



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 29, 2026 – 06:39 AM UTC

PDB ID : 3WCF / pdb_00003wcf
Title : The complex structure of HsSQS wtih ligand,BPH1218
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 : 2013-05-27
Resolution : 2.22 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

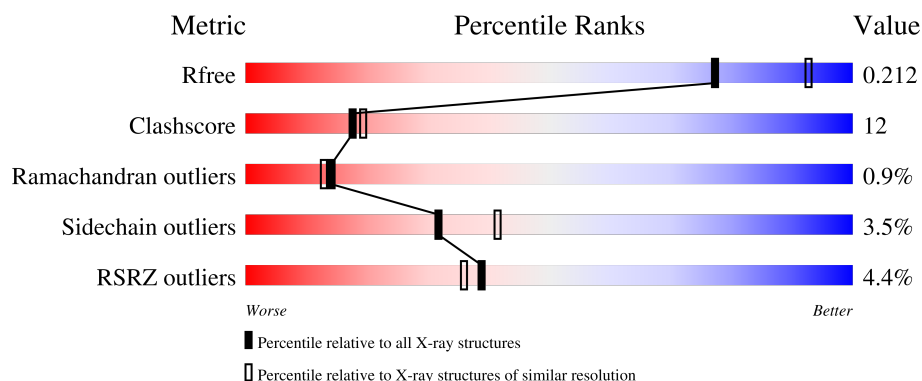
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.22 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	7682 (2.24-2.20)
Clashscore	190562	8402 (2.24-2.20)
Ramachandran outliers	187476	8303 (2.24-2.20)
Sidechain outliers	187428	8304 (2.24-2.20)
RSRZ outliers	180081	7683 (2.24-2.20)

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% 72% 18% • 7%
1	B	360	2% 71% 19% • 7%
1	C	360	2% 68% 23% • 7%
1	D	360	2% 68% 23% •• 7%
1	E	360	2% 68% 23% • 7%

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Mol	Chain	Length	Quality of chain
1	F	360	13% 57% 32% • 7%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 16821 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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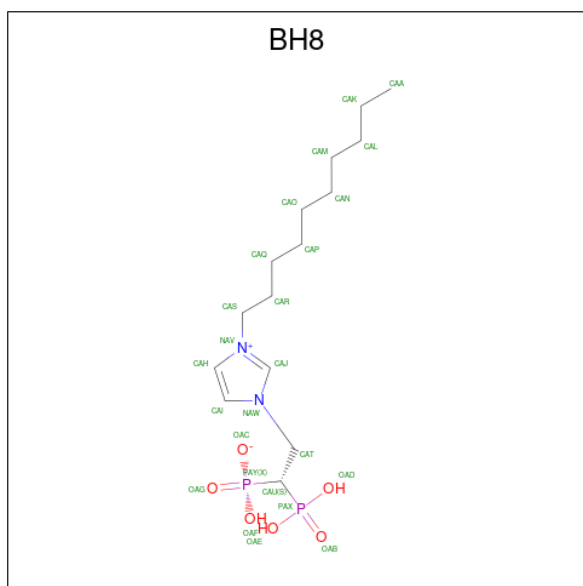
Chain	Residue	Modelled	Actual	Comment	Reference
E	16	HIS	-	expression tag	UNP P37268
E	17	HIS	-	expression tag	UNP P37268
E	18	HIS	-	expression tag	UNP P37268
E	19	HIS	-	expression tag	UNP P37268
E	20	HIS	-	expression tag	UNP P37268
E	21	SER	-	expression tag	UNP P37268
E	22	SER	-	expression tag	UNP P37268
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 [(1S)-2-(3-decyl-1H-imidazol-3-ium-1-yl)-1-phosphonoethyl]phosphate (CCD ID: BH8) (formula: C₁₅H₃₀N₂O₆P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			25	15	2	6	2		
2	B	1	Total	C	N	O	P	0	0
			25	15	2	6	2		
2	C	1	Total	C	N	O	P	0	0
			25	15	2	6	2		
2	D	1	Total	C	N	O	P	0	0
			25	15	2	6	2		
2	E	1	Total	C	N	O	P	0	0
			25	15	2	6	2		
2	F	1	Total	C	N	O	P	0	0
			25	15	2	6	2		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	151	Total	O	0	0
			151	151		
3	B	111	Total	O	0	0
			111	111		

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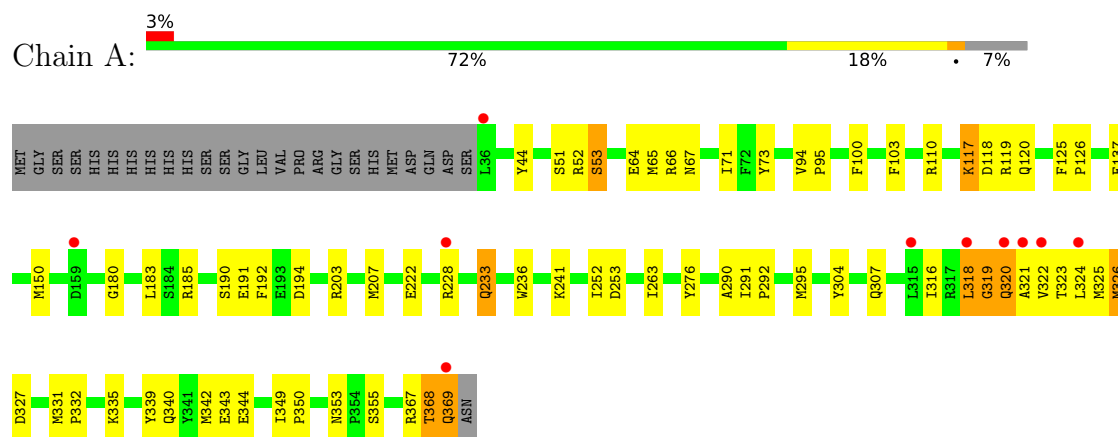
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	94	Total 94	O 94	0	0
3	D	68	Total 68	O 68	0	0
3	E	43	Total 43	O 43	0	0
3	F	28	Total 28	O 28	0	0

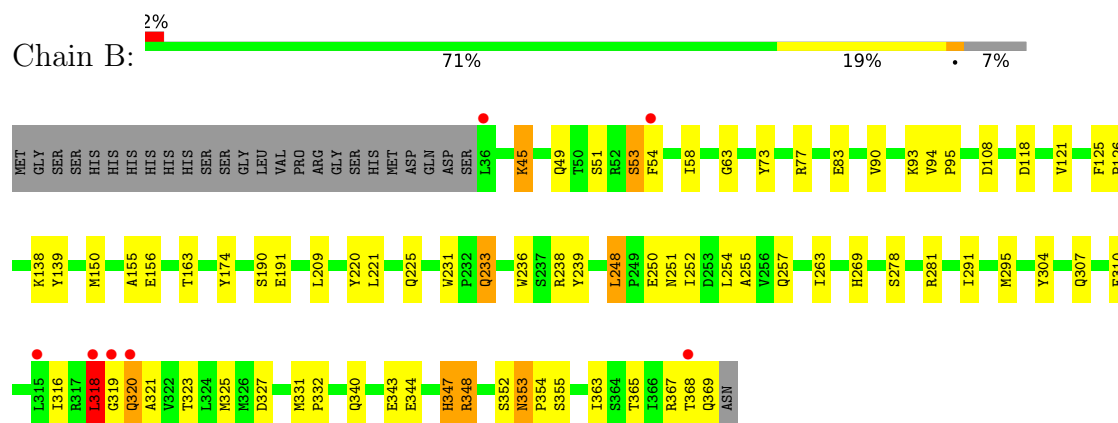
3 Residue-property plots [i](#)

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.

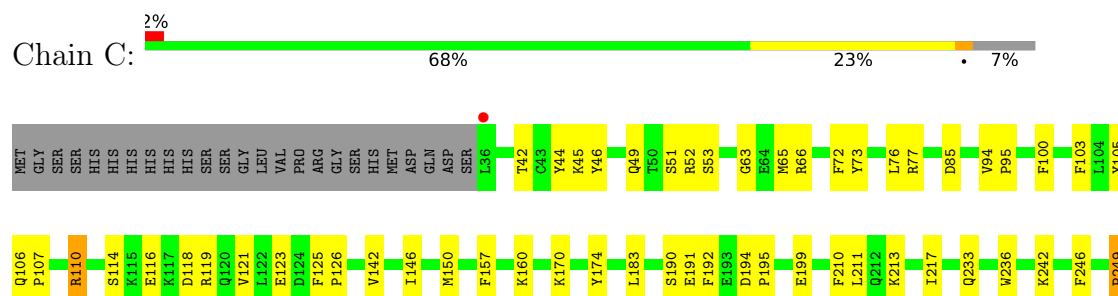
• Molecule 1: Squalene synthase

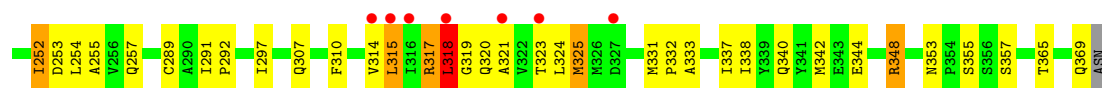


• Molecule 1: Squalene synthase

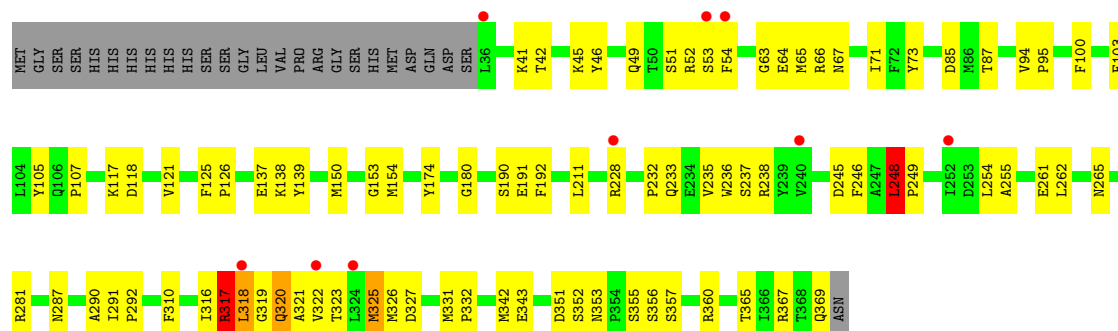


• Molecule 1: Squalene synthase

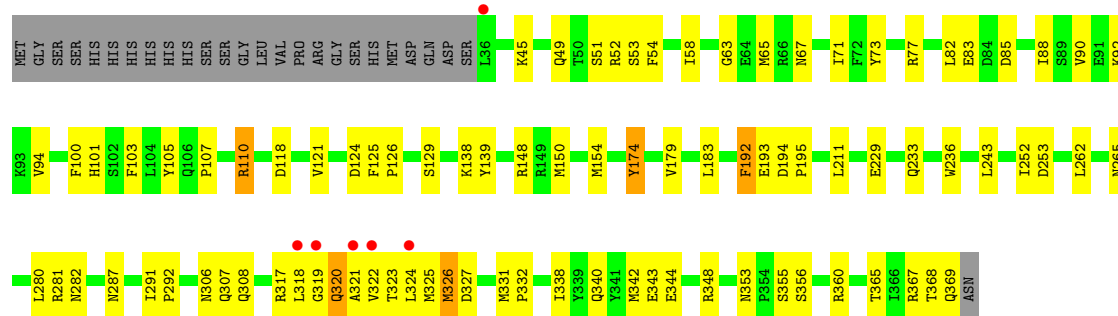




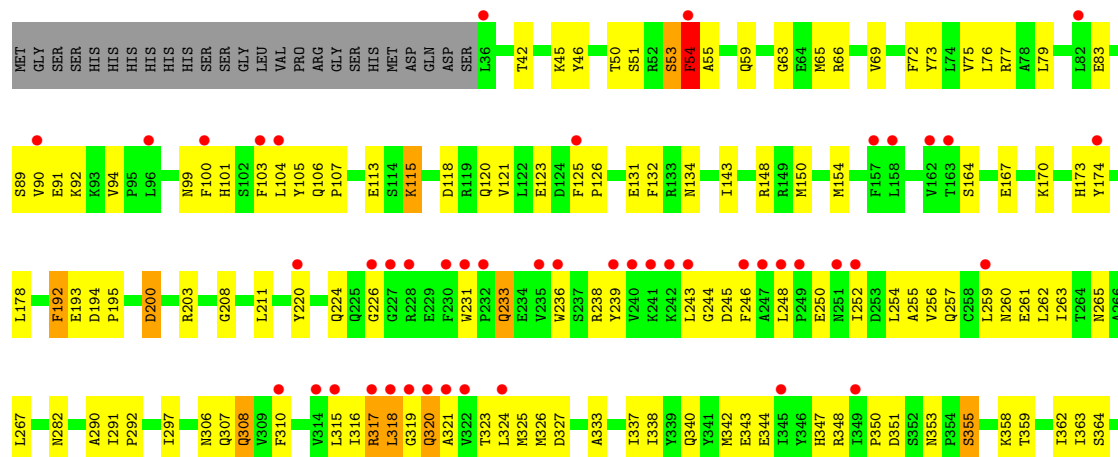
• Molecule 1: Squalene synthase



• 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	85.18Å 153.19Å 90.11Å 90.00° 92.37° 90.00°	Depositor
Resolution (Å)	25.00 – 2.22 25.00 – 2.22	Depositor EDS
% Data completeness (in resolution range)	(Not available) (25.00-2.22) 95.4 (25.00-2.22)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.43 (at 2.22Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.204 , 0.224 0.210 , 0.212	Depositor DCC
R_{free} test set	5442 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	32.7	Xtriage
Anisotropy	0.493	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 41.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.032 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	16821	wwPDB-VP
Average B, all atoms (Å ²)	48.0	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: BH8

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	0.40	0/2751	0.89	6/3724 (0.2%)
1	B	0.38	0/2751	0.90	8/3724 (0.2%)
1	C	0.38	0/2751	0.87	3/3724 (0.1%)
1	D	0.37	0/2751	0.89	6/3724 (0.2%)
1	E	0.35	0/2751	0.89	7/3724 (0.2%)
1	F	0.34	0/2751	0.89	8/3724 (0.2%)
All	All	0.37	0/16506	0.89	38/22344 (0.2%)

There are no bond length outliers.

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	317	ARG	N-CA-C	9.29	122.98	110.35
1	A	318	LEU	N-CA-C	8.40	122.50	111.75
1	A	53	SER	N-CA-C	8.03	121.88	112.72
1	F	54	PHE	N-CA-C	7.60	120.63	111.82
1	E	192	PHE	N-CA-C	-7.28	103.66	112.54
1	D	53	SER	N-CA-C	7.24	121.04	112.93
1	C	53	SER	N-CA-C	6.78	121.62	112.88
1	D	248	LEU	CA-C-N	6.69	126.95	119.32
1	D	248	LEU	C-N-CA	6.69	126.95	119.32
1	B	53	SER	N-CA-C	6.66	121.12	112.92
1	C	106	GLN	CA-C-N	6.65	128.15	119.84
1	C	106	GLN	C-N-CA	6.65	128.15	119.84
1	A	319	GLY	N-CA-C	6.27	118.84	111.63
1	F	192	PHE	N-CA-C	-6.21	105.53	113.23
1	D	254	LEU	N-CA-C	-6.13	104.67	111.36
1	F	53	SER	N-CA-C	6.09	120.78	112.68
1	B	248	LEU	CA-C-N	6.08	126.53	119.47
1	B	248	LEU	C-N-CA	6.08	126.53	119.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	368	THR	N-CA-C	5.79	118.53	111.82
1	B	353	ASN	CA-C-N	5.72	125.39	119.56
1	B	353	ASN	C-N-CA	5.72	125.39	119.56
1	A	344	GLU	N-CA-C	-5.70	105.06	111.28
1	A	290	ALA	N-CA-C	5.70	117.30	111.14
1	F	106	GLN	CA-C-N	5.67	126.93	119.84
1	F	106	GLN	C-N-CA	5.67	126.93	119.84
1	E	82	LEU	N-CA-C	-5.67	105.01	111.14
1	E	53	SER	N-CA-C	5.59	120.79	113.30
1	A	368	THR	N-CA-C	5.55	117.41	111.36
1	B	93	LYS	N-CA-C	5.51	116.97	111.07
1	F	317	ARG	N-CA-C	5.46	117.73	110.53
1	E	92	LYS	N-CA-C	-5.38	106.89	113.50
1	D	64	GLU	N-CA-C	5.30	117.75	111.33
1	B	45	LYS	N-CA-C	-5.29	105.43	111.14
1	F	290	ALA	N-CA-C	5.28	116.84	111.14
1	F	355	SER	N-CA-C	-5.26	107.24	113.97
1	B	348	ARG	N-CA-C	-5.25	106.83	114.12
1	D	290	ALA	N-CA-C	5.23	117.06	111.36
1	E	174	TYR	N-CA-C	5.00	116.73	111.28

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	65	0
1	B	2696	0	2676	60	0
1	C	2696	0	2676	68	0
1	D	2696	0	2676	61	0
1	E	2696	0	2676	71	0
1	F	2696	0	2676	97	0
2	A	25	0	30	6	0
2	B	25	0	30	0	0
2	C	25	0	30	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	25	0	30	4	0
2	E	25	0	30	7	0
2	F	25	0	30	4	0
3	A	151	0	0	5	0
3	B	111	0	0	3	0
3	C	94	0	0	0	0
3	D	68	0	0	2	0
3	E	43	0	0	0	0
3	F	28	0	0	1	0
All	All	16821	0	16236	404	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 12.

All (404) 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:F:317:ARG:HG3	1:F:318:LEU:HG	1.48	0.96
1:F:233:GLN:HA	1:F:236:TRP:NE1	1.89	0.88
1:E:326:MET:HE3	1:F:291:ILE:HD11	1.56	0.88
1:F:224:GLN:HE21	1:F:244:GLY:HA2	1.36	0.88
1:D:211:LEU:HD12	2:D:401:BH8:H21	1.57	0.87
1:C:320:GLN:O	1:C:324:LEU:HG	1.77	0.85
1:E:211:LEU:HD12	2:E:401:BH8:H21	1.58	0.85
1:E:291:ILE:HB	1:E:292:PRO:HD3	1.61	0.83
1:E:51:SER:HB2	1:E:73:TYR:CZ	2.14	0.83
1:F:343:GLU:OE1	1:F:367:ARG:NH2	2.15	0.79
1:E:322:VAL:HB	1:E:340:GLN:HE22	1.47	0.79
1:A:353:ASN:ND2	1:A:355:SER:H	1.81	0.79
1:E:100:PHE:HA	1:E:103:PHE:CD2	2.19	0.78
1:F:105:TYR:O	1:F:107:PRO:HD3	1.84	0.77
1:C:45:LYS:O	1:C:49:GLN:HG3	1.84	0.77
1:B:233:GLN:HA	1:B:236:TRP:NE1	2.00	0.77
1:B:343:GLU:OE1	1:B:367:ARG:NH2	2.17	0.77
1:C:325:MET:HE3	1:C:325:MET:HA	1.65	0.76
1:C:320:GLN:HB3	1:C:340:GLN:NE2	2.01	0.75
1:C:353:ASN:ND2	1:C:355:SER:H	1.84	0.75
1:E:105:TYR:O	1:E:107:PRO:HD3	1.86	0.74
1:E:365:THR:O	1:E:369:GLN:HG3	1.87	0.74
1:F:233:GLN:HA	1:F:236:TRP:CD1	2.23	0.73
1:F:317:ARG:HG3	1:F:318:LEU:N	2.04	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:318:LEU:HD12	1:B:319:GLY:H	1.54	0.72
1:A:319:GLY:HA3	1:A:323:THR:HG21	1.72	0.72
1:E:322:VAL:HB	1:E:340:GLN:NE2	2.04	0.72
1:F:150:MET:HG3	1:F:174:TYR:O	1.89	0.72
1:F:333:ALA:O	1:F:337:ILE:HG12	1.90	0.72
1:C:105:TYR:O	1:C:107:PRO:HD3	1.90	0.70
1:B:51:SER:HB2	1:B:73:TYR:CZ	2.27	0.69
1:D:94:VAL:HB	1:D:95:PRO:HD3	1.74	0.69
1:F:118:ASP:O	1:F:121:VAL:HG22	1.93	0.69
1:F:224:GLN:HE21	1:F:244:GLY:CA	2.05	0.69
1:F:320:GLN:HG2	1:F:321:ALA:H	1.58	0.68
1:D:353:ASN:ND2	1:D:355:SER:H	1.92	0.68
1:E:353:ASN:ND2	1:E:355:SER:H	1.92	0.68
1:E:45:LYS:O	1:E:49:GLN:HG3	1.94	0.68
1:A:325:MET:O	1:B:327:ASP:HB2	1.94	0.67
1:E:252:ILE:CG1	1:E:307:GLN:HG2	2.24	0.67
1:A:100:PHE:HA	1:A:103:PHE:CD2	2.30	0.67
1:A:353:ASN:HD22	1:A:355:SER:H	1.43	0.66
1:B:45:LYS:O	1:B:49:GLN:HG3	1.96	0.65
1:F:350:PRO:HG2	1:F:353:ASN:HB2	1.79	0.65
1:C:365:THR:O	1:C:369:GLN:HG3	1.96	0.65
1:C:291:ILE:HB	1:C:292:PRO:HD3	1.78	0.65
1:B:252:ILE:CG1	1:B:307:GLN:HG2	2.26	0.65
1:E:287:ASN:O	1:E:291:ILE:HG12	1.96	0.65
1:F:342:MET:HE2	1:F:342:MET:HA	1.77	0.65
1:C:289:CYS:SG	2:C:401:BH8:H30	2.37	0.65
1:C:100:PHE:HA	1:C:103:PHE:CD2	2.32	0.65
1:C:320:GLN:HB3	1:C:340:GLN:HE22	1.62	0.64
1:D:291:ILE:HB	1:D:292:PRO:HD3	1.79	0.64
1:F:245:ASP:HA	1:F:248:LEU:HD13	1.79	0.64
1:E:118:ASP:O	1:E:121:VAL:HG22	1.96	0.64
1:D:320:GLN:HG2	1:D:321:ALA:H	1.63	0.64
1:F:239:TYR:O	1:F:254:LEU:HD13	1.97	0.64
1:D:233:GLN:HA	1:D:236:TRP:NE1	2.12	0.64
1:B:318:LEU:CD1	1:B:319:GLY:H	2.10	0.64
1:A:252:ILE:CG1	1:A:307:GLN:HG2	2.27	0.64
1:D:232:PRO:HG2	1:D:235:VAL:HB	1.80	0.64
1:C:317:ARG:C	1:C:318:LEU:HG	2.22	0.63
1:E:325:MET:O	1:F:327:ASP:HB2	1.99	0.63
1:F:262:LEU:O	1:F:265:ASN:HB3	1.98	0.63
1:A:326:MET:HE3	1:B:291:ILE:HD11	1.79	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:233:GLN:HA	1:E:236:TRP:NE1	2.14	0.63
1:A:323:THR:HB	1:A:340:GLN:OE1	1.98	0.63
1:C:77:ARG:O	1:C:77:ARG:HD3	1.98	0.62
1:F:150:MET:O	1:F:154:MET:HG3	1.98	0.62
1:E:252:ILE:HG13	1:E:307:GLN:HG2	1.81	0.62
1:E:211:LEU:CD1	2:E:401:BH8:H21	2.28	0.62
1:B:252:ILE:HG13	1:B:307:GLN:HG2	1.82	0.62
1:E:54:PHE:O	1:E:58:ILE:HG13	2.00	0.62
1:B:278:SER:HB3	1:B:331:MET:HE2	1.81	0.62
1:D:180:GLY:HA2	2:D:401:BH8:H30	1.80	0.62
1:F:319:GLY:O	1:F:320:GLN:O	2.17	0.62
1:E:326:MET:HE3	1:F:291:ILE:CD1	2.30	0.62
1:A:94:VAL:HB	1:A:95:PRO:HD3	1.83	0.61
1:E:320:GLN:O	1:E:324:LEU:HG	2.00	0.61
1:E:125:PHE:N	1:E:126:PRO:HD2	2.16	0.61
1:D:45:LYS:O	1:D:49:GLN:HG3	2.01	0.60
1:D:85:ASP:OD2	1:D:87:THR:HB	2.01	0.60
1:D:211:LEU:CD1	2:D:401:BH8:H21	2.30	0.60
1:C:65:MET:HE2	1:C:192:PHE:CD2	2.37	0.60
1:F:320:GLN:O	1:F:324:LEU:HG	2.02	0.60
1:A:322:VAL:HB	1:A:340:GLN:NE2	2.17	0.59
1:F:224:GLN:HA	1:F:224:GLN:OE1	2.01	0.59
1:D:65:MET:HE2	1:D:192:PHE:CD2	2.37	0.59
1:A:233:GLN:HA	1:A:236:TRP:NE1	2.16	0.59
1:B:318:LEU:HD12	1:B:319:GLY:N	2.17	0.59
1:F:53:SER:HB2	2:F:401:BH8:OAF	2.02	0.59
1:F:260:ASN:HD22	1:F:353:ASN:ND2	2.00	0.59
1:C:94:VAL:HB	1:C:95:PRO:HD3	1.85	0.59
1:E:320:GLN:HG2	1:E:321:ALA:H	1.68	0.59
1:D:228:ARG:HH11	1:D:228:ARG:HG3	1.68	0.59
1:B:150:MET:HG3	1:B:174:TYR:O	2.02	0.58
1:D:325:MET:O	1:E:327:ASP:HB2	2.03	0.58
1:B:94:VAL:HB	1:B:95:PRO:HD3	1.85	0.58
1:D:343:GLU:CD	1:D:367:ARG:HH21	2.11	0.58
1:B:353:ASN:ND2	1:B:355:SER:H	2.02	0.58
1:A:343:GLU:OE1	1:A:367:ARG:NE	2.34	0.58
1:C:233:GLN:HA	1:C:236:TRP:NE1	2.19	0.58
1:A:207:MET:HE3	2:A:401:BH8:H28	1.85	0.57
1:C:110:ARG:HB3	1:C:126:PRO:HD3	1.87	0.57
1:A:343:GLU:CD	1:A:367:ARG:HH21	2.12	0.57
1:F:211:LEU:HD12	2:F:401:BH8:H15	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:179:VAL:HG11	2:E:401:BH8:H16	1.87	0.57
1:C:116:GLU:O	1:C:119:ARG:HD3	2.05	0.56
1:D:245:ASP:O	1:D:248:LEU:HB2	2.05	0.56
1:F:193:GLU:OE2	1:F:282:ASN:HB3	2.05	0.56
1:C:150:MET:HG3	1:C:174:TYR:O	2.05	0.56
1:E:319:GLY:HA3	1:E:323:THR:HG21	1.86	0.56
1:A:117:LYS:HD2	1:A:118:ASP:CG	2.30	0.56
1:F:316:ILE:N	1:F:316:ILE:HD12	2.20	0.56
1:F:100:PHE:HA	1:F:103:PHE:CD2	2.40	0.56
1:C:353:ASN:HD22	1:C:355:SER:H	1.53	0.56
1:F:211:LEU:HD12	2:F:401:BH8:H22	1.85	0.56
1:F:51:SER:HB2	1:F:73:TYR:CZ	2.41	0.56
1:C:323:THR:HB	1:C:340:GLN:OE1	2.07	0.55
1:A:320:GLN:HG2	1:A:321:ALA:H	1.71	0.55
1:E:90:VAL:O	1:E:94:VAL:HG23	2.06	0.55
1:E:110:ARG:HG2	1:E:110:ARG:NH2	2.22	0.55
1:A:194:ASP:CG	1:D:281:ARG:HH21	2.14	0.55
1:A:228:ARG:HG2	3:A:565:HOH:O	2.06	0.55
1:D:233:GLN:HA	1:D:236:TRP:CD1	2.41	0.55
1:E:110:ARG:CG	1:E:110:ARG:HH21	2.20	0.55
1:F:42:THR:HG22	1:F:46:TYR:CE2	2.42	0.55
1:D:319:GLY:HA3	1:D:323:THR:HG21	1.88	0.55
1:E:65:MET:HE2	1:E:192:PHE:CD2	2.42	0.54
1:F:323:THR:C	1:F:325:MET:H	2.15	0.54
1:D:66:ARG:NH1	1:F:367:ARG:O	2.40	0.54
1:D:287:ASN:O	1:D:291:ILE:HG12	2.07	0.54
1:D:51:SER:HB2	1:D:73:TYR:CZ	2.42	0.54
1:F:236:TRP:CG	1:F:243:LEU:HD13	2.43	0.54
1:D:365:THR:O	1:D:369:GLN:HG3	2.08	0.54
1:E:126:PRO:HA	1:E:129:SER:OG	2.08	0.54
1:A:327:ASP:HB2	1:C:325:MET:O	2.08	0.54
1:C:331:MET:HB3	1:C:332:PRO:HD3	1.90	0.54
1:F:174:TYR:HA	1:F:178:LEU:HD12	1.89	0.53
1:C:195:PRO:O	1:C:199:GLU:HG3	2.08	0.53
1:E:124:ASP:C	1:E:126:PRO:HD2	2.33	0.53
1:F:220:TYR:HB2	1:F:231:TRP:CZ2	2.44	0.53
1:A:353:ASN:HD21	1:A:355:SER:HB2	1.73	0.53
1:E:77:ARG:O	1:E:77:ARG:HD3	2.08	0.53
1:F:89:SER:OG	1:F:92:LYS:HB2	2.08	0.53
1:F:104:LEU:HD13	1:F:132:PHE:CE1	2.44	0.53
1:B:239:TYR:O	1:B:254:LEU:HD13	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:246:PHE:CD1	1:F:255:ALA:HA	2.44	0.53
1:A:51:SER:HB2	1:A:73:TYR:CZ	2.44	0.53
1:F:353:ASN:OD1	1:F:355:SER:N	2.31	0.52
1:B:325:MET:HB3	1:C:291:ILE:HD12	1.91	0.52
1:D:100:PHE:HA	1:D:103:PHE:CD2	2.45	0.52
1:E:331:MET:HB3	1:E:332:PRO:HD3	1.90	0.52
1:F:131:GLU:HA	1:F:134:ASN:HD22	1.74	0.52
1:A:320:GLN:CG	1:A:321:ALA:H	2.23	0.52
1:B:320:GLN:HG2	1:B:321:ALA:H	1.74	0.52
1:E:51:SER:HB2	1:E:73:TYR:OH	2.08	0.52
1:B:239:TYR:HE1	1:B:257:GLN:HB3	1.74	0.52
1:C:72:PHE:CZ	1:C:76:LEU:HD11	2.45	0.52
1:E:233:GLN:HA	1:E:236:TRP:CE2	2.44	0.52
1:B:190:SER:O	1:B:191:GLU:HB2	2.09	0.52
1:D:343:GLU:OE2	1:D:367:ARG:NE	2.37	0.52
1:B:365:THR:O	1:B:369:GLN:HG3	2.09	0.52
1:E:233:GLN:HA	1:E:236:TRP:CD1	2.44	0.52
1:E:325:MET:HA	1:E:325:MET:HE3	1.92	0.52
1:D:118:ASP:O	1:D:121:VAL:HG22	2.09	0.52
1:D:138:LYS:HE3	1:D:139:TYR:CZ	2.45	0.52
1:D:320:GLN:CG	1:D:321:ALA:H	2.23	0.52
1:F:72:PHE:CD1	1:F:143:ILE:HG23	2.45	0.52
1:B:233:GLN:HA	1:B:236:TRP:CD1	2.45	0.51
1:D:255:ALA:HB1	1:D:310:PHE:CZ	2.44	0.51
1:E:150:MET:O	1:E:154:MET:HG3	2.10	0.51
1:C:118:ASP:O	1:C:121:VAL:HG22	2.09	0.51
1:E:67:ASN:O	1:E:71:ILE:HG12	2.10	0.51
1:E:110:ARG:HG2	1:E:110:ARG:HH21	1.76	0.51
1:E:353:ASN:HD22	1:E:355:SER:H	1.56	0.51
1:F:101:HIS:CD2	1:F:148:ARG:HG3	2.46	0.51
1:D:105:TYR:O	1:D:107:PRO:HD3	2.09	0.51
1:F:75:VAL:HG23	1:F:76:LEU:N	2.25	0.51
1:C:51:SER:HB2	1:C:73:TYR:CZ	2.44	0.51
1:C:319:GLY:HA3	1:C:323:THR:HG21	1.93	0.51
1:F:79:LEU:HB2	1:F:100:PHE:CE2	2.45	0.51
1:A:318:LEU:HD12	1:A:318:LEU:N	2.26	0.51
1:B:255:ALA:HB1	1:B:310:PHE:CZ	2.46	0.51
1:D:190:SER:O	1:D:191:GLU:HB2	2.10	0.51
1:D:262:LEU:O	1:D:265:ASN:HB3	2.11	0.51
1:A:325:MET:HE3	1:A:325:MET:HA	1.91	0.51
1:C:353:ASN:HD21	1:C:355:SER:HB2	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:347:HIS:O	1:B:347:HIS:CG	2.63	0.51
1:A:233:GLN:HA	1:A:236:TRP:CD1	2.46	0.50
1:F:54:PHE:N	1:F:54:PHE:CD1	2.79	0.50
1:B:155:ALA:HB3	3:B:535:HOH:O	2.10	0.50
1:D:117:LYS:HG3	1:D:118:ASP:N	2.27	0.50
1:D:153:GLY:HA3	1:D:174:TYR:CG	2.46	0.50
1:C:183:LEU:CD1	2:C:401:BH8:H26	2.42	0.50
1:D:66:ARG:NH2	1:F:367:ARG:O	2.44	0.50
1:F:115:LYS:HE2	1:F:115:LYS:HA	1.94	0.50
1:D:125:PHE:N	1:D:126:PRO:CD	2.75	0.50
1:F:46:TYR:CE1	1:F:120:GLN:HG2	2.46	0.50
1:A:343:GLU:HG2	3:A:562:HOH:O	2.11	0.50
1:A:353:ASN:ND2	1:A:355:SER:HB2	2.27	0.49
1:E:100:PHE:HA	1:E:103:PHE:CE2	2.46	0.49
1:B:118:ASP:O	1:B:121:VAL:HG22	2.12	0.49
1:F:54:PHE:N	1:F:54:PHE:HD1	2.10	0.49
1:F:132:PHE:C	1:F:132:PHE:CD2	2.90	0.49
1:A:183:LEU:HD12	2:A:401:BH8:H23	1.94	0.49
1:D:238:ARG:HH22	1:D:261:GLU:CD	2.19	0.49
1:A:52:ARG:HG3	1:A:53:SER:N	2.28	0.49
1:C:42:THR:HG22	1:C:46:TYR:CE2	2.47	0.49
1:C:344:GLU:O	1:C:348:ARG:HD2	2.13	0.49
1:D:351:ASP:N	1:D:351:ASP:OD1	2.45	0.49
1:B:51:SER:HB2	1:B:73:TYR:OH	2.13	0.48
1:C:320:GLN:HG2	1:C:321:ALA:H	1.76	0.48
1:F:323:THR:HB	1:F:340:GLN:OE1	2.13	0.48
1:E:262:LEU:O	1:E:265:ASN:HB3	2.13	0.48
1:A:44:TYR:OH	1:A:66:ARG:HG2	2.12	0.48
1:C:325:MET:HA	1:C:325:MET:CE	2.39	0.48
1:F:263:ILE:O	1:F:267:LEU:HG	2.13	0.48
1:D:137:GLU:HB2	3:D:505:HOH:O	2.13	0.48
1:A:295:MET:HE3	1:C:321:ALA:CB	2.43	0.48
1:B:325:MET:CB	1:C:291:ILE:HD12	2.43	0.48
1:C:315:LEU:HD23	1:C:315:LEU:H	1.79	0.48
1:E:150:MET:HG3	1:E:174:TYR:O	2.14	0.48
1:F:208:GLY:HA2	2:F:401:BH8:H21	1.96	0.48
1:B:316:ILE:N	1:B:316:ILE:HD12	2.29	0.48
1:F:90:VAL:O	1:F:94:VAL:HG23	2.14	0.48
1:D:255:ALA:HB1	1:D:310:PHE:CE2	2.49	0.48
1:A:64:GLU:HG3	3:A:605:HOH:O	2.13	0.48
1:D:153:GLY:HA3	1:D:174:TYR:CD1	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:320:GLN:CG	1:F:321:ALA:H	2.23	0.48
1:B:220:TYR:HB2	1:B:231:TRP:CZ2	2.48	0.47
1:C:125:PHE:N	1:C:126:PRO:CD	2.78	0.47
1:B:138:LYS:HE3	1:B:139:TYR:CZ	2.50	0.47
1:B:209:LEU:HD23	1:B:269:HIS:CE1	2.49	0.47
1:E:343:GLU:OE2	1:E:367:ARG:NE	2.44	0.47
1:C:210:PHE:CE2	1:C:297:ILE:HG13	2.50	0.47
1:D:41:LYS:HD3	1:F:364:SER:OG	2.14	0.47
1:D:318:LEU:HD12	1:D:319:GLY:H	1.79	0.47
1:A:125:PHE:N	1:A:126:PRO:CD	2.77	0.47
1:D:331:MET:HB3	1:D:332:PRO:HD3	1.96	0.47
1:A:335:LYS:HB3	1:A:339:TYR:CE2	2.49	0.47
1:B:163:THR:HA	1:B:233:GLN:HB3	1.96	0.47
1:B:344:GLU:O	1:B:348:ARG:HG3	2.15	0.47
2:E:401:BH8:OAE	2:E:401:BH8:H9	2.15	0.47
1:B:248:LEU:HB2	1:B:251:ASN:HD22	1.80	0.47
1:D:42:THR:HG22	1:D:46:TYR:CE2	2.50	0.47
1:F:238:ARG:HH22	1:F:261:GLU:CD	2.23	0.47
1:D:318:LEU:HD12	1:D:319:GLY:N	2.30	0.46
1:E:194:ASP:OD1	1:E:195:PRO:HD2	2.15	0.46
1:C:252:ILE:HD11	1:C:307:GLN:HB3	1.98	0.46
1:B:125:PHE:N	1:B:126:PRO:CD	2.78	0.46
1:D:150:MET:HG3	1:D:174:TYR:O	2.16	0.46
1:E:193:GLU:OE2	1:E:282:ASN:HB3	2.16	0.46
1:F:224:GLN:NE2	1:F:244:GLY:CA	2.78	0.46
1:A:291:ILE:HB	1:A:292:PRO:HD3	1.98	0.46
1:E:73:TYR:CD1	1:E:73:TYR:C	2.93	0.46
1:F:308:GLN:HA	1:F:308:GLN:OE1	2.14	0.46
1:F:65:MET:O	1:F:69:VAL:HG23	2.14	0.46
1:F:72:PHE:HD1	1:F:143:ILE:HG23	1.80	0.46
1:A:183:LEU:CD1	2:A:401:BH8:H23	2.45	0.46
1:D:356:SER:O	1:D:360:ARG:HG3	2.16	0.46
1:F:320:GLN:HG3	3:F:518:HOH:O	2.15	0.46
1:E:323:THR:C	1:E:325:MET:H	2.24	0.46
1:F:170:LYS:O	1:F:173:HIS:HB3	2.16	0.46
1:B:90:VAL:HG23	3:B:571:HOH:O	2.14	0.45
1:E:85:ASP:HB3	1:E:88:ILE:HD12	1.99	0.45
1:A:320:GLN:O	1:A:324:LEU:HG	2.16	0.45
1:D:248:LEU:HD12	1:D:248:LEU:HA	1.72	0.45
1:D:150:MET:O	1:D:154:MET:HG3	2.16	0.45
1:E:280:LEU:O	1:E:281:ARG:HD3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:356:SER:HB2	1:E:360:ARG:HH12	1.81	0.45
1:C:85:ASP:OD1	1:C:114:SER:HA	2.17	0.45
1:A:65:MET:HE2	1:A:192:PHE:CD2	2.51	0.45
1:A:222:GLU:HB2	3:A:556:HOH:O	2.16	0.45
1:C:190:SER:O	1:C:191:GLU:HB2	2.16	0.45
1:B:320:GLN:NE2	1:C:52:ARG:HG2	2.31	0.45
1:E:101:HIS:CE1	1:E:148:ARG:HD3	2.52	0.45
1:E:338:ILE:O	1:E:342:MET:HG2	2.17	0.45
1:C:44:TYR:OH	1:C:66:ARG:HG2	2.16	0.45
1:C:233:GLN:HA	1:C:236:TRP:CD1	2.52	0.45
1:F:233:GLN:HE21	1:F:233:GLN:HB3	1.62	0.45
1:F:245:ASP:O	1:F:248:LEU:HB2	2.17	0.45
1:A:323:THR:HG23	1:A:324:LEU:N	2.32	0.44
1:E:229:GLU:HB3	1:E:243:LEU:HD23	1.99	0.44
1:B:221:LEU:HG	1:B:225:GLN:OE1	2.17	0.44
1:C:252:ILE:CG2	1:C:253:ASP:N	2.80	0.44
1:C:255:ALA:HB1	1:C:310:PHE:CZ	2.52	0.44
1:C:315:LEU:H	1:C:315:LEU:CD2	2.27	0.44
1:F:306:ASN:ND2	1:F:308:GLN:HB2	2.32	0.44
1:B:353:ASN:HA	1:B:354:PRO:HD3	1.80	0.44
1:A:190:SER:O	1:A:191:GLU:HB2	2.17	0.44
1:E:327:ASP:C	1:E:327:ASP:OD1	2.58	0.44
1:B:233:GLN:HB3	1:B:233:GLN:HE21	1.52	0.44
1:B:331:MET:HB3	1:B:332:PRO:HD3	2.00	0.44
1:D:67:ASN:O	1:D:71:ILE:HG12	2.18	0.44
1:F:255:ALA:HB1	1:F:310:PHE:CE2	2.52	0.44
1:A:252:ILE:HG23	1:A:253:ASP:N	2.33	0.44
1:F:255:ALA:HB1	1:F:310:PHE:CZ	2.53	0.44
1:F:256:VAL:O	1:F:259:LEU:HB3	2.18	0.44
1:A:252:ILE:HG12	1:A:307:GLN:HG2	1.98	0.43
1:A:318:LEU:N	1:A:318:LEU:CD1	2.81	0.43
1:B:73:TYR:CD1	1:B:73:TYR:C	2.95	0.43
1:A:180:GLY:HA2	2:A:401:BH8:H24	2.00	0.43
1:F:224:GLN:C	1:F:226:GLY:H	2.26	0.43
2:D:401:BH8:H7	3:D:556:HOH:O	2.18	0.43
1:E:320:GLN:CG	1:E:321:ALA:H	2.30	0.43
1:F:65:MET:HE2	1:F:192:PHE:CD2	2.52	0.43
1:B:239:TYR:CE1	1:B:257:GLN:HB3	2.53	0.43
1:E:356:SER:HB2	1:E:360:ARG:NH1	2.33	0.43
1:A:207:MET:CE	2:A:401:BH8:H28	2.49	0.43
1:D:51:SER:HB2	1:D:73:TYR:OH	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:252:ILE:HG12	1:E:307:GLN:HG2	1.99	0.43
1:A:319:GLY:CA	1:A:323:THR:HG21	2.46	0.43
1:C:157:PHE:CE2	1:C:160:LYS:HE3	2.53	0.43
1:F:148:ARG:HG2	1:F:148:ARG:HH21	1.83	0.43
1:F:252:ILE:CG1	1:F:307:GLN:HG2	2.49	0.43
1:B:51:SER:OG	1:B:54:PHE:HB2	2.19	0.42
1:B:238:ARG:HH11	1:B:238:ARG:CB	2.32	0.42
1:D:51:SER:OG	1:D:54:PHE:HB2	2.19	0.42
1:F:323:THR:C	1:F:325:MET:N	2.75	0.42
1:A:207:MET:HB3	2:A:401:BH8:H27	2.00	0.42
1:A:295:MET:HE3	1:C:321:ALA:HB3	2.00	0.42
1:B:363:ILE:O	1:B:367:ARG:HG3	2.18	0.42
1:C:333:ALA:O	1:C:337:ILE:HG13	2.19	0.42
1:D:233:GLN:O	1:D:237:SER:HB3	2.19	0.42
1:F:194:ASP:CG	1:F:195:PRO:HD2	2.44	0.42
1:A:318:LEU:CD1	1:A:318:LEU:H	2.33	0.42
1:A:326:MET:HE3	1:B:291:ILE:CD1	2.48	0.42
1:B:281:ARG:NH2	1:F:194:ASP:OD2	2.49	0.42
1:B:367:ARG:O	1:C:66:ARG:NH1	2.47	0.42
1:D:342:MET:HE2	1:D:342:MET:HA	2.02	0.42
1:A:316:ILE:HD12	1:A:316:ILE:N	2.34	0.42
1:C:254:LEU:HD23	1:C:257:GLN:NE2	2.35	0.42
1:F:75:VAL:CG2	1:F:76:LEU:N	2.83	0.42
1:F:164:SER:O	1:F:167:GLU:HB2	2.20	0.42
1:F:200:ASP:OD1	1:F:203:ARG:HG3	2.20	0.42
1:A:353:ASN:HD22	1:A:355:SER:N	2.15	0.42
1:B:344:GLU:O	1:B:348:ARG:NH1	2.53	0.42
1:C:183:LEU:HD12	2:C:401:BH8:H26	2.01	0.42
1:C:213:LYS:O	1:C:217:ILE:HG13	2.20	0.42
1:F:343:GLU:OE1	1:F:343:GLU:HA	2.20	0.42
1:F:359:THR:O	1:F:362:ILE:HG22	2.20	0.42
1:D:246:PHE:CD1	1:D:255:ALA:HA	2.54	0.42
1:E:319:GLY:CA	1:E:323:THR:HG21	2.48	0.42
1:A:263:ILE:CD1	1:A:304:TYR:HA	2.50	0.42
1:B:263:ILE:HD12	1:B:304:TYR:HA	2.02	0.42
1:C:211:LEU:HD12	2:C:401:BH8:H20	2.02	0.42
1:D:100:PHE:HA	1:D:103:PHE:CE2	2.55	0.42
1:E:306:ASN:OD1	1:E:308:GLN:HB2	2.20	0.42
1:E:326:MET:HA	1:F:327:ASP:HB2	2.02	0.42
1:F:297:ILE:CD1	1:F:338:ILE:HG12	2.50	0.42
1:A:252:ILE:HG13	1:A:307:GLN:HG2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:342:MET:HE2	1:A:342:MET:HA	2.02	0.41
1:D:317:ARG:H	1:D:317:ARG:HG2	1.35	0.41
1:D:322:VAL:O	1:D:326:MET:HG3	2.20	0.41
1:E:252:ILE:HG23	1:E:253:ASP:N	2.34	0.41
1:A:185:ARG:NH1	3:A:530:HOH:O	2.53	0.41
1:A:331:MET:HB3	1:A:332:PRO:HD3	2.02	0.41
1:C:252:ILE:HD12	1:C:307:GLN:HG2	2.02	0.41
1:A:349:ILE:HA	1:A:350:PRO:HD3	1.87	0.41
1:F:50:THR:HB	1:F:77:ARG:NH1	2.34	0.41
1:F:239:TYR:HE1	1:F:257:GLN:HE21	1.68	0.41
1:A:67:ASN:O	1:A:71:ILE:HG12	2.21	0.41
1:C:146:ILE:HG22	1:C:150:MET:HE3	2.01	0.41
1:C:210:PHE:HE2	1:C:297:ILE:HG13	1.86	0.41
1:C:353:ASN:ND2	1:C:355:SER:HB2	2.35	0.41
1:E:101:HIS:ND1	1:E:148:ARG:HB2	2.36	0.41
1:E:138:LYS:HE3	1:E:139:TYR:CZ	2.54	0.41
1:D:353:ASN:HD21	1:D:355:SER:HB2	1.85	0.41
1:F:348:ARG:O	1:F:350:PRO:HD3	2.20	0.41
1:B:368:THR:O	1:B:368:THR:HG22	2.20	0.41
1:C:320:GLN:CG	1:C:321:ALA:H	2.33	0.41
1:F:99:ASN:O	1:F:100:PHE:C	2.62	0.41
1:A:368:THR:O	1:A:369:GLN:C	2.63	0.41
1:B:291:ILE:O	1:B:295:MET:HG3	2.21	0.41
1:F:125:PHE:N	1:F:126:PRO:CD	2.84	0.41
1:F:327:ASP:OD1	1:F:327:ASP:C	2.62	0.41
1:A:276:TYR:CD1	1:A:276:TYR:C	2.98	0.41
1:B:316:ILE:N	1:B:316:ILE:CD1	2.83	0.41
1:D:320:GLN:HG2	1:D:321:ALA:N	2.33	0.41
1:F:362:ILE:HG23	1:F:363:ILE:N	2.36	0.41
1:A:65:MET:HE2	1:A:192:PHE:CG	2.55	0.41
1:A:150:MET:HE3	1:A:150:MET:HB2	1.96	0.41
1:B:54:PHE:O	1:B:58:ILE:HG13	2.21	0.41
1:B:344:GLU:HG2	1:B:348:ARG:HH12	1.86	0.41
1:C:77:ARG:HD3	1:C:77:ARG:C	2.46	0.41
1:C:246:PHE:CE1	1:C:255:ALA:HA	2.56	0.41
1:E:183:LEU:CD1	2:E:401:BH8:H25	2.51	0.41
1:E:183:LEU:HD12	2:E:401:BH8:H25	2.03	0.41
1:A:52:ARG:CG	1:A:53:SER:N	2.83	0.41
1:C:119:ARG:O	1:C:123:GLU:HG3	2.21	0.41
1:B:320:GLN:CG	1:B:321:ALA:H	2.33	0.40
1:E:344:GLU:O	1:E:348:ARG:HG3	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:77:ARG:O	1:F:77:ARG:HD3	2.21	0.40
1:C:51:SER:HB2	1:C:73:TYR:CE1	2.56	0.40
1:F:355:SER:O	1:F:358:LYS:N	2.54	0.40
1:B:108:ASP:HB2	3:B:566:HOH:O	2.21	0.40
1:C:338:ILE:O	1:C:342:MET:HG2	2.22	0.40
1:B:238:ARG:HB2	1:B:238:ARG:NH1	2.37	0.40
1:E:183:LEU:HD12	2:E:401:BH8:CAL	2.51	0.40
1:F:55:ALA:O	1:F:59:GLN:HG3	2.21	0.40
1:C:142:VAL:O	1:C:146:ILE:HG12	2.21	0.40
1:E:356:SER:O	1:E:360:ARG:HG3	2.20	0.40
1:F:291:ILE:HB	1:F:292:PRO:HD3	2.03	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%)	321 (97%)	10 (3%)	1 (0%)	36	41
1	B	332/360 (92%)	320 (96%)	9 (3%)	3 (1%)	14	13
1	C	332/360 (92%)	314 (95%)	14 (4%)	4 (1%)	10	8
1	D	332/360 (92%)	314 (95%)	15 (4%)	3 (1%)	14	13
1	E	332/360 (92%)	311 (94%)	18 (5%)	3 (1%)	14	13
1	F	332/360 (92%)	310 (93%)	18 (5%)	4 (1%)	10	8
All	All	1992/2160 (92%)	1890 (95%)	84 (4%)	18 (1%)	14	13

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	318	LEU

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Mol	Chain	Res	Type
1	C	317	ARG
1	D	317	ARG
1	F	320	GLN
1	B	320	GLN
1	E	320	GLN
1	E	318	LEU
1	A	320	GLN
1	D	320	GLN
1	C	63	GLY
1	C	249	PRO
1	C	318	LEU
1	D	63	GLY
1	F	66	ARG
1	F	63	GLY
1	F	200	ASP
1	E	63	GLY
1	B	63	GLY

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	297/320 (93%)	287 (97%)	10 (3%)	32	42
1	B	297/320 (93%)	286 (96%)	11 (4%)	30	39
1	C	297/320 (93%)	285 (96%)	12 (4%)	28	36
1	D	297/320 (93%)	287 (97%)	10 (3%)	32	42
1	E	297/320 (93%)	293 (99%)	4 (1%)	61	75
1	F	297/320 (93%)	281 (95%)	16 (5%)	20	23
All	All	1782/1920 (93%)	1719 (96%)	63 (4%)	32	41

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	110	ARG

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Mol	Chain	Res	Type
1	A	117	LYS
1	A	119	ARG
1	A	120	GLN
1	A	137	GLU
1	A	203	ARG
1	A	233	GLN
1	A	241	LYS
1	A	326	MET
1	A	369	GLN
1	B	53	SER
1	B	77	ARG
1	B	83	GLU
1	B	156	GLU
1	B	233	GLN
1	B	250	GLU
1	B	318	LEU
1	B	323	THR
1	B	340	GLN
1	B	347	HIS
1	B	352	SER
1	C	110	ARG
1	C	170	LYS
1	C	194	ASP
1	C	242	LYS
1	C	249	PRO
1	C	252	ILE
1	C	314	VAL
1	C	315	LEU
1	C	318	LEU
1	C	325	MET
1	C	348	ARG
1	C	357	SER
1	D	52	ARG
1	D	248	LEU
1	D	249	PRO
1	D	316	ILE
1	D	317	ARG
1	D	318	LEU
1	D	325	MET
1	D	327	ASP
1	D	352	SER
1	D	357	SER

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Mol	Chain	Res	Type
1	E	52	ARG
1	E	83	GLU
1	E	110	ARG
1	E	326	MET
1	F	45	LYS
1	F	54	PHE
1	F	83	GLU
1	F	91	GLU
1	F	113	GLU
1	F	115	LYS
1	F	123	GLU
1	F	233	GLN
1	F	250	GLU
1	F	308	GLN
1	F	315	LEU
1	F	318	LEU
1	F	326	MET
1	F	344	GLU
1	F	347	HIS
1	F	351	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (56) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	48	ASN
1	A	59	GLN
1	A	67	ASN
1	A	120	GLN
1	A	225	GLN
1	A	233	GLN
1	A	260	ASN
1	A	283	GLN
1	A	305	ASN
1	A	353	ASN
1	A	369	GLN
1	B	48	ASN
1	B	59	GLN
1	B	67	ASN
1	B	212	GLN
1	B	224	GLN
1	B	225	GLN
1	B	233	GLN

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Mol	Chain	Res	Type
1	B	251	ASN
1	B	257	GLN
1	B	283	GLN
1	B	320	GLN
1	C	99	ASN
1	C	120	GLN
1	C	225	GLN
1	C	233	GLN
1	C	260	ASN
1	C	283	GLN
1	C	305	ASN
1	C	353	ASN
1	D	48	ASN
1	D	59	GLN
1	D	212	GLN
1	D	224	GLN
1	D	225	GLN
1	D	257	GLN
1	D	283	GLN
1	D	293	GLN
1	D	353	ASN
1	E	215	ASN
1	E	224	GLN
1	E	225	GLN
1	E	293	GLN
1	F	48	ASN
1	F	59	GLN
1	F	101	HIS
1	F	120	GLN
1	F	134	ASN
1	F	166	GLN
1	F	224	GLN
1	F	225	GLN
1	F	233	GLN
1	F	257	GLN
1	F	283	GLN
1	F	293	GLN
1	F	369	GLN

5.3.3 RNA ⓘ

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 oligosaccharides 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	BH8	F	401	-	24,25,25	2.34	5 (20%)	31,34,34	0.97	2 (6%)
2	BH8	A	401	-	24,25,25	2.35	5 (20%)	31,34,34	1.00	2 (6%)
2	BH8	B	401	-	24,25,25	2.34	5 (20%)	31,34,34	0.97	2 (6%)
2	BH8	D	401	-	24,25,25	2.35	5 (20%)	31,34,34	1.02	2 (6%)
2	BH8	C	401	-	24,25,25	2.34	5 (20%)	31,34,34	1.04	2 (6%)
2	BH8	E	401	-	24,25,25	2.35	5 (20%)	31,34,34	1.02	2 (6%)

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	BH8	F	401	-	-	2/24/26/26	0/1/1/1
2	BH8	A	401	-	-	4/24/26/26	0/1/1/1
2	BH8	B	401	-	-	8/24/26/26	0/1/1/1
2	BH8	D	401	-	-	9/24/26/26	0/1/1/1
2	BH8	C	401	-	-	8/24/26/26	0/1/1/1
2	BH8	E	401	-	-	8/24/26/26	0/1/1/1

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	401	BH8	PAY-OAG	7.09	1.61	1.49
2	A	401	BH8	PAY-OAG	7.08	1.61	1.49
2	B	401	BH8	PAY-OAG	7.08	1.61	1.49
2	C	401	BH8	PAY-OAG	7.07	1.61	1.49
2	D	401	BH8	PAY-OAG	7.06	1.61	1.49
2	F	401	BH8	PAY-OAG	7.04	1.61	1.49
2	E	401	BH8	PAX-OAE	4.19	1.61	1.54
2	B	401	BH8	PAX-OAD	4.18	1.61	1.54
2	E	401	BH8	PAX-OAD	4.17	1.61	1.54
2	B	401	BH8	PAY-OAF	4.17	1.61	1.54
2	D	401	BH8	PAX-OAD	4.16	1.61	1.54
2	A	401	BH8	PAX-OAD	4.16	1.61	1.54
2	F	401	BH8	PAX-OAD	4.15	1.61	1.54
2	B	401	BH8	PAX-OAE	4.14	1.61	1.54
2	E	401	BH8	PAY-OAF	4.14	1.61	1.54
2	A	401	BH8	PAY-OAF	4.13	1.61	1.54
2	D	401	BH8	PAX-OAE	4.13	1.61	1.54
2	C	401	BH8	PAX-OAE	4.12	1.61	1.54
2	C	401	BH8	PAX-OAD	4.12	1.61	1.54
2	A	401	BH8	PAX-OAE	4.11	1.61	1.54
2	F	401	BH8	PAX-OAE	4.10	1.61	1.54
2	F	401	BH8	PAY-OAF	4.10	1.61	1.54
2	D	401	BH8	PAY-OAF	4.10	1.61	1.54
2	C	401	BH8	PAY-OAF	4.09	1.61	1.54
2	A	401	BH8	PAY-OAC	-4.08	1.48	1.54
2	E	401	BH8	PAY-OAC	-4.08	1.48	1.54
2	D	401	BH8	PAY-OAC	-4.03	1.48	1.54
2	B	401	BH8	PAY-OAC	-4.02	1.48	1.54
2	F	401	BH8	PAY-OAC	-4.01	1.48	1.54
2	C	401	BH8	PAY-OAC	-4.00	1.48	1.54

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	BH8	OAF-PAY-OAG	-2.97	105.99	113.45
2	F	401	BH8	OAF-PAY-OAG	-2.93	106.11	113.45
2	D	401	BH8	OAF-PAY-OAG	-2.89	106.19	113.45
2	C	401	BH8	OAF-PAY-OAG	-2.88	106.22	113.45
2	E	401	BH8	OAF-PAY-OAG	-2.85	106.30	113.45
2	B	401	BH8	OAF-PAY-OAG	-2.69	106.71	113.45
2	D	401	BH8	OAG-PAY-CAU	-2.43	105.39	112.29
2	F	401	BH8	OAG-PAY-CAU	-2.30	105.75	112.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	BH8	OAG-PAY-CAU	-2.29	105.79	112.29
2	C	401	BH8	OAG-PAY-CAU	-2.24	105.92	112.29
2	E	401	BH8	OAG-PAY-CAU	-2.21	106.01	112.29
2	A	401	BH8	OAG-PAY-CAU	-2.02	106.56	112.29

There are no chirality outliers.

All (39) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	401	BH8	PAY-CAU-PAX-OAD
2	B	401	BH8	CAT-CAU-PAX-OAD
2	B	401	BH8	PAY-CAU-PAX-OAE
2	B	401	BH8	CAT-CAU-PAX-OAE
2	B	401	BH8	PAY-CAU-PAX-OAB
2	B	401	BH8	CAT-CAU-PAX-OAB
2	C	401	BH8	PAY-CAU-PAX-OAD
2	C	401	BH8	CAT-CAU-PAX-OAD
2	C	401	BH8	PAY-CAU-PAX-OAE
2	C	401	BH8	PAY-CAU-PAX-OAB
2	C	401	BH8	CAT-CAU-PAX-OAB
2	D	401	BH8	PAY-CAU-PAX-OAD
2	D	401	BH8	CAT-CAU-PAX-OAD
2	D	401	BH8	PAY-CAU-PAX-OAB
2	D	401	BH8	CAT-CAU-PAX-OAB
2	E	401	BH8	PAY-CAU-PAX-OAD
2	E	401	BH8	CAT-CAU-PAX-OAD
2	E	401	BH8	PAY-CAU-PAX-OAB
2	E	401	BH8	CAT-CAU-PAX-OAB
2	E	401	BH8	CAQ-CAR-CAS-NAV
2	A	401	BH8	CAK-CAL-CAM-CAN
2	F	401	BH8	CAQ-CAR-CAS-NAV
2	A	401	BH8	CAL-CAM-CAN-CAO
2	D	401	BH8	CAK-CAL-CAM-CAN
2	A	401	BH8	CAP-CAQ-CAR-CAS
2	D	401	BH8	PAY-CAU-PAX-OAE
2	E	401	BH8	PAY-CAU-PAX-OAE
2	C	401	BH8	CAR-CAS-NAV-CAH
2	B	401	BH8	CAU-CAT-NAW-CAI
2	C	401	BH8	CAU-CAT-NAW-CAI
2	D	401	BH8	CAU-CAT-NAW-CAI
2	F	401	BH8	CAU-CAT-NAW-CAI
2	C	401	BH8	CAT-CAU-PAX-OAE

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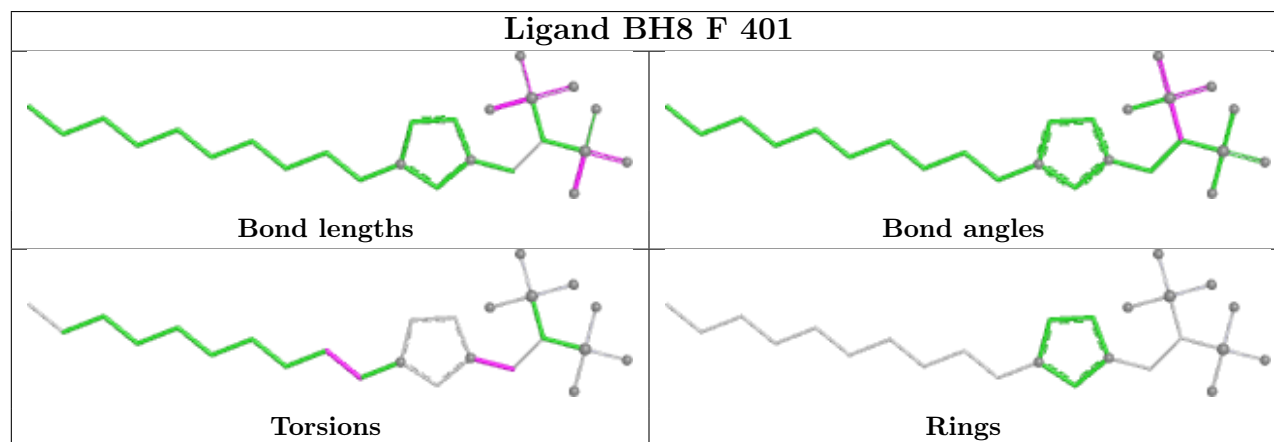
Mol	Chain	Res	Type	Atoms
2	D	401	BH8	CAT-CAU-PAX-OAE
2	E	401	BH8	CAT-CAU-PAX-OAE
2	B	401	BH8	CAN-CAO-CAP-CAQ
2	E	401	BH8	CAA-CAK-CAL-CAM
2	A	401	BH8	CAR-CAS-NAV-CAJ
2	D	401	BH8	CAT-CAU-PAY-OAG

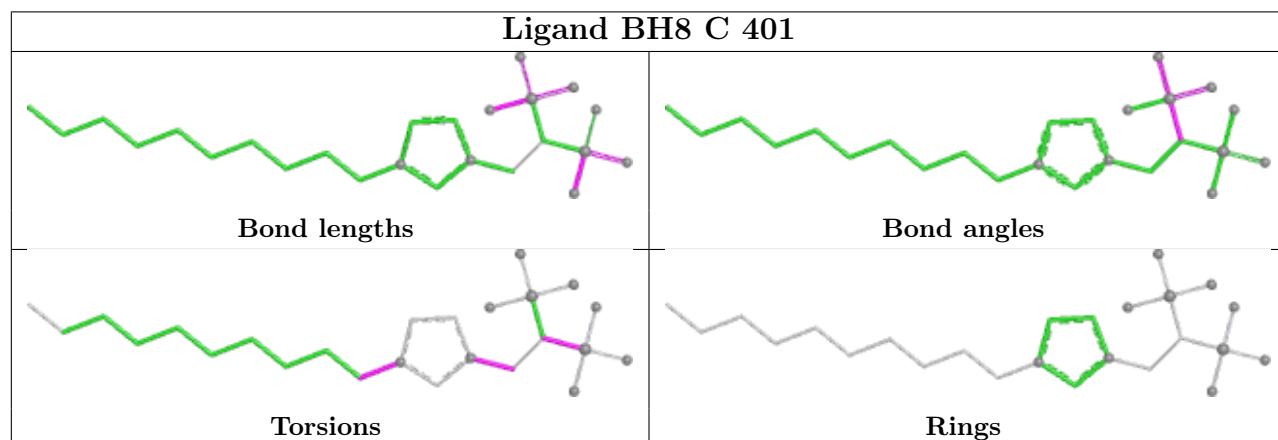
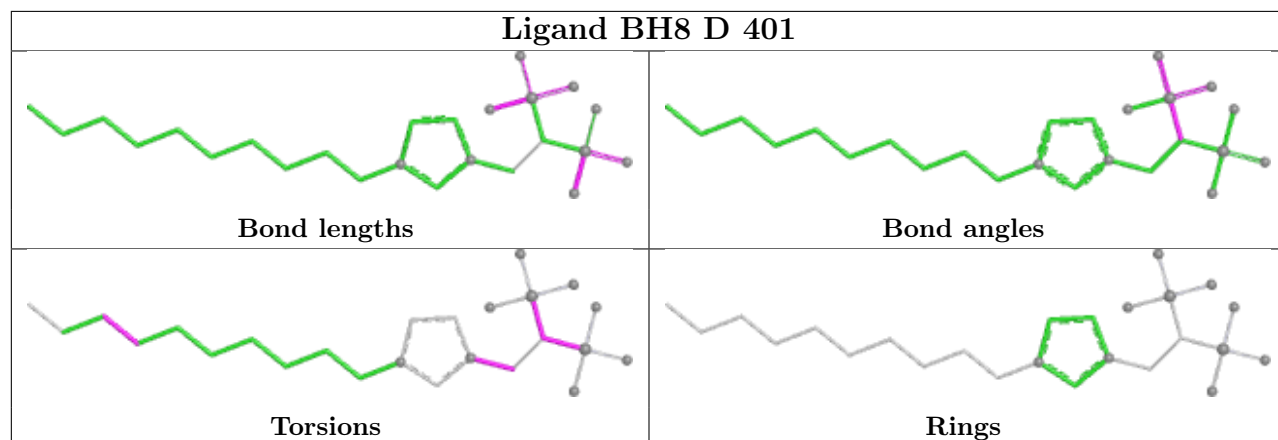
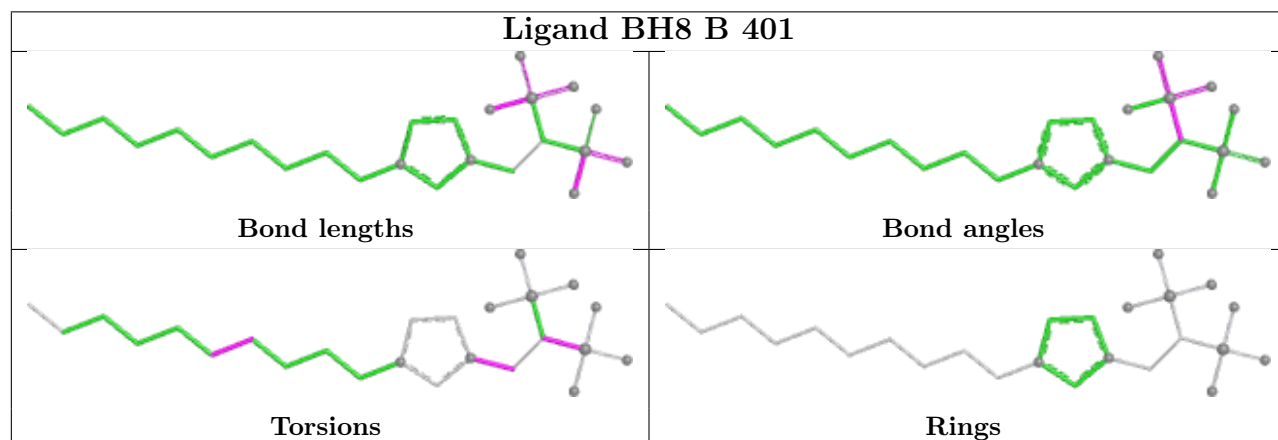
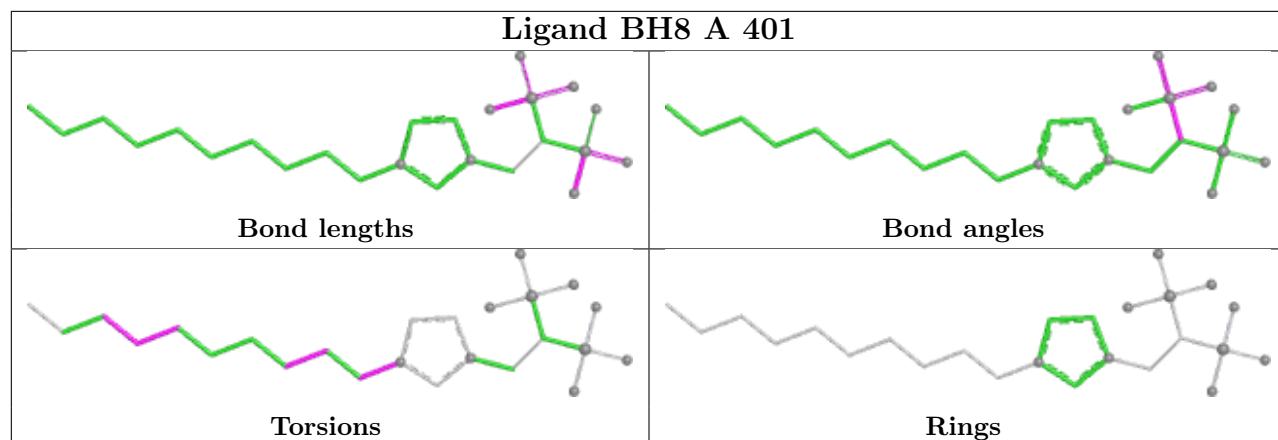
There are no ring outliers.

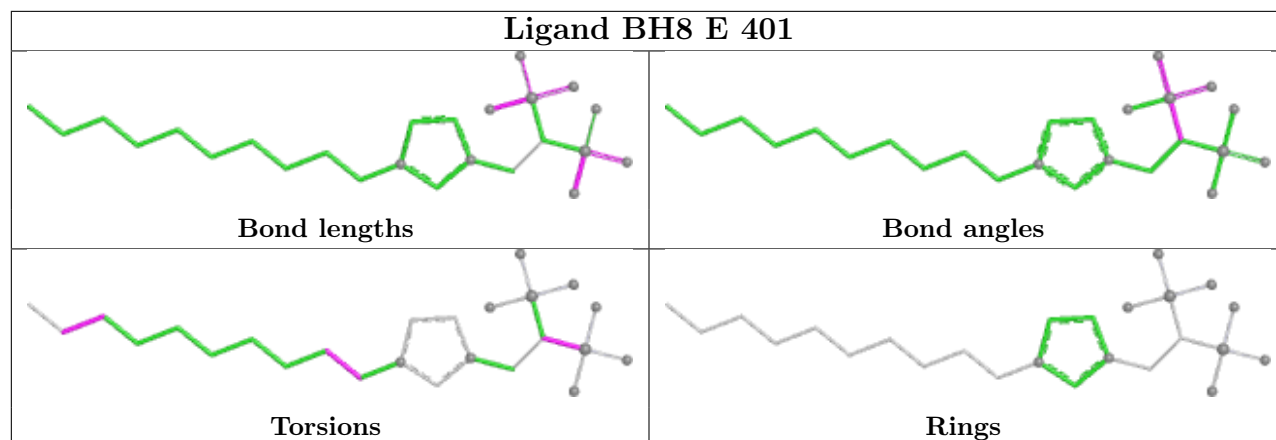
5 monomers are involved in 25 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	401	BH8	4	0
2	A	401	BH8	6	0
2	D	401	BH8	4	0
2	C	401	BH8	4	0
2	E	401	BH8	7	0

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







5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	334/360 (92%)	-0.28	10 (2%) 52 50	17, 32, 64, 105	0
1	B	334/360 (92%)	-0.09	7 (2%) 63 61	22, 35, 74, 108	0
1	C	334/360 (92%)	-0.08	8 (2%) 59 57	25, 38, 70, 101	0
1	D	334/360 (92%)	0.12	9 (2%) 56 53	24, 45, 77, 104	0
1	E	334/360 (92%)	0.32	6 (1%) 67 65	30, 49, 86, 109	0
1	F	334/360 (92%)	0.83	48 (14%) 6 4	27, 66, 107, 119	0
All	All	2004/2160 (92%)	0.14	88 (4%) 39 36	17, 43, 94, 119	0

All (88) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	318	LEU	4.7
1	B	318	LEU	4.6
1	D	36	LEU	4.3
1	F	54	PHE	4.1
1	C	318	LEU	3.9
1	F	230	PHE	3.9
1	F	318	LEU	3.8
1	A	318	LEU	3.7
1	C	315	LEU	3.6
1	C	36	LEU	3.4
1	E	36	LEU	3.3
1	F	162	VAL	3.3
1	E	324	LEU	3.2
1	D	324	LEU	3.1
1	F	158	LEU	3.1
1	B	54	PHE	3.1
1	F	36	LEU	3.0
1	F	232	PRO	2.9
1	F	104	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
1	F	247	ALA	2.9
1	F	315	LEU	2.9
1	E	319	GLY	2.8
1	B	320	GLN	2.8
1	D	54	PHE	2.8
1	A	315	LEU	2.7
1	F	322	VAL	2.6
1	F	248	LEU	2.6
1	D	318	LEU	2.6
1	B	36	LEU	2.6
1	F	321	ALA	2.6
1	B	368	THR	2.5
1	F	241	LYS	2.5
1	F	252	ILE	2.5
1	F	243	LEU	2.5
1	B	319	GLY	2.5
1	F	319	GLY	2.5
1	F	103	PHE	2.5
1	F	90	VAL	2.5
1	F	231	TRP	2.5
1	C	314	VAL	2.4
1	F	310	PHE	2.4
1	F	236	TRP	2.4
1	B	315	LEU	2.4
1	F	220	TYR	2.4
1	F	228	ARG	2.4
1	F	226	GLY	2.4
1	D	322	VAL	2.4
1	F	235	VAL	2.4
1	F	314	VAL	2.4
1	A	324	LEU	2.4
1	F	96	LEU	2.4
1	F	246	PHE	2.3
1	D	228	ARG	2.3
1	F	240	VAL	2.3
1	D	252	ILE	2.3
1	F	324	LEU	2.2
1	F	249	PRO	2.2
1	F	227	GLY	2.2
1	F	242	LYS	2.2
1	F	320	GLN	2.2
1	F	369	GLN	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	316	ILE	2.2
1	C	323	THR	2.2
1	F	174	TYR	2.1
1	F	259	LEU	2.1
1	F	251	ASN	2.1
1	A	369	GLN	2.1
1	D	53	SER	2.1
1	E	322	VAL	2.1
1	F	100	PHE	2.1
1	F	125	PHE	2.1
1	F	317	ARG	2.1
1	F	163	THR	2.0
1	E	321	ALA	2.0
1	A	320	GLN	2.0
1	A	36	LEU	2.0
1	F	82	LEU	2.0
1	F	349	ILE	2.0
1	F	239	TYR	2.0
1	A	228	ARG	2.0
1	A	322	VAL	2.0
1	D	240	VAL	2.0
1	F	157	PHE	2.0
1	A	321	ALA	2.0
1	C	321	ALA	2.0
1	A	159	ASP	2.0
1	C	327	ASP	2.0
1	F	345	ILE	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,

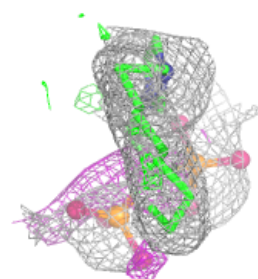
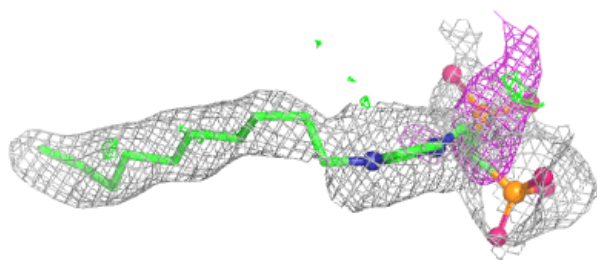
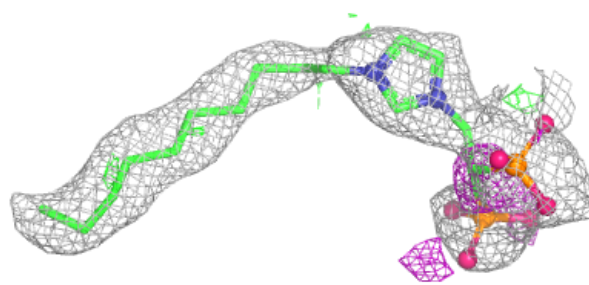
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	BH8	D	401	25/25	0.70	0.18	48,86,98,98	0
2	BH8	A	401	25/25	0.71	0.21	41,84,95,96	0
2	BH8	F	401	25/25	0.76	0.16	57,89,99,99	0
2	BH8	E	401	25/25	0.82	0.15	58,86,93,93	0
2	BH8	C	401	25/25	0.83	0.14	39,71,80,81	0
2	BH8	B	401	25/25	0.85	0.12	32,60,68,68	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.

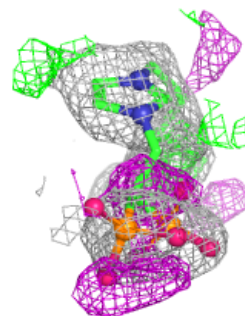
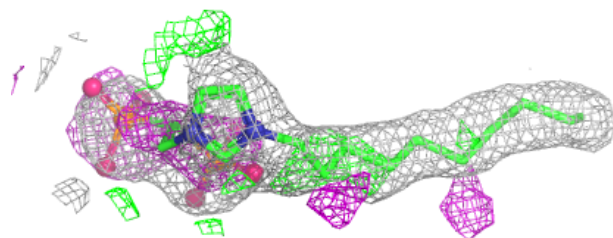
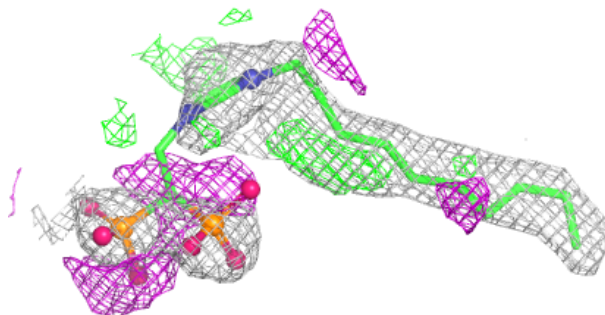
Electron density around BH8 D 401:

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

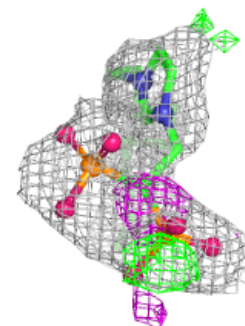
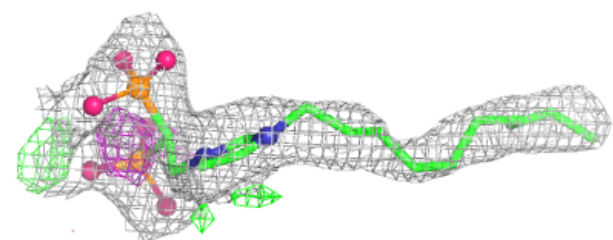
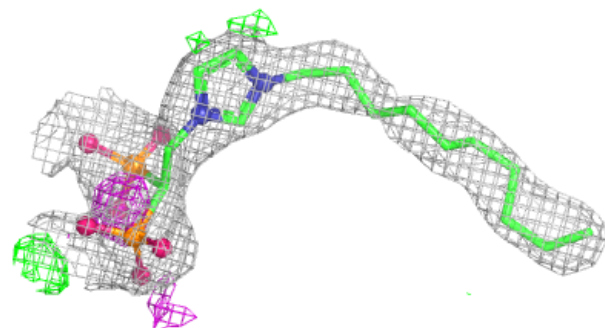


Electron density around BH8 A 401:

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

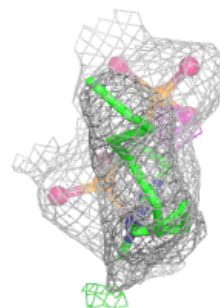
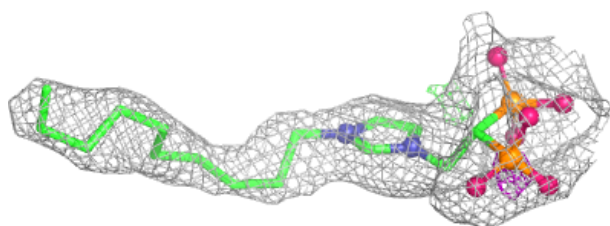
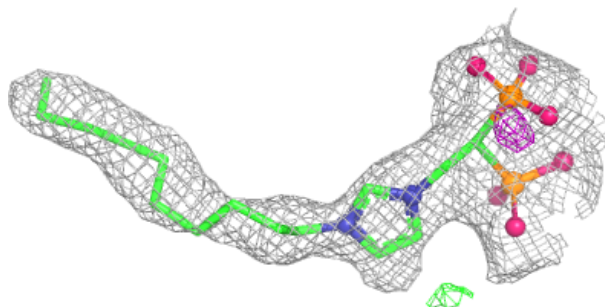
**Electron density around BH8 F 401:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

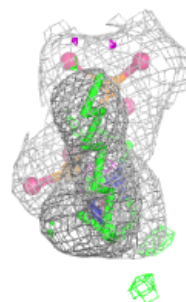
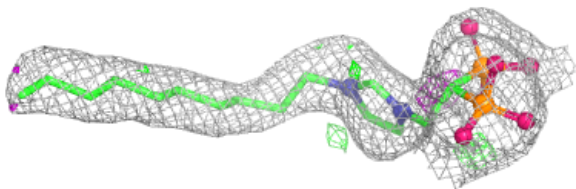
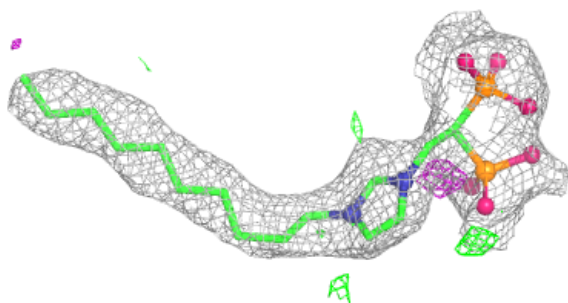


Electron density around BH8 E 401:

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and green (positive)

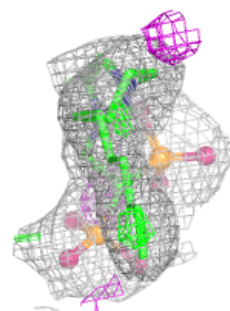
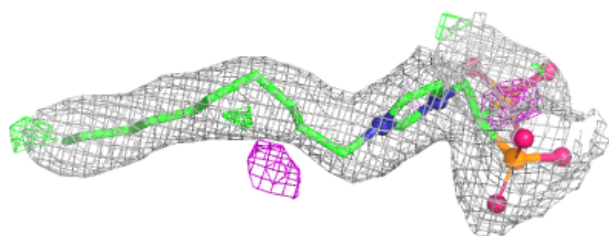
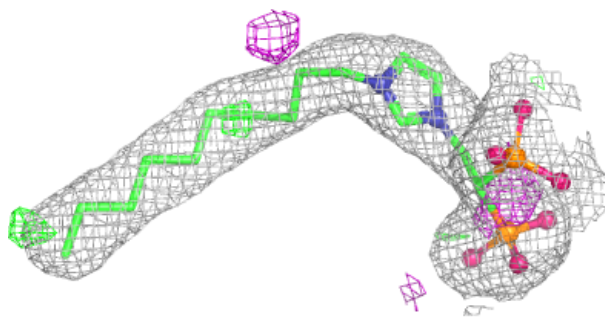
**Electron density around BH8 C 401:**

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



Electron density around BH8 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)



6.5 Other polymers ⓘ

There are no such residues in this entry.