



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 03:46 PM UTC

PDB ID : 3WCL / pdb_00003wcl
Title : The complex structure of HsSQS wtih ligand,BPH1344
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 : 2.24 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 ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

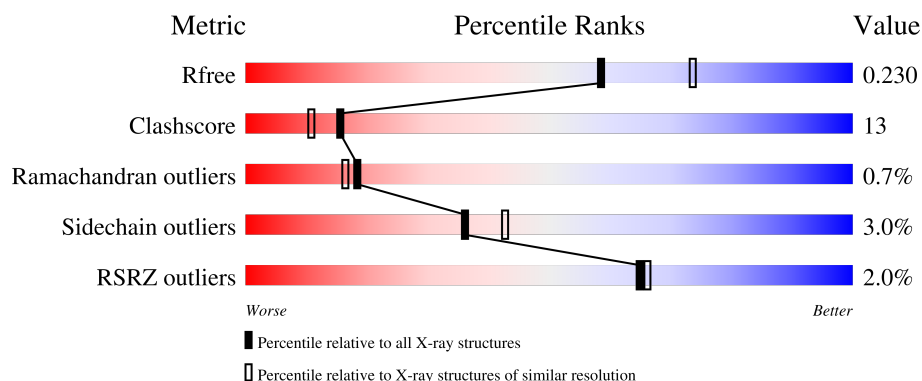
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.24 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 $\leq 5\%$. 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	3% 73% 18% • 7%
1	B	360	2% 72% 20% • 7%
1	C	360	70% 21% • 7%
1	D	360	0% 64% 27% • 7%
1	E	360	2% 66% 24% • 7%

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Mol	Chain	Length	Quality of chain
1	F	360	3% 62% 27% • • 7%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 17580 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
1	A	334	Total	C	N	O	S	0	0	0
			2696	1717	458	503	18			
1	B	334	Total	C	N	O	S	0	0	0
			2696	1717	458	503	18			
1	C	334	Total	C	N	O	S	0	0	0
			2696	1717	458	503	18			
1	D	334	Total	C	N	O	S	0	0	0
			2696	1717	458	503	18			
1	E	334	Total	C	N	O	S	0	0	0
			2696	1717	458	503	18			
1	F	334	Total	C	N	O	S	0	0	0
			2696	1717	458	503	18			

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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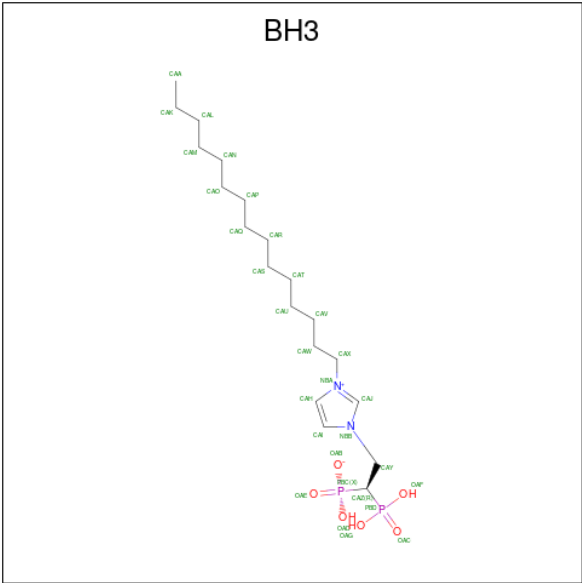
Chain	Residue	Modelled	Actual	Comment	Reference
E	16	HIS	-	expression tag	UNP P37268
E	17	HIS	-	expression tag	UNP P37268
E	18	HIS	-	expression tag	UNP P37268
E	19	HIS	-	expression tag	UNP P37268
E	20	HIS	-	expression tag	UNP P37268
E	21	SER	-	expression tag	UNP P37268
E	22	SER	-	expression tag	UNP P37268
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 hydrogen [(1R)-2-(3-pentadecyl-1H-imidazol-3-ium-1-yl)-1-phosphonoethyl]phosphonate (CCD ID: BH3) (formula: C₂₀H₄₀N₂O₆P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			30	20	2	6	2		

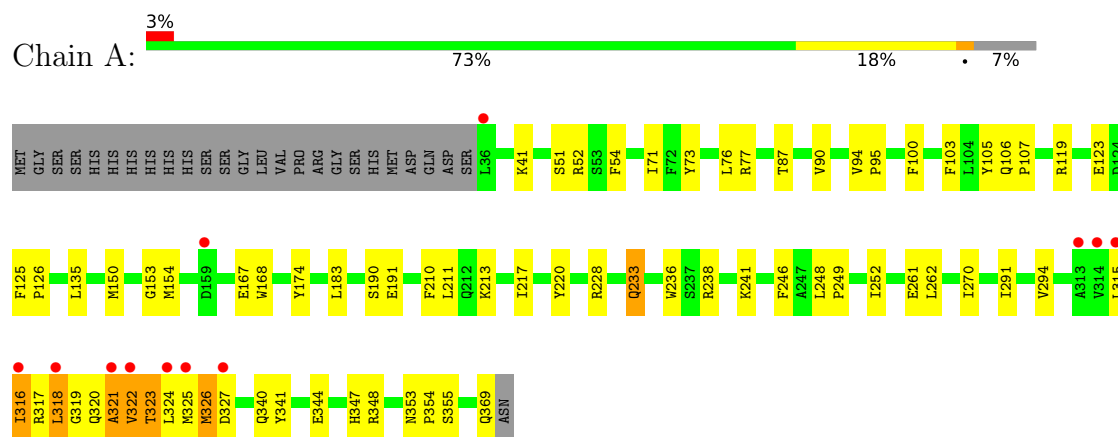
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	284	Total	O	0	0
			284	284		
3	B	285	Total	O	0	0
			285	285		
3	C	285	Total	O	0	0
			285	285		
3	D	201	Total	O	0	0
			201	201		
3	E	157	Total	O	0	0
			157	157		
3	F	162	Total	O	0	0
			162	162		

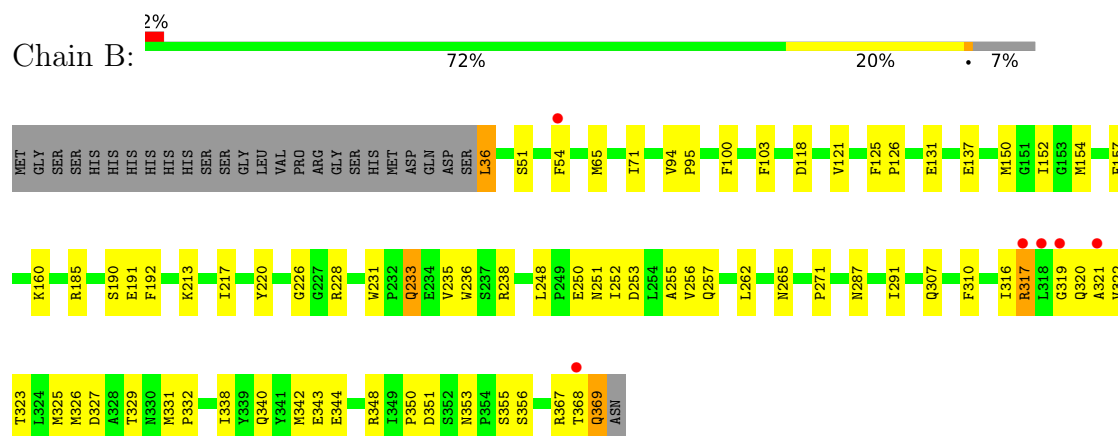
3 Residue-property plots [i](#)

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.

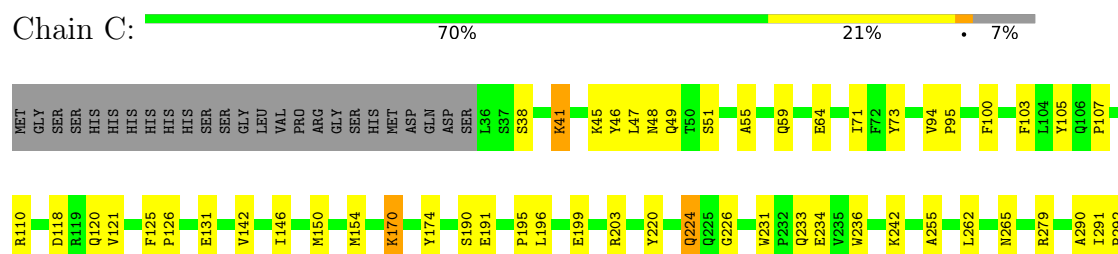
• Molecule 1: Squalene synthase



• Molecule 1: Squalene synthase

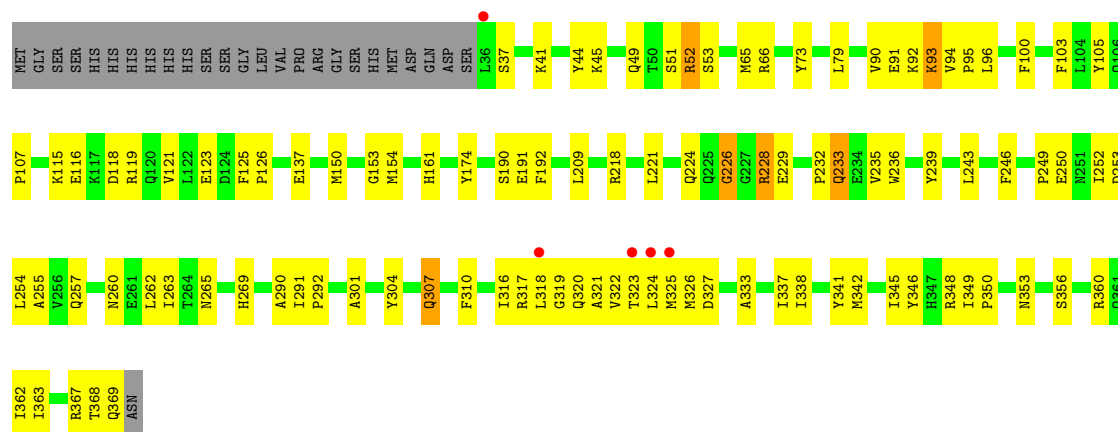


• Molecule 1: Squalene synthase

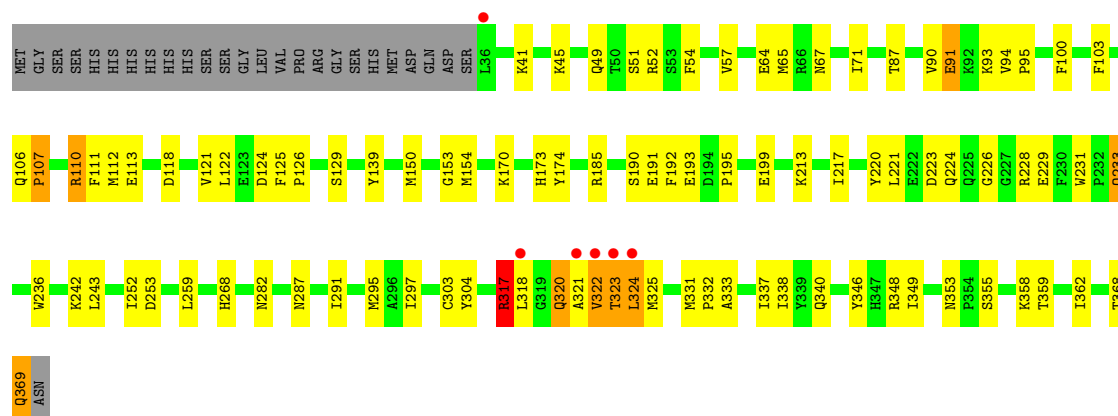




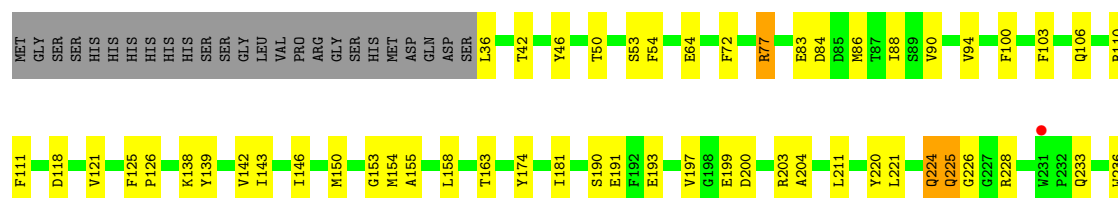
• Molecule 1: Squalene synthase

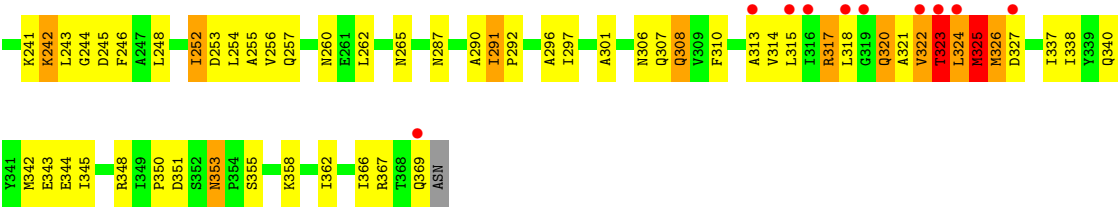


• Molecule 1: Squalene synthase



• Molecule 1: Squalene synthase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	85.25Å 153.30Å 90.93Å 90.00° 91.47° 90.00°	Depositor
Resolution (Å)	25.00 – 2.24 25.00 – 2.24	Depositor EDS
% Data completeness (in resolution range)	(Not available) (25.00-2.24) 95.1 (25.00-2.24)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.06 (at 2.24Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.188 , 0.229 0.189 , 0.230	Depositor DCC
R_{free} test set	5365 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	29.8	Xtriage
Anisotropy	0.624	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 45.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.029 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	17580	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.18% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: BH3

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	0.44	0/2751	0.90	6/3724 (0.2%)
1	B	0.42	0/2751	0.87	0/3724
1	C	0.41	0/2751	0.87	2/3724 (0.1%)
1	D	0.40	0/2751	0.85	2/3724 (0.1%)
1	E	0.39	0/2751	0.87	5/3724 (0.1%)
1	F	0.37	0/2751	0.87	12/3724 (0.3%)
All	All	0.40	0/16506	0.87	27/22344 (0.1%)

There are no bond length outliers.

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	290	ALA	N-CA-C	6.52	118.18	111.14
1	F	353	ASN	CA-C-N	6.37	126.06	119.82
1	F	353	ASN	C-N-CA	6.37	126.06	119.82
1	F	54	PHE	N-CA-C	6.31	121.02	112.88
1	A	323	THR	N-CA-C	-6.21	105.53	113.23
1	E	106	GLN	CA-C-N	6.05	127.40	119.84
1	E	106	GLN	C-N-CA	6.05	127.40	119.84
1	F	248	LEU	CA-C-N	5.94	126.36	119.47
1	F	248	LEU	C-N-CA	5.94	126.36	119.47
1	E	87	THR	N-CA-C	-5.89	106.09	113.28
1	D	290	ALA	N-CA-C	5.82	117.42	111.14
1	E	221	LEU	N-CA-C	5.80	117.29	110.97
1	F	290	ALA	N-CA-C	5.69	117.28	111.14
1	A	317	ARG	N-CA-C	5.61	118.72	110.59
1	F	291	ILE	CB-CA-C	-5.60	108.38	113.70
1	C	320	GLN	N-CA-C	5.57	118.17	108.76
1	D	93	LYS	N-CA-C	5.46	117.04	111.14
1	F	324	LEU	N-CA-C	-5.38	106.75	113.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	106	GLN	CA-C-N	5.34	126.52	119.84
1	F	106	GLN	C-N-CA	5.34	126.52	119.84
1	F	325	MET	N-CA-C	-5.33	105.62	113.61
1	A	106	GLN	CA-C-N	5.28	126.44	119.84
1	A	106	GLN	C-N-CA	5.28	126.44	119.84
1	A	87	THR	N-CA-C	-5.25	106.87	113.28
1	E	323	THR	N-CA-C	-5.24	106.73	113.02
1	A	168	TRP	N-CA-C	-5.11	105.63	111.14
1	F	204	ALA	N-CA-C	-5.07	105.83	111.36

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	2696	0	2676	64	0
1	B	2696	0	2676	72	0
1	C	2696	0	2676	68	0
1	D	2696	0	2676	77	0
1	E	2696	0	2676	75	0
1	F	2696	0	2676	101	0
2	A	30	0	40	13	0
3	A	284	0	0	3	0
3	B	285	0	0	5	0
3	C	285	0	0	2	0
3	D	201	0	0	2	0
3	E	157	0	0	8	0
3	F	162	0	0	6	0
All	All	17580	0	16096	433	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 13.

All (433) 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:325:MET:HE1	3:E:555:HOH:O	1.18	1.30
1:F:301:ALA:HA	1:F:345:ILE:HD11	1.27	1.12
1:E:229:GLU:HG2	1:E:243:LEU:HD23	1.41	1.01
1:B:326:MET:HE3	1:C:291:ILE:HD11	1.46	0.94
1:F:233:GLN:HA	1:F:236:TRP:NE1	1.83	0.94
1:F:320:GLN:NE2	1:F:322:VAL:H	1.65	0.93
1:D:301:ALA:HA	1:D:345:ILE:HD11	1.53	0.90
1:A:318:LEU:HD11	1:A:324:LEU:HD11	1.50	0.90
1:A:316:ILE:HD13	1:A:316:ILE:H	1.36	0.90
1:B:317:ARG:NH1	1:B:348:ARG:HH12	1.70	0.89
1:D:260:ASN:HD22	1:D:353:ASN:ND2	1.70	0.89
1:B:317:ARG:HH12	1:B:348:ARG:NH1	1.71	0.87
1:A:322:VAL:HB	1:A:340:GLN:HE22	1.38	0.86
1:F:260:ASN:HD22	1:F:353:ASN:ND2	1.73	0.86
1:B:317:ARG:HH12	1:B:348:ARG:HH12	1.20	0.84
1:B:226:GLY:HA3	1:B:228:ARG:HH12	1.42	0.84
1:F:301:ALA:HA	1:F:345:ILE:CD1	2.08	0.83
2:A:401:BH3:H16	3:A:783:HOH:O	1.79	0.83
1:A:320:GLN:O	1:A:323:THR:HG22	1.78	0.82
1:F:320:GLN:HE22	1:F:322:VAL:H	1.28	0.81
1:C:323:THR:HG23	1:C:337:ILE:HD12	1.65	0.78
1:D:252:ILE:HD11	1:D:307:GLN:HB3	1.65	0.78
1:D:37:SER:O	1:D:41:LYS:HG3	1.84	0.77
1:B:353:ASN:ND2	1:B:355:SER:H	1.82	0.77
1:E:268:HIS:HA	1:E:362:ILE:HD11	1.65	0.77
1:A:77:ARG:HD2	2:A:401:BH3:H11	1.66	0.76
1:C:45:LYS:O	1:C:49:GLN:HG3	1.84	0.76
1:F:291:ILE:HB	1:F:292:PRO:HD3	1.68	0.76
1:A:213:LYS:O	1:A:217:ILE:HG12	1.86	0.76
1:D:325:MET:CE	3:E:555:HOH:O	1.92	0.76
1:F:287:ASN:O	1:F:291:ILE:HG12	1.87	0.75
1:B:233:GLN:HA	1:B:236:TRP:NE1	2.01	0.74
1:F:350:PRO:HG2	1:F:353:ASN:HB2	1.70	0.74
1:E:325:MET:O	1:F:327:ASP:HB2	1.88	0.74
1:D:368:THR:O	1:D:369:GLN:HB2	1.86	0.73
1:C:362:ILE:O	1:C:366:ILE:HD13	1.88	0.73
1:F:320:GLN:HE22	1:F:322:VAL:HG23	1.53	0.73
1:C:170:LYS:HD2	1:C:174:TYR:CE2	2.25	0.72
1:B:235:VAL:HA	1:B:238:ARG:HH21	1.55	0.72
1:E:320:GLN:O	1:E:324:LEU:HB2	1.90	0.72
1:D:325:MET:SD	3:E:555:HOH:O	2.39	0.72
1:B:226:GLY:HA3	1:B:228:ARG:NH1	2.04	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:321:ALA:O	1:A:322:VAL:HG23	1.89	0.71
1:D:233:GLN:HA	1:D:236:TRP:NE1	2.05	0.71
1:A:322:VAL:HB	1:A:340:GLN:NE2	2.06	0.70
1:A:353:ASN:ND2	1:A:355:SER:H	1.89	0.70
1:E:353:ASN:ND2	1:E:355:SER:H	1.89	0.70
1:C:297:ILE:CD1	1:C:338:ILE:HG12	2.21	0.70
1:A:326:MET:HA	1:B:327:ASP:HB2	1.75	0.69
1:D:45:LYS:O	1:D:49:GLN:HG3	1.93	0.69
1:C:94:VAL:HB	1:C:95:PRO:HD3	1.74	0.69
1:F:100:PHE:HA	1:F:103:PHE:CD2	2.28	0.69
1:A:315:LEU:O	1:A:315:LEU:HD12	1.92	0.69
1:E:323:THR:HB	1:E:340:GLN:OE1	1.93	0.68
1:E:323:THR:OG1	1:E:337:ILE:HG12	1.93	0.68
1:B:36:LEU:N	1:B:36:LEU:HD12	2.08	0.68
1:D:119:ARG:O	1:D:123:GLU:HG3	1.94	0.68
1:C:353:ASN:ND2	1:C:355:SER:H	1.91	0.68
1:E:45:LYS:O	1:E:49:GLN:HG3	1.93	0.68
1:D:229:GLU:HG3	1:D:243:LEU:HD23	1.76	0.67
1:E:195:PRO:O	1:E:199:GLU:HG3	1.94	0.67
1:D:116:GLU:O	1:D:119:ARG:HD3	1.95	0.67
1:E:94:VAL:HB	1:E:95:PRO:HD3	1.76	0.67
1:C:318:LEU:HD12	1:C:320:GLN:HG3	1.77	0.67
1:E:125:PHE:N	1:E:126:PRO:HD2	2.10	0.67
1:D:94:VAL:HB	1:D:95:PRO:HD3	1.77	0.67
1:E:320:GLN:H	1:E:323:THR:CG2	2.08	0.67
1:F:260:ASN:HD22	1:F:353:ASN:HD22	1.41	0.67
1:E:320:GLN:H	1:E:323:THR:HG23	1.59	0.66
1:E:368:THR:O	1:E:369:GLN:HB2	1.93	0.66
1:B:343:GLU:OE1	1:B:367:ARG:NH2	2.28	0.66
1:F:306:ASN:OD1	1:F:308:GLN:HB2	1.96	0.66
1:D:229:GLU:HG3	1:D:243:LEU:CD2	2.26	0.65
1:B:348:ARG:HG3	1:B:348:ARG:HH11	1.61	0.65
1:F:150:MET:HG3	1:F:174:TYR:O	1.96	0.64
1:A:294:VAL:HG11	1:C:325:MET:HE2	1.80	0.64
1:B:320:GLN:HB3	1:B:340:GLN:HE22	1.63	0.64
1:D:228:ARG:HB2	1:D:228:ARG:NH1	2.13	0.64
1:C:105:TYR:O	1:C:107:PRO:HD3	1.97	0.64
1:A:211:LEU:HD12	2:A:401:BH3:CAO	2.29	0.63
1:E:100:PHE:HA	1:E:103:PHE:CD2	2.33	0.63
1:D:100:PHE:HA	1:D:103:PHE:CD2	2.34	0.63
1:A:233:GLN:HA	1:A:236:TRP:NE1	2.13	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:118:ASP:O	1:E:121:VAL:HG22	1.97	0.63
1:D:65:MET:HE2	1:D:192:PHE:CD2	2.34	0.62
1:B:317:ARG:HD2	3:B:555:HOH:O	1.99	0.62
1:F:90:VAL:O	1:F:94:VAL:HG23	1.98	0.62
1:F:343:GLU:CD	1:F:367:ARG:HH21	2.08	0.62
1:E:57:VAL:HG22	3:E:410:HOH:O	1.99	0.62
1:C:51:SER:HB2	1:C:73:TYR:CE1	2.35	0.62
1:A:347:HIS:HB2	3:B:677:HOH:O	2.00	0.61
1:A:316:ILE:H	1:A:316:ILE:CD1	2.13	0.61
1:D:252:ILE:HD11	1:D:307:GLN:CB	2.30	0.61
1:A:322:VAL:O	1:A:326:MET:HG2	2.00	0.61
1:E:110:ARG:HH11	1:E:112:MET:HE2	1.65	0.60
1:F:322:VAL:O	1:F:323:THR:HG23	2.01	0.60
1:B:325:MET:HE2	1:C:295:MET:HG3	1.81	0.60
1:B:326:MET:HE3	1:C:291:ILE:CD1	2.25	0.60
1:D:327:ASP:HB2	1:F:325:MET:O	2.01	0.60
1:D:150:MET:HG3	1:D:174:TYR:O	2.01	0.60
1:F:236:TRP:CG	1:F:243:LEU:HD13	2.36	0.60
1:E:323:THR:HG21	3:E:556:HOH:O	2.00	0.59
1:A:325:MET:O	1:B:327:ASP:HB2	2.03	0.59
1:B:271:PRO:HG2	3:B:674:HOH:O	2.01	0.59
1:E:64:GLU:HG2	3:E:548:HOH:O	2.03	0.59
1:C:308:GLN:OE1	1:C:311:LYS:HD2	2.03	0.59
1:D:291:ILE:HB	1:D:292:PRO:HD3	1.84	0.59
1:F:83:GLU:OE1	1:F:154:MET:HE2	2.03	0.59
1:B:320:GLN:HB3	1:B:340:GLN:NE2	2.18	0.59
1:E:91:GLU:H	1:E:91:GLU:CD	2.11	0.59
1:E:233:GLN:HA	1:E:236:TRP:NE1	2.17	0.59
1:F:262:LEU:O	1:F:265:ASN:HB3	2.03	0.59
1:F:64:GLU:HG3	3:F:422:HOH:O	2.02	0.59
1:C:150:MET:HG3	1:C:174:TYR:O	2.03	0.58
1:B:319:GLY:HA3	1:B:323:THR:HG21	1.85	0.58
1:E:321:ALA:HA	1:E:324:LEU:HB2	1.84	0.58
1:A:294:VAL:CG1	1:C:325:MET:HE2	2.33	0.58
1:A:326:MET:HA	1:B:327:ASP:CB	2.33	0.58
1:E:268:HIS:CG	1:E:362:ILE:HD12	2.39	0.58
1:F:255:ALA:HB1	1:F:310:PHE:CZ	2.39	0.58
1:F:221:LEU:O	1:F:224:GLN:HB2	2.04	0.58
1:D:368:THR:HG21	1:E:41:LYS:HA	1.86	0.58
1:F:323:THR:C	1:F:325:MET:H	2.10	0.58
1:E:268:HIS:ND1	1:E:362:ILE:HD12	2.18	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:94:VAL:HB	1:A:95:PRO:HD3	1.85	0.57
1:D:301:ALA:CA	1:D:345:ILE:HD11	2.29	0.57
1:A:52:ARG:HG2	2:A:401:BH3:OAB	2.04	0.57
1:E:322:VAL:HG21	1:F:53:SER:O	2.05	0.57
1:F:118:ASP:O	1:F:121:VAL:HG22	2.04	0.57
1:C:348:ARG:O	1:C:350:PRO:HD3	2.05	0.57
1:A:100:PHE:HA	1:A:103:PHE:CD2	2.40	0.57
1:C:55:ALA:O	1:C:59:GLN:HG3	2.05	0.57
1:B:322:VAL:HB	1:B:340:GLN:OE1	2.05	0.56
1:F:322:VAL:O	1:F:322:VAL:HG12	2.04	0.56
1:C:195:PRO:O	1:C:199:GLU:HG3	2.06	0.56
1:C:320:GLN:O	1:C:321:ALA:HB3	2.05	0.56
1:C:326:MET:HE2	1:C:333:ALA:CB	2.36	0.56
1:D:318:LEU:HD23	1:D:319:GLY:N	2.21	0.56
1:D:233:GLN:HA	1:D:236:TRP:CD1	2.41	0.56
1:E:331:MET:HB3	1:E:332:PRO:HD3	1.87	0.56
1:D:52:ARG:CD	1:D:53:SER:H	2.19	0.55
1:F:163:THR:HA	1:F:233:GLN:HB3	1.89	0.55
1:C:64:GLU:HG3	3:C:429:HOH:O	2.07	0.55
1:F:224:GLN:OE1	1:F:244:GLY:HA2	2.06	0.55
1:B:331:MET:HB3	1:B:332:PRO:HD3	1.88	0.55
1:B:348:ARG:NH1	1:B:348:ARG:HG3	2.21	0.55
1:F:242:LYS:HB2	1:F:245:ASP:OD2	2.07	0.55
1:A:51:SER:HB2	1:A:73:TYR:CZ	2.42	0.55
1:A:51:SER:OG	1:A:54:PHE:HB2	2.07	0.55
1:B:36:LEU:HB2	1:D:137:GLU:OE1	2.07	0.55
1:E:223:ASP:OD1	1:E:228:ARG:NH2	2.40	0.55
1:E:368:THR:HG22	1:F:36:LEU:HD11	1.88	0.55
1:A:325:MET:C	1:B:327:ASP:HB2	2.32	0.54
1:B:65:MET:HE2	1:B:192:PHE:CD2	2.43	0.54
1:C:262:LEU:O	1:C:265:ASN:HB3	2.07	0.54
1:E:185:ARG:HD2	3:E:438:HOH:O	2.06	0.54
1:C:118:ASP:O	1:C:121:VAL:HG22	2.08	0.54
1:E:368:THR:O	1:E:369:GLN:CB	2.55	0.54
1:F:320:GLN:NE2	1:F:322:VAL:N	2.47	0.54
1:B:325:MET:HE1	1:C:294:VAL:CG1	2.37	0.54
1:B:353:ASN:HD22	1:B:355:SER:H	1.50	0.54
1:F:314:VAL:HG13	3:F:501:HOH:O	2.06	0.54
1:E:353:ASN:HD22	1:E:355:SER:H	1.55	0.54
1:A:291:ILE:HD11	1:C:326:MET:HE3	1.90	0.54
1:D:118:ASP:O	1:D:121:VAL:HG22	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:210:PHE:CZ	1:A:270:ILE:HD11	2.43	0.54
1:D:92:LYS:O	1:D:96:LEU:HG	2.08	0.54
1:C:318:LEU:CD1	1:C:320:GLN:HG3	2.38	0.53
1:D:115:LYS:HD3	3:D:516:HOH:O	2.07	0.53
1:F:211:LEU:HD23	1:F:296:ALA:HB2	1.91	0.53
1:B:325:MET:HE1	1:C:294:VAL:HG12	1.90	0.53
1:D:356:SER:O	1:D:360:ARG:HG3	2.08	0.53
1:E:333:ALA:O	1:E:337:ILE:HG13	2.09	0.53
1:F:320:GLN:HE22	1:F:322:VAL:CG2	2.19	0.53
2:A:401:BH3:OAG	2:A:401:BH3:H9	2.09	0.53
1:C:315:LEU:O	1:C:315:LEU:HD12	2.08	0.53
1:F:86:MET:C	1:F:88:ILE:H	2.17	0.53
1:F:233:GLN:HA	1:F:236:TRP:HE1	1.68	0.53
1:A:211:LEU:HD12	2:A:401:BH3:H29	1.90	0.53
1:F:252:ILE:HD12	1:F:307:GLN:HE21	1.72	0.53
1:D:345:ILE:O	1:D:349:ILE:HG13	2.09	0.53
1:F:200:ASP:OD2	1:F:203:ARG:HB2	2.08	0.53
1:F:252:ILE:HD13	1:F:252:ILE:O	2.08	0.53
1:C:297:ILE:HD12	1:C:338:ILE:HG12	1.90	0.53
1:D:348:ARG:O	1:D:350:PRO:HD3	2.09	0.52
1:E:233:GLN:HA	1:E:236:TRP:CE2	2.44	0.52
1:F:233:GLN:HA	1:F:236:TRP:CD1	2.44	0.52
1:F:343:GLU:OE1	1:F:367:ARG:NH2	2.41	0.52
1:B:256:VAL:HG21	1:B:307:GLN:HG2	1.90	0.52
1:F:190:SER:O	1:F:191:GLU:HB2	2.08	0.52
1:C:255:ALA:HB1	1:C:310:PHE:CZ	2.44	0.52
1:D:51:SER:HB2	1:D:73:TYR:CZ	2.45	0.52
1:A:318:LEU:HG	1:A:319:GLY:N	2.25	0.52
1:F:242:LYS:HE2	3:F:451:HOH:O	2.10	0.52
1:F:317:ARG:HG3	1:F:318:LEU:N	2.25	0.52
1:F:321:ALA:O	1:F:325:MET:HB2	2.09	0.51
1:B:125:PHE:N	1:B:126:PRO:CD	2.73	0.51
1:C:333:ALA:O	1:C:337:ILE:HG12	2.09	0.51
1:E:304:TYR:CE2	1:E:349:ILE:HD13	2.45	0.51
1:D:224:GLN:C	1:D:226:GLY:H	2.18	0.51
1:A:353:ASN:HD22	1:A:355:SER:H	1.57	0.51
1:E:268:HIS:HA	1:E:362:ILE:CD1	2.38	0.51
1:C:47:LEU:O	1:C:51:SER:HB3	2.11	0.51
1:F:50:THR:HB	1:F:77:ARG:HD2	1.93	0.51
1:E:259:LEU:HD21	1:E:303:CYS:O	2.10	0.51
1:C:100:PHE:HA	1:C:103:PHE:CD2	2.45	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:369:GLN:HG3	3:F:531:HOH:O	2.12	0.50
1:B:213:LYS:O	1:B:217:ILE:HG12	2.11	0.50
1:F:88:ILE:HD11	1:F:111:PHE:HZ	1.76	0.50
1:F:314:VAL:HG12	1:F:315:LEU:N	2.27	0.50
1:F:321:ALA:C	1:F:323:THR:H	2.20	0.50
1:F:355:SER:HA	1:F:358:LYS:HE2	1.92	0.50
1:A:150:MET:O	1:A:154:MET:HG3	2.11	0.50
1:C:224:GLN:HE21	1:C:224:GLN:CA	2.25	0.50
1:D:255:ALA:HB1	1:D:310:PHE:CZ	2.47	0.50
1:E:111:PHE:HB3	1:E:122:LEU:HB3	1.93	0.50
1:E:325:MET:C	1:F:327:ASP:HB2	2.37	0.50
1:A:326:MET:HA	1:B:327:ASP:OD2	2.12	0.50
1:D:239:TYR:O	1:D:254:LEU:HD13	2.12	0.50
1:D:125:PHE:N	1:D:126:PRO:CD	2.75	0.49
1:D:291:ILE:HD11	1:F:326:MET:HE3	1.94	0.49
1:F:72:PHE:CD1	1:F:143:ILE:HG23	2.46	0.49
1:B:262:LEU:O	1:B:265:ASN:HB3	2.11	0.49
1:D:341:TYR:O	1:D:345:ILE:HD13	2.12	0.49
1:F:84:ASP:HA	1:F:228:ARG:HH21	1.77	0.49
1:F:320:GLN:O	1:F:324:LEU:HG	2.12	0.49
1:C:125:PHE:N	1:C:126:PRO:CD	2.76	0.49
1:D:338:ILE:O	1:D:342:MET:HG2	2.13	0.49
1:A:316:ILE:HG12	1:A:316:ILE:O	2.11	0.49
1:F:84:ASP:HA	1:F:228:ARG:NH2	2.28	0.49
1:E:297:ILE:CD1	1:E:338:ILE:HG12	2.43	0.49
1:F:110:ARG:HD3	3:F:426:HOH:O	2.12	0.49
1:E:228:ARG:HG3	1:E:228:ARG:HH11	1.78	0.49
1:D:161:HIS:CD2	1:D:229:GLU:HB3	2.46	0.49
1:F:252:ILE:HG23	1:F:253:ASP:N	2.27	0.49
1:E:297:ILE:HD12	1:E:338:ILE:HG12	1.93	0.49
1:B:252:ILE:HG23	1:B:253:ASP:N	2.28	0.48
1:F:317:ARG:HG3	1:F:318:LEU:H	1.77	0.48
1:A:125:PHE:N	1:A:126:PRO:CD	2.76	0.48
1:B:190:SER:O	1:B:191:GLU:HB2	2.12	0.48
1:C:71:ILE:HD11	1:C:131:GLU:HB3	1.94	0.48
1:D:52:ARG:HD3	1:D:53:SER:H	1.77	0.48
1:E:51:SER:OG	1:E:54:PHE:HB2	2.13	0.48
1:A:211:LEU:HD12	2:A:401:BH3:H30	1.94	0.48
1:A:316:ILE:HD13	1:A:316:ILE:N	2.17	0.48
1:B:150:MET:O	1:B:154:MET:HG3	2.12	0.48
1:D:327:ASP:HB2	1:F:326:MET:HA	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:342:MET:HE2	1:F:342:MET:HA	1.95	0.48
2:A:401:BH3:OAG	2:A:401:BH3:CAI	2.61	0.48
1:B:157:PHE:CE2	1:B:160:LYS:HE3	2.49	0.48
1:B:217:ILE:HD11	1:B:262:LEU:HB3	1.94	0.48
1:E:252:ILE:HG23	1:E:253:ASP:N	2.29	0.48
1:D:320:GLN:H	1:D:323:THR:HB	1.78	0.48
1:E:110:ARG:NH1	1:E:112:MET:HE2	2.27	0.48
1:A:71:ILE:HD12	1:A:135:LEU:HD21	1.95	0.48
2:A:401:BH3:H38	3:A:761:HOH:O	2.12	0.48
1:D:228:ARG:HB2	1:D:228:ARG:CZ	2.44	0.48
1:E:64:GLU:HG3	1:E:139:TYR:OH	2.14	0.48
1:F:369:GLN:HG3	1:F:369:GLN:O	2.14	0.48
1:B:185:ARG:HD2	3:B:673:HOH:O	2.14	0.48
1:F:358:LYS:HD2	3:F:477:HOH:O	2.13	0.48
1:A:341:TYR:O	1:A:344:GLU:HB2	2.13	0.48
1:B:287:ASN:O	1:B:291:ILE:HD13	2.14	0.48
1:F:252:ILE:CG2	1:F:253:ASP:N	2.76	0.48
1:E:126:PRO:HA	1:E:129:SER:OG	2.14	0.48
1:C:224:GLN:HA	1:C:224:GLN:NE2	2.29	0.47
1:C:233:GLN:HA	1:C:236:TRP:NE1	2.29	0.47
1:C:150:MET:O	1:C:154:MET:HG3	2.14	0.47
1:E:320:GLN:HB3	1:E:340:GLN:NE2	2.29	0.47
1:B:118:ASP:O	1:B:121:VAL:HG22	2.14	0.47
1:D:190:SER:O	1:D:191:GLU:HB2	2.14	0.47
1:B:325:MET:O	1:C:327:ASP:HB2	2.15	0.47
1:B:368:THR:O	1:B:369:GLN:HB3	2.14	0.47
1:D:232:PRO:HG2	1:D:235:VAL:HB	1.97	0.47
1:D:246:PHE:CD1	1:D:255:ALA:HA	2.49	0.47
1:E:224:GLN:C	1:E:226:GLY:H	2.23	0.47
1:F:125:PHE:N	1:F:126:PRO:CD	2.78	0.47
1:F:297:ILE:HD12	1:F:338:ILE:HG12	1.95	0.47
1:F:343:GLU:OE1	1:F:367:ARG:NE	2.46	0.47
1:B:368:THR:O	1:B:368:THR:HG22	2.14	0.47
1:B:233:GLN:HA	1:B:236:TRP:CD1	2.50	0.47
1:E:124:ASP:C	1:E:126:PRO:HD2	2.39	0.46
1:E:295:MET:HE2	1:E:295:MET:HA	1.96	0.46
1:F:252:ILE:HD13	1:F:256:VAL:HG23	1.97	0.46
1:B:94:VAL:HB	1:B:95:PRO:HD3	1.97	0.46
1:E:346:TYR:O	1:E:349:ILE:HG12	2.15	0.46
1:C:196:LEU:HD21	1:C:279:ARG:CZ	2.45	0.46
1:F:323:THR:HG22	1:F:337:ILE:HG12	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:228:ARG:HG2	1:B:228:ARG:HH11	1.81	0.46
1:D:218:ARG:HH21	1:D:218:ARG:HG3	1.79	0.46
1:D:320:GLN:N	1:D:323:THR:HB	2.31	0.46
1:A:190:SER:O	1:A:191:GLU:HB2	2.14	0.46
1:E:125:PHE:N	1:E:126:PRO:CD	2.77	0.46
1:A:119:ARG:O	1:A:123:GLU:HG3	2.16	0.46
1:D:253:ASP:O	1:D:257:GLN:HG3	2.15	0.46
1:A:318:LEU:HD12	1:A:319:GLY:H	1.80	0.46
1:B:100:PHE:HA	1:B:103:PHE:CD2	2.51	0.46
1:E:153:GLY:HA3	1:E:174:TYR:CG	2.50	0.46
1:E:268:HIS:CE1	1:E:362:ILE:HD12	2.51	0.46
1:A:318:LEU:CG	1:A:319:GLY:N	2.79	0.46
1:B:220:TYR:HB2	1:B:231:TRP:CZ2	2.51	0.46
1:B:343:GLU:OE1	1:C:48:ASN:HB3	2.15	0.46
1:D:316:ILE:N	1:D:316:ILE:HD12	2.31	0.46
1:A:76:LEU:HB3	2:A:401:BH3:H18	1.98	0.45
1:B:36:LEU:N	1:B:36:LEU:CD1	2.78	0.45
1:B:368:THR:O	1:B:369:GLN:CB	2.64	0.45
1:C:242:LYS:NZ	1:C:242:LYS:HB3	2.32	0.45
1:F:323:THR:C	1:F:325:MET:N	2.74	0.45
1:E:90:VAL:O	1:E:94:VAL:HG23	2.16	0.45
1:E:317:ARG:H	1:E:317:ARG:HG2	1.53	0.45
1:F:138:LYS:HE3	1:F:139:TYR:CZ	2.51	0.45
1:F:224:GLN:C	1:F:226:GLY:H	2.24	0.45
1:B:152:ILE:HD11	3:B:507:HOH:O	2.15	0.45
1:E:358:LYS:H	1:E:358:LYS:HD2	1.80	0.45
1:A:326:MET:CA	1:B:327:ASP:HB2	2.45	0.45
1:A:344:GLU:O	1:A:348:ARG:HG3	2.17	0.45
1:C:224:GLN:C	1:C:226:GLY:H	2.24	0.45
1:A:327:ASP:HB3	1:C:325:MET:O	2.17	0.45
1:E:67:ASN:O	1:E:71:ILE:HG12	2.17	0.45
1:A:228:ARG:NH1	3:A:731:HOH:O	2.49	0.45
1:A:238:ARG:NH2	1:A:261:GLU:OE2	2.44	0.45
1:B:233:GLN:HE21	1:B:233:GLN:HB3	1.57	0.45
1:C:338:ILE:O	1:C:342:MET:HG2	2.17	0.45
1:C:353:ASN:HD21	1:C:355:SER:HB2	1.81	0.45
1:E:268:HIS:CA	1:E:362:ILE:HD11	2.41	0.45
1:D:333:ALA:O	1:D:337:ILE:HG13	2.17	0.45
1:F:153:GLY:HA3	1:F:174:TYR:CD1	2.52	0.45
1:F:320:GLN:HE22	1:F:322:VAL:N	2.06	0.44
1:B:353:ASN:HD22	1:B:355:SER:N	2.14	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:297:ILE:HD13	1:C:338:ILE:HG12	1.95	0.44
1:D:346:TYR:HA	1:D:349:ILE:CD1	2.47	0.44
1:F:42:THR:HG22	1:F:46:TYR:CE2	2.53	0.44
1:D:362:ILE:HG23	1:D:363:ILE:N	2.32	0.44
1:E:150:MET:O	1:E:154:MET:HG3	2.17	0.44
1:F:220:TYR:CZ	1:F:246:PHE:HB2	2.52	0.44
1:B:344:GLU:O	1:B:348:ARG:HG3	2.18	0.44
1:E:213:LYS:O	1:E:217:ILE:HG13	2.17	0.44
1:B:338:ILE:O	1:B:342:MET:HG2	2.16	0.44
1:D:153:GLY:HA3	1:D:174:TYR:CG	2.53	0.44
1:F:252:ILE:HG23	1:F:253:ASP:OD2	2.17	0.44
1:A:105:TYR:O	1:A:107:PRO:HD3	2.18	0.44
1:C:190:SER:O	1:C:191:GLU:HB2	2.18	0.44
1:B:248:LEU:HB2	1:B:251:ASN:HD22	1.83	0.44
1:C:51:SER:HB2	1:C:73:TYR:CZ	2.52	0.43
1:C:320:GLN:O	1:C:321:ALA:CB	2.65	0.43
1:D:262:LEU:O	1:D:265:ASN:HB3	2.18	0.43
1:D:320:GLN:OE1	1:D:321:ALA:N	2.51	0.43
1:D:52:ARG:HH21	1:D:52:ARG:HG3	1.83	0.43
1:D:105:TYR:O	1:D:107:PRO:HD3	2.18	0.43
1:F:323:THR:HA	1:F:326:MET:HG2	1.99	0.43
1:B:316:ILE:N	1:B:316:ILE:HD12	2.34	0.43
1:B:321:ALA:O	1:B:325:MET:HG2	2.19	0.43
1:F:323:THR:O	1:F:326:MET:O	2.37	0.43
1:F:220:TYR:OH	1:F:246:PHE:HB2	2.19	0.43
1:F:252:ILE:HD13	1:F:252:ILE:C	2.44	0.43
1:A:41:LYS:HE3	3:C:679:HOH:O	2.18	0.43
1:A:90:VAL:O	1:A:94:VAL:HG23	2.19	0.43
1:D:326:MET:HE1	3:E:411:HOH:O	2.16	0.43
1:A:52:ARG:CZ	2:A:401:BH3:OAC	2.67	0.43
1:A:107:PRO:O	1:A:126:PRO:HB3	2.19	0.43
1:C:38:SER:HA	1:C:41:LYS:HG3	2.00	0.43
1:D:233:GLN:HE21	1:D:233:GLN:HB3	1.56	0.43
1:E:190:SER:O	1:E:191:GLU:HB2	2.18	0.43
1:C:46:TYR:CE1	1:C:120:GLN:HG3	2.53	0.43
1:E:170:LYS:O	1:E:173:HIS:HB3	2.19	0.43
1:A:52:ARG:NH1	2:A:401:BH3:OAC	2.51	0.42
1:E:287:ASN:O	1:E:291:ILE:HD13	2.19	0.42
1:F:181:ILE:HD12	1:F:181:ILE:N	2.34	0.42
1:A:217:ILE:HD11	1:A:262:LEU:HB3	2.01	0.42
1:A:153:GLY:HA3	1:A:174:TYR:CG	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:297:ILE:CD1	1:F:338:ILE:HG12	2.49	0.42
1:A:327:ASP:HB3	1:C:326:MET:HA	2.02	0.42
1:E:220:TYR:HB2	1:E:231:TRP:CZ2	2.55	0.42
1:F:253:ASP:O	1:F:257:GLN:HG3	2.19	0.42
1:A:353:ASN:HA	1:A:354:PRO:HD3	1.82	0.42
1:F:252:ILE:HD11	1:F:307:GLN:HG2	2.00	0.42
1:B:51:SER:OG	1:B:54:PHE:HB2	2.19	0.42
1:D:209:LEU:HD23	1:D:269:HIS:CE1	2.54	0.42
1:E:193:GLU:OE2	1:E:282:ASN:HB3	2.20	0.42
1:E:317:ARG:HH21	1:E:348:ARG:NH1	2.17	0.42
1:C:291:ILE:HB	1:C:292:PRO:HD3	2.01	0.42
1:C:318:LEU:HD12	1:C:319:GLY:O	2.19	0.42
1:E:65:MET:HE2	1:E:192:PHE:CD2	2.55	0.42
1:F:150:MET:HE3	1:F:150:MET:HB2	1.78	0.42
1:D:250:GLU:OE2	1:D:250:GLU:N	2.49	0.42
1:C:73:TYR:CD1	1:C:73:TYR:C	2.97	0.42
1:C:343:GLU:CD	1:C:367:ARG:HH21	2.27	0.42
1:F:86:MET:C	1:F:88:ILE:N	2.77	0.42
1:B:327:ASP:OD2	1:B:329:THR:HG23	2.19	0.42
1:F:110:ARG:HG2	1:F:110:ARG:HH21	1.84	0.42
1:D:137:GLU:HB2	3:D:451:HOH:O	2.19	0.41
1:F:155:ALA:HA	1:F:158:LEU:HG	2.02	0.41
1:F:320:GLN:HE22	1:F:322:VAL:CB	2.32	0.41
1:B:255:ALA:HB1	1:B:310:PHE:CZ	2.54	0.41
1:C:326:MET:HE2	1:C:333:ALA:HB2	2.02	0.41
1:D:322:VAL:HG13	1:E:291:ILE:HG21	2.01	0.41
1:F:242:LYS:HB2	1:F:245:ASP:CG	2.45	0.41
1:F:362:ILE:O	1:F:366:ILE:HG13	2.21	0.41
1:B:353:ASN:HD22	1:B:356:SER:H	1.67	0.41
1:D:263:ILE:CD1	1:D:304:TYR:HA	2.50	0.41
1:D:324:LEU:HD12	1:D:324:LEU:C	2.46	0.41
1:F:142:VAL:O	1:F:146:ILE:HD13	2.19	0.41
1:F:225:GLN:HE21	1:F:225:GLN:HB2	1.49	0.41
1:E:150:MET:HG3	1:E:174:TYR:O	2.20	0.41
1:F:344:GLU:O	1:F:348:ARG:HG3	2.20	0.41
1:B:253:ASP:O	1:B:257:GLN:HG3	2.19	0.41
1:C:319:GLY:O	1:C:324:LEU:HD21	2.20	0.41
1:D:229:GLU:HG3	1:D:243:LEU:HD22	2.03	0.41
1:D:90:VAL:HA	1:D:93:LYS:HB3	2.03	0.41
1:A:233:GLN:HA	1:A:236:TRP:CD1	2.55	0.41
1:C:220:TYR:HB2	1:C:231:TRP:CZ2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:79:LEU:HB2	1:D:100:PHE:CE2	2.55	0.41
1:B:327:ASP:OD2	1:B:329:THR:CG2	2.69	0.41
1:A:183:LEU:CD1	2:A:401:BH3:H35	2.51	0.40
1:B:348:ARG:O	1:B:350:PRO:HD3	2.21	0.40
1:C:326:MET:HG3	1:C:337:ILE:HD11	2.03	0.40
1:F:242:LYS:O	1:F:245:ASP:HB2	2.21	0.40
1:A:318:LEU:CG	1:A:319:GLY:H	2.34	0.40
1:B:71:ILE:HD11	1:B:131:GLU:HB3	2.02	0.40
1:C:110:ARG:HH21	1:C:110:ARG:HG3	1.85	0.40
1:C:142:VAL:O	1:C:146:ILE:HG12	2.21	0.40
1:C:331:MET:HB3	1:C:332:PRO:HD3	2.03	0.40
1:D:301:ALA:CB	1:D:345:ILE:HD11	2.51	0.40
1:E:349:ILE:HD12	1:E:359:THR:HG21	2.03	0.40
1:A:220:TYR:OH	1:A:246:PHE:HB2	2.20	0.40
1:D:52:ARG:HD2	1:D:53:SER:H	1.85	0.40
1:D:150:MET:O	1:D:154:MET:HG3	2.21	0.40
1:C:242:LYS:NZ	1:C:242:LYS:CB	2.84	0.40
1:D:44:TYR:OH	1:D:66:ARG:HG2	2.22	0.40
1:E:304:TYR:CZ	1:E:349:ILE:HD13	2.57	0.40
1:F:193:GLU:HB3	1:F:197:VAL:HG21	2.02	0.40
1:A:248:LEU:HA	1:A:249:PRO:HD3	1.88	0.40
1:E:93:LYS:HE3	1:E:93:LYS:HB2	1.93	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	332/360 (92%)	320 (96%)	9 (3%)	3 (1%)	14	11
1	B	332/360 (92%)	316 (95%)	16 (5%)	0	100	100
1	C	332/360 (92%)	316 (95%)	15 (4%)	1 (0%)	36	38

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	332/360 (92%)	312 (94%)	17 (5%)	3 (1%)	14	11
1	E	332/360 (92%)	317 (96%)	12 (4%)	3 (1%)	14	11
1	F	332/360 (92%)	307 (92%)	21 (6%)	4 (1%)	10	6
All	All	1992/2160 (92%)	1888 (95%)	90 (4%)	14 (1%)	18	16

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	318	LEU
1	A	322	VAL
1	D	317	ARG
1	F	317	ARG
1	A	321	ALA
1	C	321	ALA
1	E	318	LEU
1	F	322	VAL
1	F	313	ALA
1	F	323	THR
1	E	107	PRO
1	E	317	ARG
1	D	249	PRO
1	D	226	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	297/320 (93%)	290 (98%)	7 (2%)	43	51
1	B	297/320 (93%)	290 (98%)	7 (2%)	43	51
1	C	297/320 (93%)	291 (98%)	6 (2%)	48	58
1	D	297/320 (93%)	290 (98%)	7 (2%)	43	51
1	E	297/320 (93%)	285 (96%)	12 (4%)	28	32
1	F	297/320 (93%)	282 (95%)	15 (5%)	21	21

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	1782/1920 (93%)	1728 (97%)	54 (3%)	36	43

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	167	GLU
1	A	233	GLN
1	A	241	LYS
1	A	252	ILE
1	A	316	ILE
1	A	326	MET
1	A	369	GLN
1	B	36	LEU
1	B	137	GLU
1	B	233	GLN
1	B	250	GLU
1	B	317	ARG
1	B	351	ASP
1	B	369	GLN
1	C	41	LYS
1	C	170	LYS
1	C	203	ARG
1	C	224	GLN
1	C	234	GLU
1	C	326	MET
1	D	52	ARG
1	D	91	GLU
1	D	221	LEU
1	D	228	ARG
1	D	233	GLN
1	D	307	GLN
1	D	367	ARG
1	E	52	ARG
1	E	91	GLU
1	E	107	PRO
1	E	110	ARG
1	E	113	GLU
1	E	233	GLN
1	E	242	LYS
1	E	317	ARG
1	E	320	GLN
1	E	322	VAL

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Mol	Chain	Res	Type
1	E	324	LEU
1	E	369	GLN
1	F	77	ARG
1	F	199	GLU
1	F	224	GLN
1	F	225	GLN
1	F	241	LYS
1	F	242	LYS
1	F	252	ILE
1	F	254	LEU
1	F	308	GLN
1	F	320	GLN
1	F	323	THR
1	F	325	MET
1	F	326	MET
1	F	340	GLN
1	F	351	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (54) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	48	ASN
1	A	59	GLN
1	A	161	HIS
1	A	224	GLN
1	A	225	GLN
1	A	233	GLN
1	A	320	GLN
1	A	340	GLN
1	B	98	HIS
1	B	120	GLN
1	B	224	GLN
1	B	225	GLN
1	B	233	GLN
1	B	251	ASN
1	B	260	ASN
1	B	305	ASN
1	B	308	GLN
1	B	340	GLN
1	B	353	ASN
1	B	369	GLN
1	C	98	HIS

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Mol	Chain	Res	Type
1	C	140	GLN
1	C	225	GLN
1	C	233	GLN
1	C	257	GLN
1	C	260	ASN
1	C	293	GLN
1	C	305	ASN
1	C	347	HIS
1	C	353	ASN
1	D	67	ASN
1	D	161	HIS
1	D	224	GLN
1	D	233	GLN
1	D	251	ASN
1	D	308	GLN
1	D	353	ASN
1	E	48	ASN
1	E	59	GLN
1	E	120	GLN
1	E	225	GLN
1	E	233	GLN
1	E	308	GLN
1	F	48	ASN
1	F	59	GLN
1	F	224	GLN
1	F	225	GLN
1	F	257	GLN
1	F	293	GLN
1	F	307	GLN
1	F	308	GLN
1	F	320	GLN
1	F	347	HIS
1	F	353	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	BH3	A	401	-	29,30,30	2.14	5 (17%)	36,39,39	0.98	3 (8%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	BH3	A	401	-	-	6/29/31/31	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	BH3	PBC-OAE	7.10	1.61	1.49
2	A	401	BH3	PBC-OAD	4.14	1.61	1.54
2	A	401	BH3	PBD-OAF	4.13	1.61	1.54
2	A	401	BH3	PBD-OAG	4.13	1.61	1.54
2	A	401	BH3	PBC-OAB	-4.04	1.48	1.54

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	BH3	OAD-PBC-OAE	-3.01	105.91	113.45
2	A	401	BH3	OAB-PBC-CAZ	2.15	113.59	106.88
2	A	401	BH3	OAE-PBC-CAZ	-2.07	106.42	112.29

There are no chirality outliers.

All (6) torsion outliers are listed below:

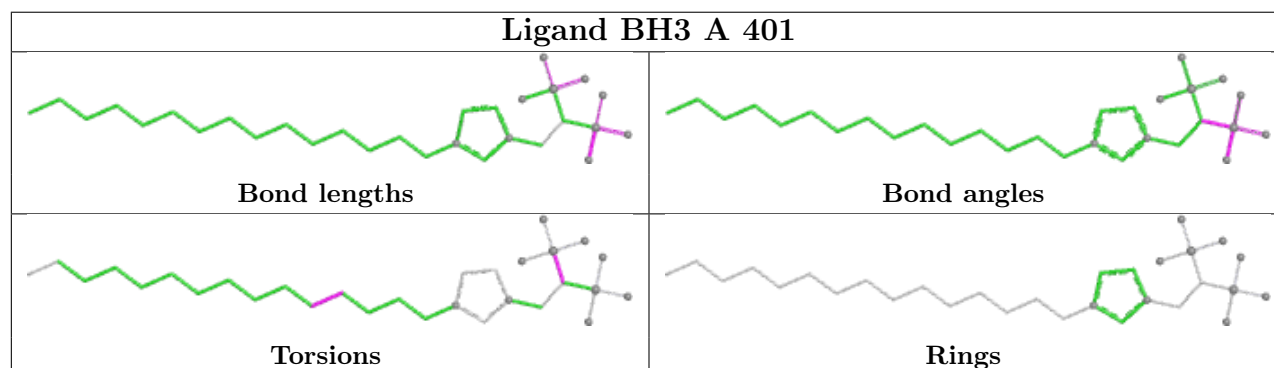
Mol	Chain	Res	Type	Atoms
2	A	401	BH3	CAY-CAZ-PBD-OAF
2	A	401	BH3	PBC-CAZ-PBD-OAF
2	A	401	BH3	PBC-CAZ-PBD-OAC
2	A	401	BH3	CAY-CAZ-PBD-OAC
2	A	401	BH3	CAY-CAZ-PBD-OAG
2	A	401	BH3	CAS-CAT-CAU-CAV

There are no ring outliers.

1 monomer is involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	BH3	13	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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	334/360 (92%)	-0.39	12 (3%) 46 46	18, 30, 62, 112	0
1	B	334/360 (92%)	-0.23	6 (1%) 67 69	19, 32, 72, 102	0
1	C	334/360 (92%)	-0.33	1 (0%) 90 91	19, 33, 68, 89	0
1	D	334/360 (92%)	-0.11	5 (1%) 72 73	21, 40, 74, 112	0
1	E	334/360 (92%)	-0.04	6 (1%) 67 69	28, 44, 74, 108	0
1	F	334/360 (92%)	0.22	11 (3%) 49 49	23, 51, 87, 126	0
All	All	2004/2160 (92%)	-0.15	41 (2%) 65 66	18, 38, 79, 126	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	318	LEU	4.2
1	A	322	VAL	4.1
1	A	315	LEU	3.9
1	F	369	GLN	3.6
1	E	318	LEU	3.2
1	D	36	LEU	3.1
1	F	323	THR	2.9
1	A	314	VAL	2.9
1	A	321	ALA	2.8
1	F	316	ILE	2.8
1	A	318	LEU	2.8
1	F	327	ASP	2.8
1	E	36	LEU	2.7
1	B	318	LEU	2.7
1	B	321	ALA	2.6
1	E	321	ALA	2.6
1	B	319	GLY	2.6
1	F	319	GLY	2.6
1	A	36	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	D	318	LEU	2.5
1	B	368	THR	2.4
1	A	324	LEU	2.4
1	B	54	PHE	2.4
1	D	325	MET	2.4
1	D	324	LEU	2.4
1	F	313	ALA	2.4
1	E	323	THR	2.4
1	A	313	ALA	2.3
1	F	315	LEU	2.3
1	A	316	ILE	2.3
1	F	231	TRP	2.2
1	E	322	VAL	2.2
1	A	325	MET	2.2
1	B	317	ARG	2.1
1	E	324	LEU	2.1
1	F	324	LEU	2.1
1	D	323	THR	2.1
1	A	159	ASP	2.1
1	C	325	MET	2.0
1	F	322	VAL	2.0
1	A	327	ASP	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

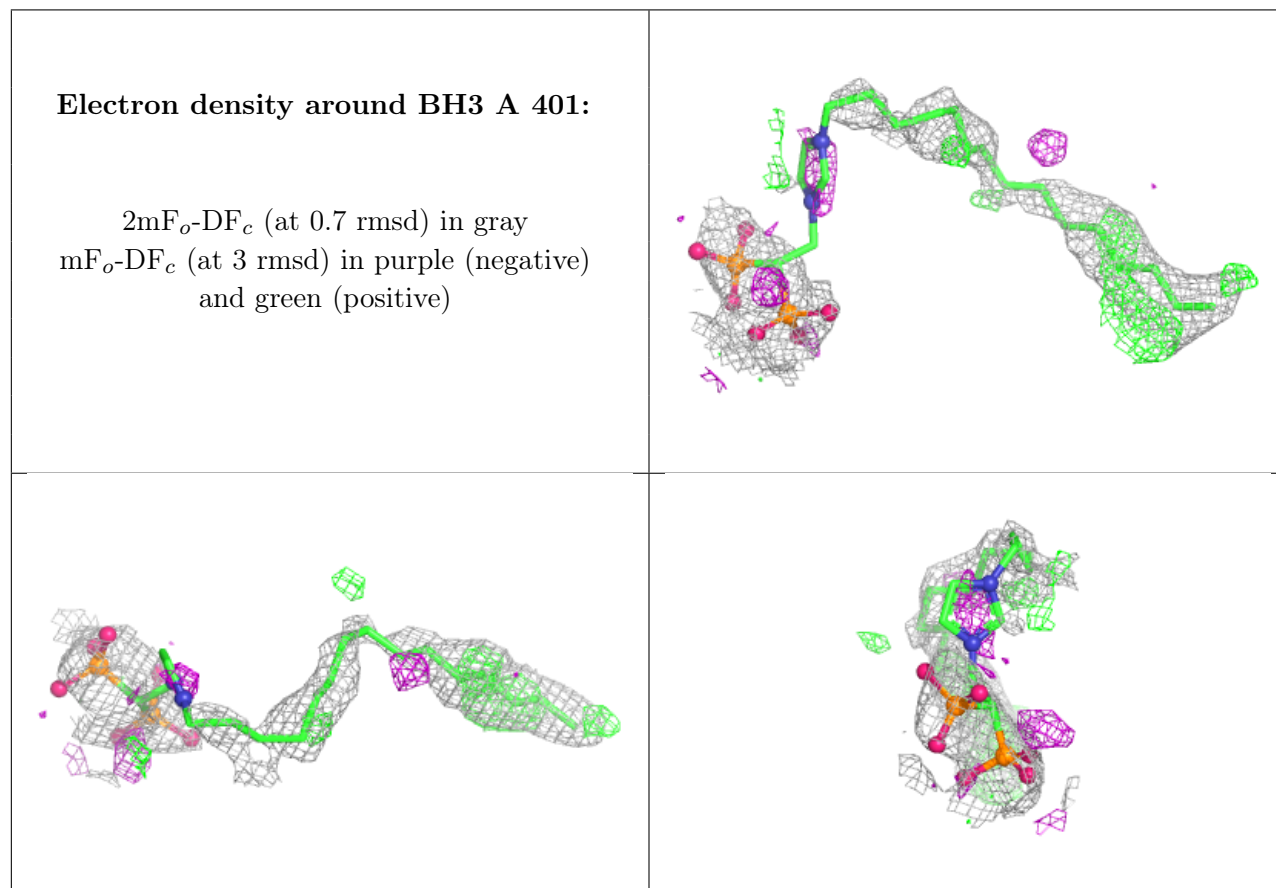
There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	BH3	A	401	30/30	0.65	0.21	48,109,125,125	0

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



6.5 Other polymers [i](#)

There are no such residues in this entry.