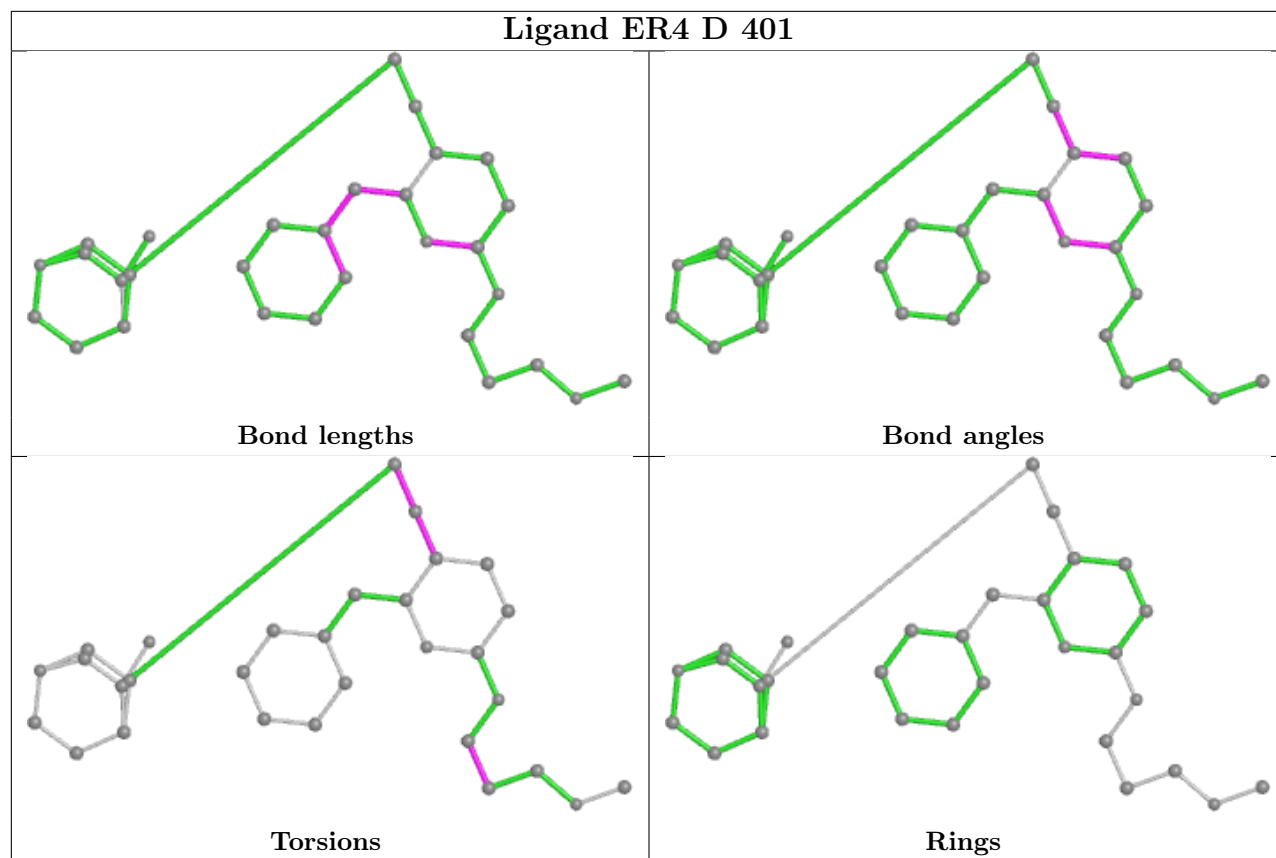
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	901	ER4	1	0
2	F	401	ER4	1	0
2	B	401	ER4	1	0
2	D	401	ER4	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.









## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.



## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	329/360 (91%)	-0.08	14 (4%) 35 36	13, 24, 50, 78	0
1	B	332/360 (92%)	0.26	24 (7%) 15 16	17, 28, 64, 79	0
1	C	332/360 (92%)	0.21	20 (6%) 21 22	18, 29, 57, 76	0
1	D	325/360 (90%)	0.26	25 (7%) 13 13	19, 33, 67, 80	0
1	E	330/360 (91%)	0.41	19 (5%) 23 24	23, 38, 65, 80	0
1	F	328/360 (91%)	0.90	59 (17%) 1 1	22, 49, 78, 90	0
All	All	1976/2160 (91%)	0.33	161 (8%) 12 12	13, 33, 70, 90	0

All (161) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	324	LEU	8.9
1	B	315	LEU	8.7
1	C	36	LEU	7.5
1	A	315	LEU	7.4
1	C	313	ALA	7.3
1	B	313	ALA	7.1
1	F	36	LEU	6.8
1	C	315	LEU	6.0
1	E	324	LEU	6.0
1	E	323	THR	5.9
1	F	248	LEU	5.9
1	F	323	THR	5.4
1	B	312	GLY	5.4
1	B	316	ILE	5.3
1	B	34	ASP	5.3
1	F	322	VAL	5.2
1	A	316	ILE	5.1
1	E	315	LEU	4.9
1	C	316	ILE	4.8

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	A	324	LEU	4.6
1	A	36	LEU	4.6
1	C	312	GLY	4.5
1	D	323	THR	4.5
1	B	35	SER	4.4
1	B	314	VAL	4.4
1	D	159	ASP	4.4
1	F	351	ASP	4.3
1	D	324	LEU	4.3
1	E	313	ALA	4.2
1	D	37	SER	4.1
1	F	37	SER	4.1
1	A	313	ALA	4.1
1	F	313	ALA	4.1
1	C	317	ARG	4.0
1	F	250	GLU	3.9
1	F	312	GLY	3.9
1	F	52	ARG	3.9
1	D	347	HIS	3.8
1	F	89	SER	3.7
1	E	316	ILE	3.7
1	B	323	THR	3.7
1	F	76	LEU	3.6
1	C	324	LEU	3.6
1	F	241	LYS	3.6
1	E	159	ASP	3.6
1	F	159	ASP	3.6
1	B	36	LEU	3.6
1	D	322	VAL	3.6
1	C	159	ASP	3.5
1	F	75	VAL	3.4
1	A	323	THR	3.4
1	D	352	SER	3.2
1	F	231	TRP	3.2
1	F	369	GLN	3.2
1	F	87	THR	3.2
1	F	314	VAL	3.1
1	B	241	LYS	3.1
1	D	241	LYS	3.1
1	D	52	ARG	3.1
1	C	248	LEU	3.1
1	B	259	LEU	3.1

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	F	79	LEU	3.0
1	F	242	LYS	3.0
1	F	347	HIS	3.0
1	F	221	LEU	3.0
1	F	311	LYS	3.0
1	E	370	ASN	3.0
1	B	159	ASP	3.0
1	C	52	ARG	2.9
1	F	69	VAL	2.9
1	C	37	SER	2.9
1	B	52	ARG	2.8
1	A	314	VAL	2.8
1	F	217	ILE	2.8
1	D	251	ASN	2.8
1	B	352	SER	2.8
1	F	91	GLU	2.8
1	F	115	LYS	2.8
1	C	323	THR	2.8
1	B	347	HIS	2.7
1	F	85	ASP	2.7
1	B	263	ILE	2.7
1	E	91	GLU	2.7
1	E	37	SER	2.7
1	F	92	LYS	2.6
1	F	327	ASP	2.6
1	F	163	THR	2.6
1	B	324	LEU	2.6
1	F	358	LYS	2.6
1	F	325	MET	2.6
1	D	351	ASP	2.6
1	F	161	HIS	2.6
1	B	242	LYS	2.6
1	C	225	GLN	2.6
1	D	327	ASP	2.6
1	F	137	GLU	2.5
1	F	352	SER	2.5
1	F	170	LYS	2.5
1	F	307	GLN	2.5
1	A	325	MET	2.5
1	F	118	ASP	2.5
1	B	228	ARG	2.4
1	C	228	ARG	2.4

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	E	194	ASP	2.4
1	D	228	ARG	2.4
1	D	369	GLN	2.4
1	F	86	MET	2.4
1	A	248	LEU	2.4
1	D	309	VAL	2.4
1	D	225	GLN	2.4
1	D	263	ILE	2.4
1	B	256	VAL	2.3
1	A	327	ASP	2.3
1	E	312	GLY	2.3
1	B	368	THR	2.3
1	A	52	ARG	2.3
1	C	226	GLY	2.3
1	F	263	ILE	2.3
1	B	321	ALA	2.3
1	F	240	VAL	2.3
1	A	159	ASP	2.3
1	E	92	LYS	2.3
1	F	237	SER	2.3
1	D	325	MET	2.3
1	E	347	HIS	2.3
1	F	228	ARG	2.3
1	D	91	GLU	2.3
1	B	358	LYS	2.2
1	F	235	VAL	2.2
1	F	108	ASP	2.2
1	A	76	LEU	2.2
1	E	321	ALA	2.2
1	F	110	ARG	2.2
1	D	370	ASN	2.2
1	A	347	HIS	2.2
1	F	53	SER	2.2
1	F	90	VAL	2.2
1	C	241	LYS	2.2
1	D	248	LEU	2.2
1	F	158	LEU	2.2
1	E	38	SER	2.2
1	B	250	GLU	2.2
1	C	322	VAL	2.2
1	D	86	MET	2.2
1	E	86	MET	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	249	PRO	2.2
1	D	224	GLN	2.2
1	C	250	GLU	2.1
1	C	314	VAL	2.1
1	F	167	GLU	2.1
1	F	310	PHE	2.1
1	E	87	THR	2.1
1	C	227	GLY	2.1
1	E	75	VAL	2.1
1	D	368	THR	2.1
1	F	246	PHE	2.1
1	F	113	GLU	2.0
1	F	239	TYR	2.0
1	F	236	TRP	2.0
1	E	281	ARG	2.0
1	F	259	LEU	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](#)

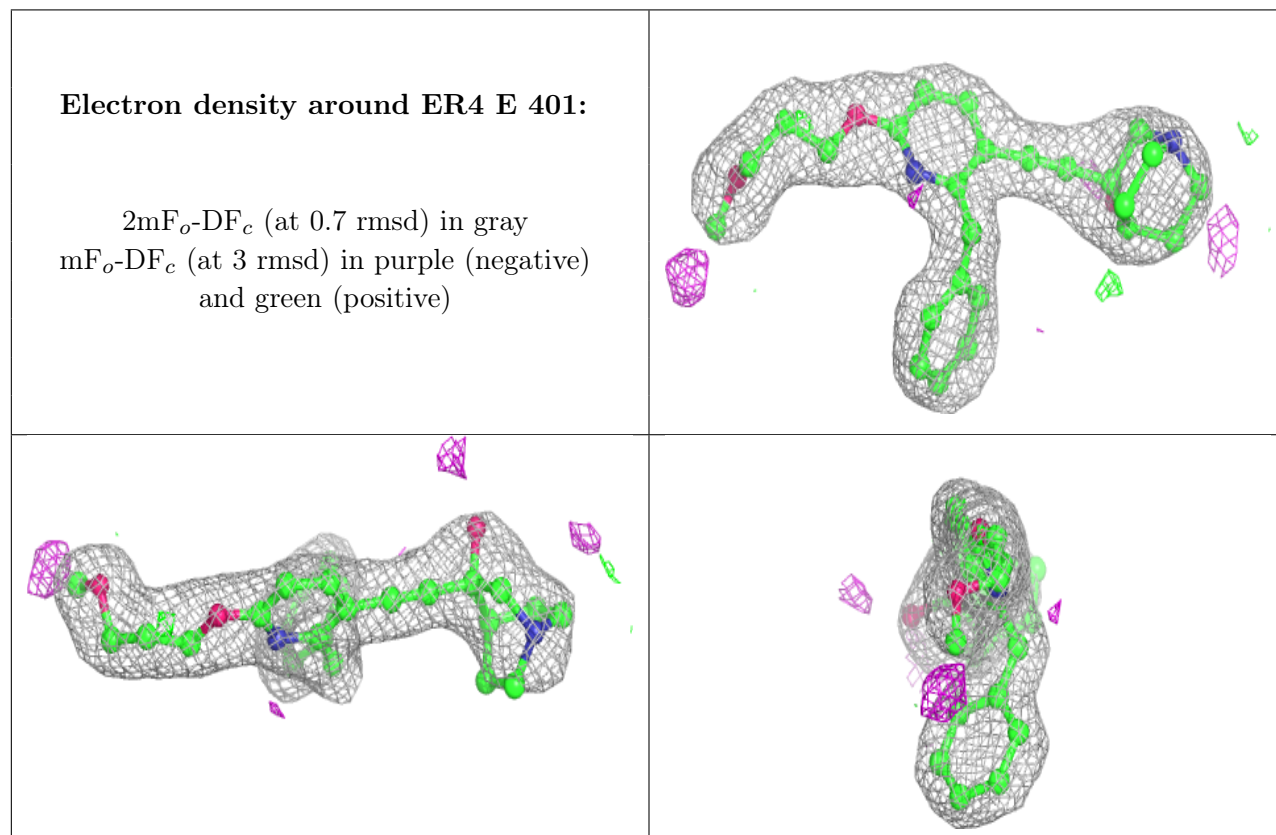
There are no monosaccharides in this entry.

## 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<sup>th</sup> 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.

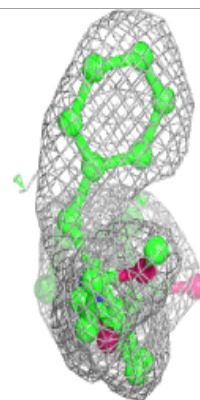
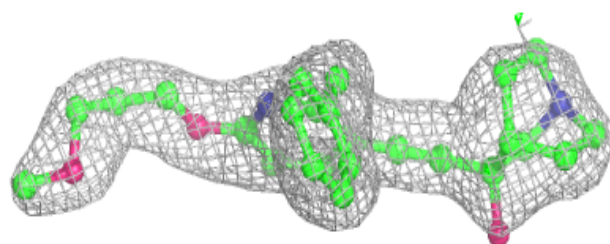
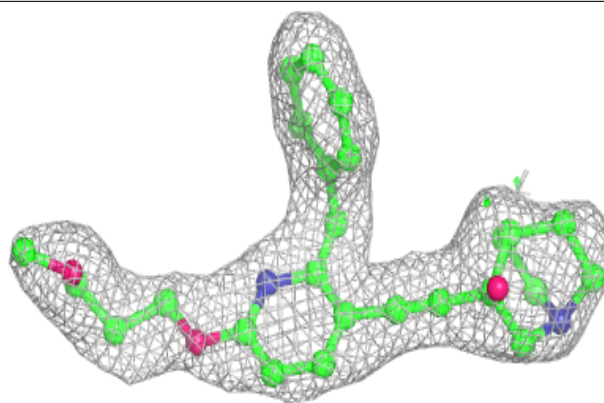
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	ER4	E	401	30/30	0.90	0.17	33,38,58,59	0
2	ER4	F	401	30/30	0.91	0.19	39,48,61,61	0
2	ER4	B	401	30/30	0.92	0.15	19,29,50,51	0
2	ER4	D	401	30/30	0.94	0.14	24,30,44,44	0
2	ER4	C	401	30/30	0.95	0.16	23,29,48,48	0
2	ER4	A	901	30/30	0.96	0.14	20,25,44,45	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.

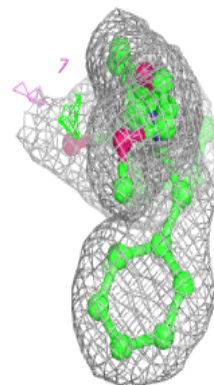
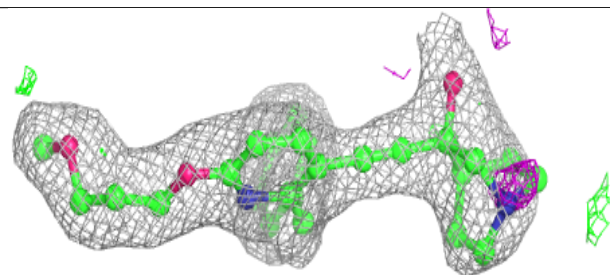
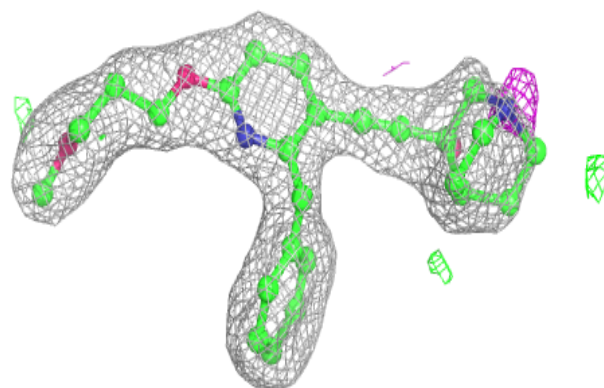


**Electron density around ER4 F 401:**

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

**Electron density around ER4 B 401:**

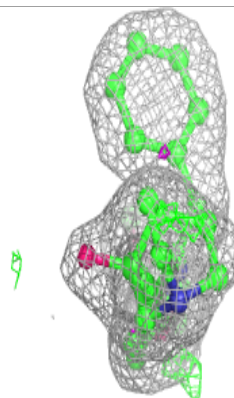
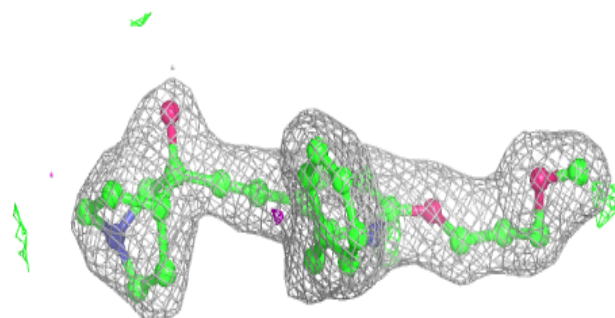
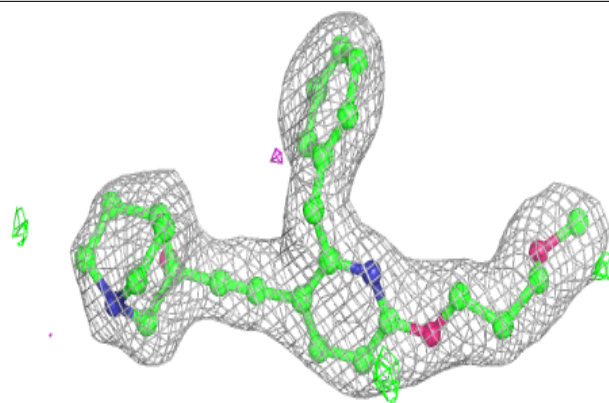
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



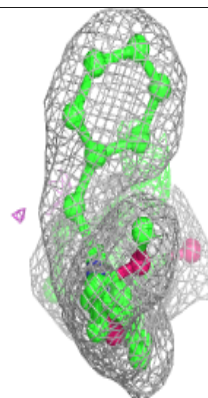
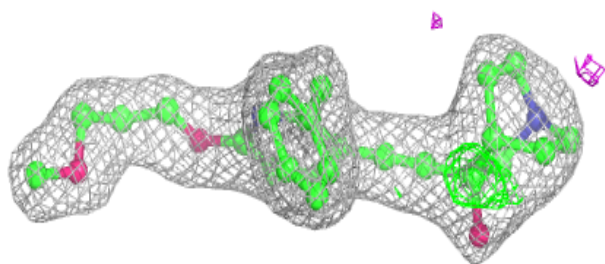
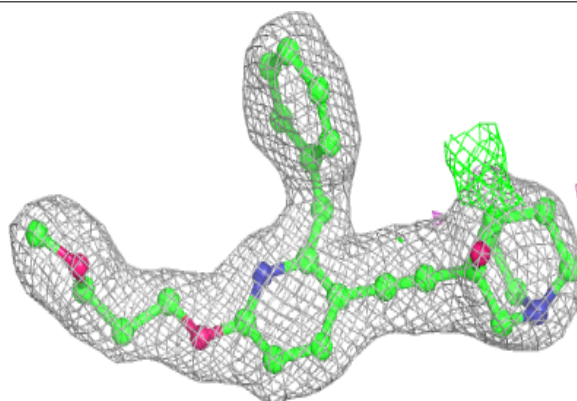


**Electron density around ER4 D 401:**

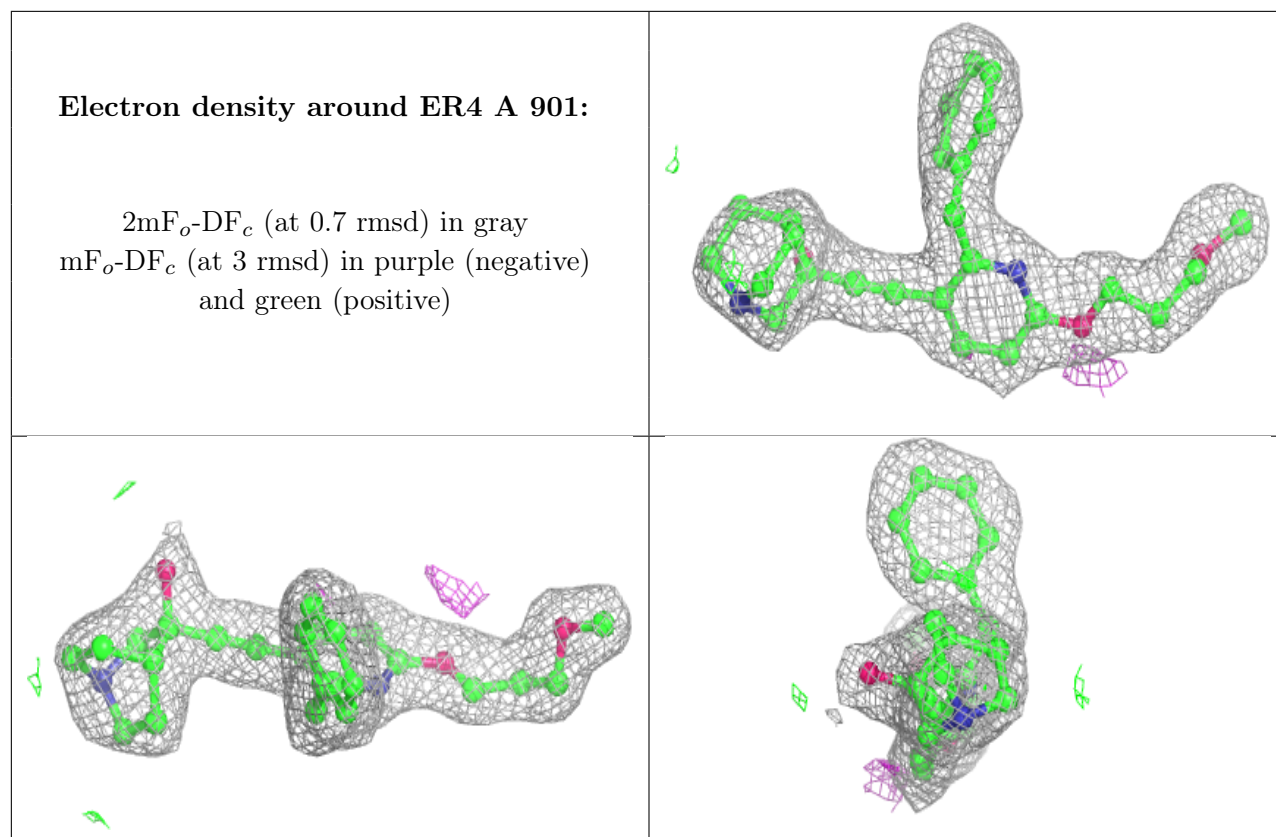
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around ER4 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)







## 6.5 Other polymers [i](#)

There are no such residues in this entry.