



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:39 PM UTC

PDB ID : 3WCH / pdb_00003wch
Title : The complex structure of HsSQS with ligand BPH1237
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 : 2.50 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

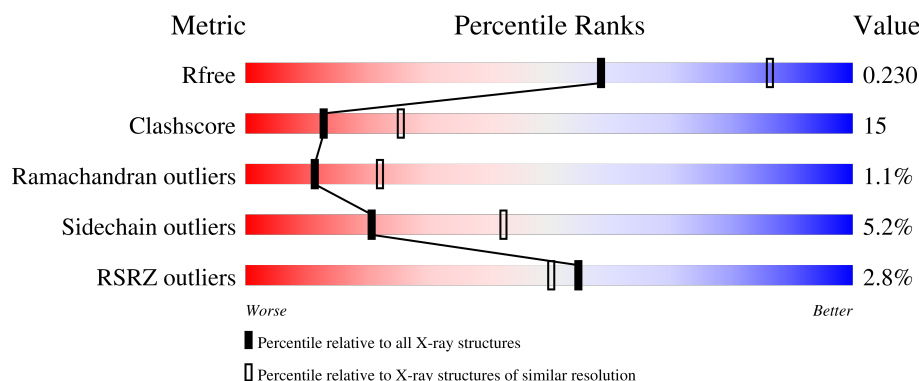
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	 2% 66% 25% 7%
1	B	360	 2% 63% 27% 7%
1	C	360	 % 67% 22% 7%
1	D	360	 3% 61% 27% 7%
1	E	360	 4% 66% 24% 7%

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Mol	Chain	Length	Quality of chain
1	F	360	4% 56% 31% 5% 7%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 16820 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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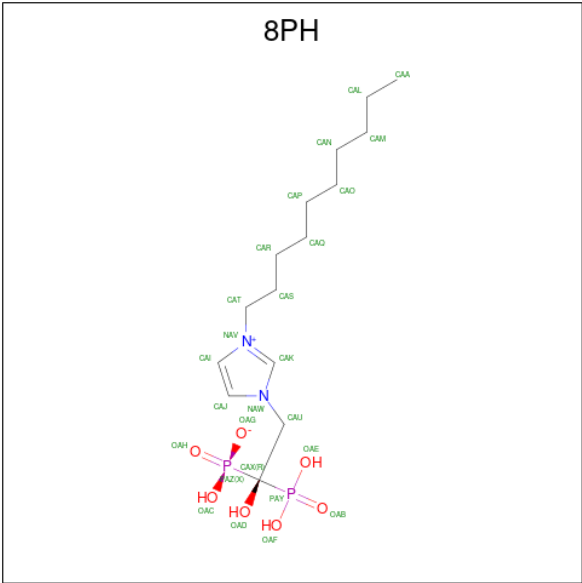
Chain	Residue	Modelled	Actual	Comment	Reference
E	16	HIS	-	expression tag	UNP P37268
E	17	HIS	-	expression tag	UNP P37268
E	18	HIS	-	expression tag	UNP P37268
E	19	HIS	-	expression tag	UNP P37268
E	20	HIS	-	expression tag	UNP P37268
E	21	SER	-	expression tag	UNP P37268
E	22	SER	-	expression tag	UNP P37268
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-decyl-1H-imidazol-3-ium-1-yl)-1-hydroxy-1-phosphonoethyl]phosphonate (CCD ID: 8PH) (formula: C₁₅H₃₀N₂O₇P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 26	C 15	N 2	O 7	P 2	0	0
2	A	1	Total C 8 8					0	0
2	A	1	Total 18	C 7	N 2	O 7	P 2	0	0
2	B	1	Total 26	C 15	N 2	O 7	P 2	0	0
2	B	1	Total C 9 9					0	0
2	C	1	Total 26	C 15	N 2	O 7	P 2	0	0
2	D	1	Total 26	C 15	N 2	O 7	P 2	0	0
2	E	1	Total 26	C 15	N 2	O 7	P 2	0	0
2	F	1	Total 26	C 15	N 2	O 7	P 2	0	0
2	F	1	Total 26	C 15	N 2	O 7	P 2	0	0

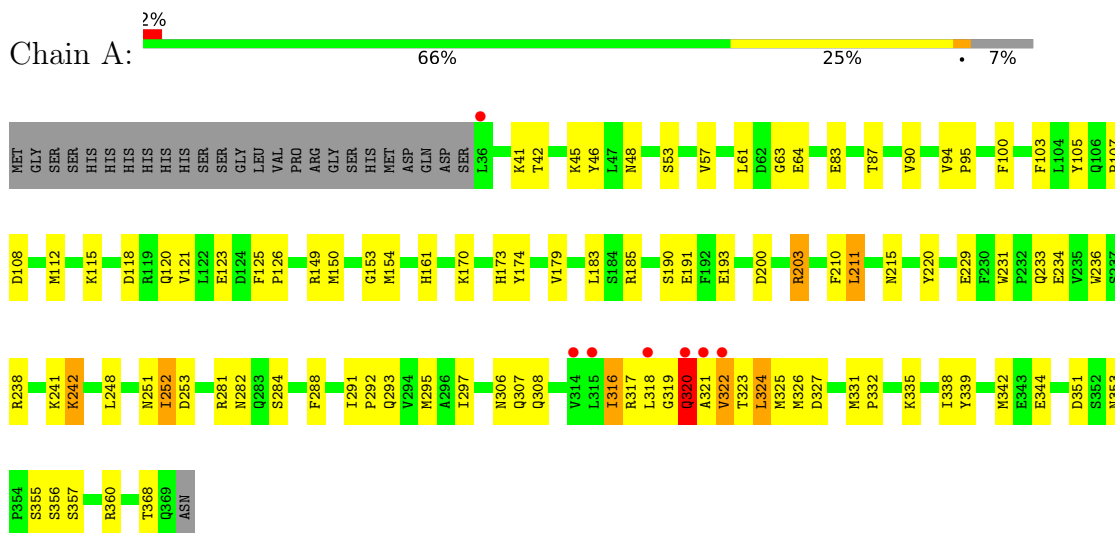
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	99	Total 99	O 99	0	0
3	B	89	Total 89	O 89	0	0
3	C	86	Total 86	O 86	0	0
3	D	62	Total 62	O 62	0	0
3	E	50	Total 50	O 50	0	0
3	F	41	Total 41	O 41	0	0

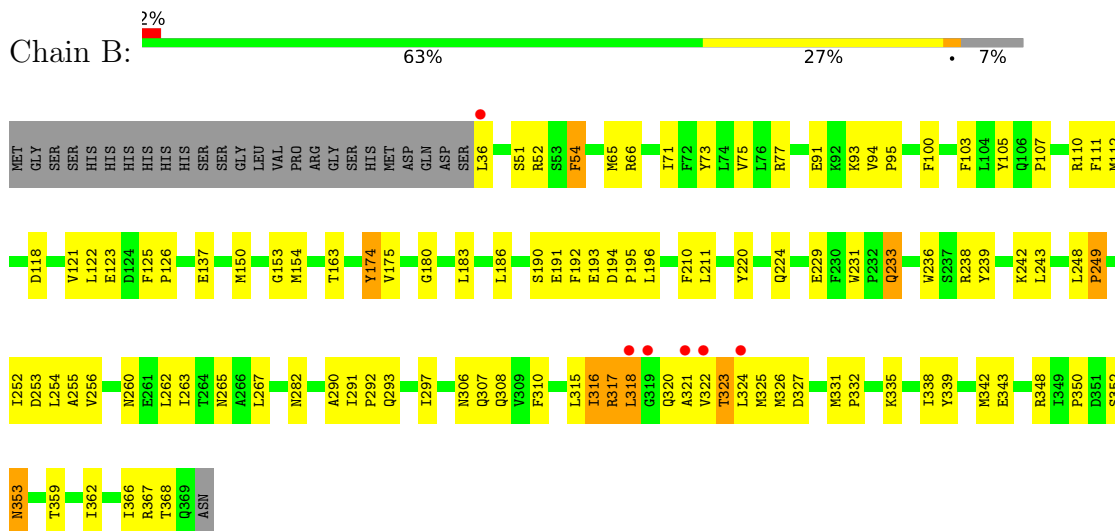
3 Residue-property plots

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

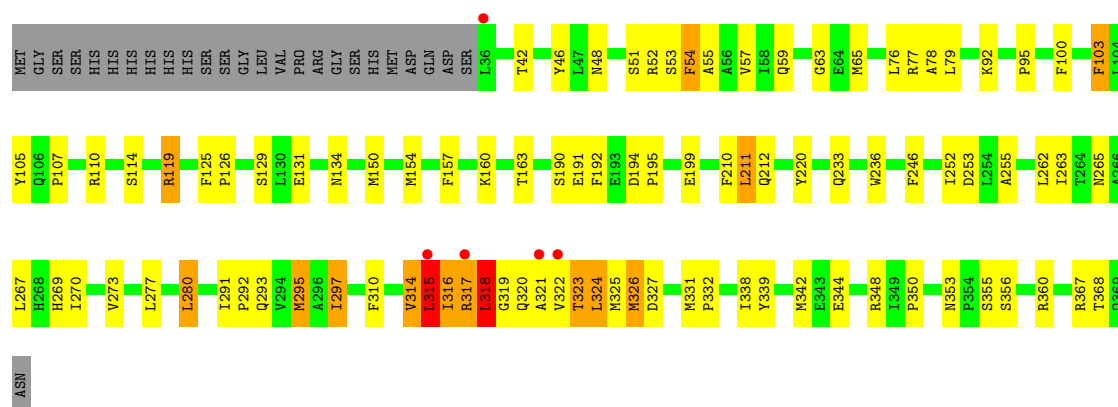


- Molecule 1: Squalene synthase

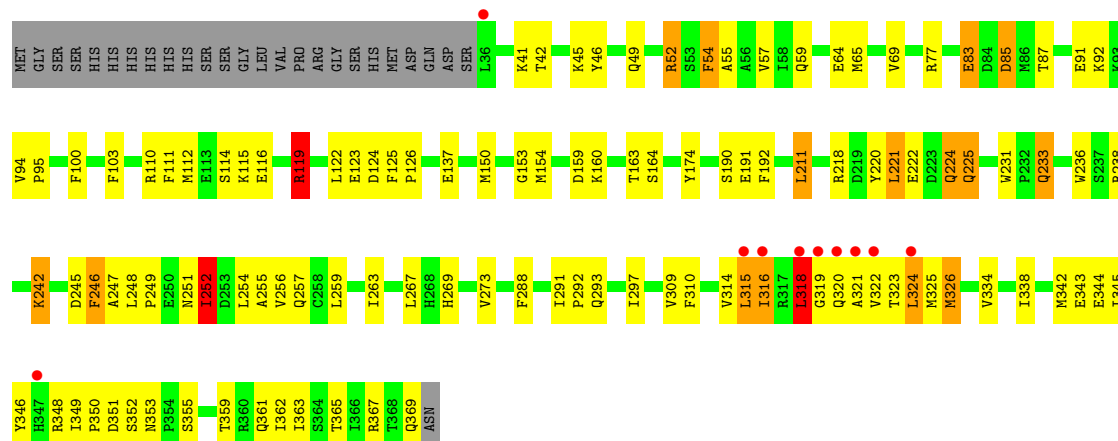


- Molecule 1: Squalene synthase

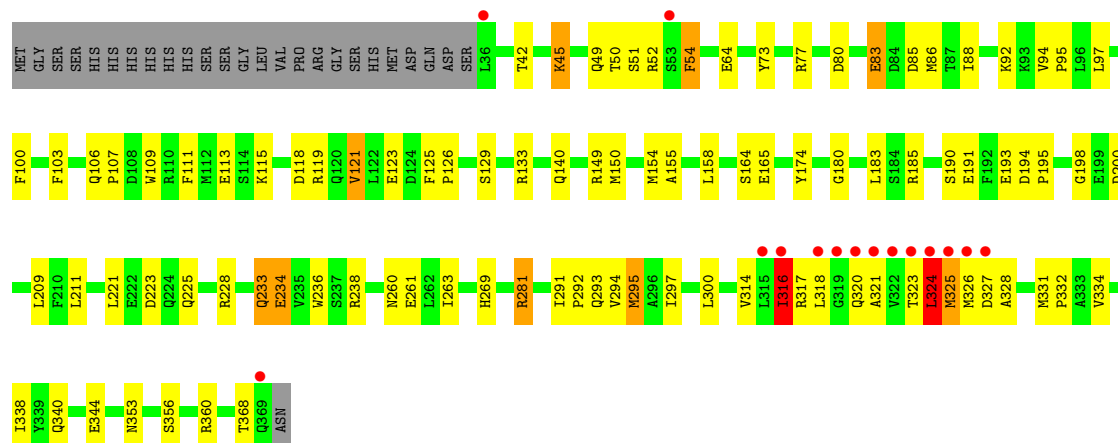




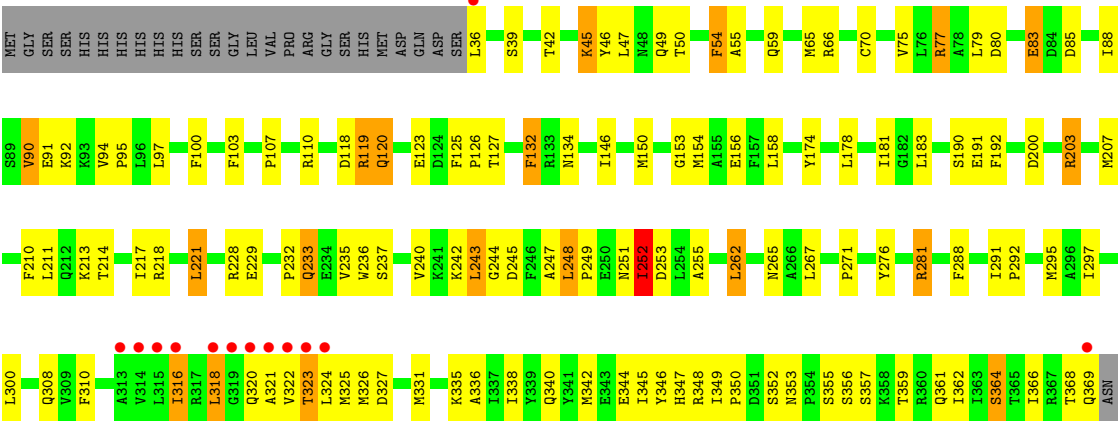
• Molecule 1: Squalene synthase



• 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	85.63Å 153.55Å 90.40Å 90.00° 91.18° 90.00°	Depositor
Resolution (Å)	25.00 – 2.50 25.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	(Not available) (25.00-2.50) 97.0 (25.00-2.50)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.60 (at 2.50Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.192 , 0.250 0.202 , 0.230	Depositor DCC
R_{free} test set	3967 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	36.6	Xtriage
Anisotropy	0.498	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 49.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.032 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	16820	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.88% 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: 8PH

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	0.45	0/2751	0.89	7/3724 (0.2%)
1	B	0.43	0/2751	0.94	13/3724 (0.3%)
1	C	0.43	0/2751	0.94	8/3724 (0.2%)
1	D	0.41	0/2751	0.90	7/3724 (0.2%)
1	E	0.41	0/2751	0.95	14/3724 (0.4%)
1	F	0.38	0/2751	0.90	6/3724 (0.2%)
All	All	0.42	0/16506	0.92	55/22344 (0.2%)

There are no bond length outliers.

All (55) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	323	THR	N-CA-C	-10.76	98.78	114.39
1	B	54	PHE	N-CA-C	9.79	124.62	112.87
1	C	317	ARG	N-CA-C	9.48	123.60	109.95
1	D	54	PHE	N-CA-C	9.48	124.67	113.20
1	C	324	LEU	N-CA-C	-9.25	101.52	113.17
1	A	324	LEU	N-CA-C	-7.53	104.12	113.38
1	E	106	GLN	CA-C-N	7.46	129.16	119.84
1	E	106	GLN	C-N-CA	7.46	129.16	119.84
1	C	314	VAL	N-CA-C	7.38	120.41	108.97
1	E	324	LEU	N-CA-C	-7.22	104.20	113.16
1	B	317	ARG	N-CA-C	7.07	119.97	110.35
1	E	317	ARG	N-CA-C	6.86	119.68	110.35
1	A	53	SER	N-CA-C	6.67	121.41	113.28
1	B	324	LEU	N-CA-C	-6.39	105.48	113.28
1	F	323	THR	N-CA-C	-6.27	105.94	113.97
1	B	186	LEU	N-CA-C	-6.18	104.08	111.69
1	D	64	GLU	N-CA-C	6.14	118.76	111.33
1	D	324	LEU	N-CA-C	-6.10	105.32	112.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	248	LEU	CA-C-N	6.10	127.46	119.84
1	B	248	LEU	C-N-CA	6.10	127.46	119.84
1	B	323	THR	N-CA-C	-6.07	105.41	114.64
1	E	314	VAL	N-CA-C	6.01	116.77	107.99
1	E	316	ILE	N-CA-C	5.96	121.73	109.34
1	C	368	THR	N-CA-C	5.93	120.05	112.34
1	E	45	LYS	N-CA-C	-5.92	104.91	111.36
1	C	315	LEU	N-CA-C	5.90	123.37	110.80
1	E	64	GLU	N-CA-C	5.67	118.24	111.71
1	D	315	LEU	N-CA-C	5.62	118.83	110.52
1	B	353	ASN	CA-C-N	5.57	125.66	119.87
1	B	353	ASN	C-N-CA	5.57	125.66	119.87
1	A	45	LYS	N-CA-C	-5.54	105.16	111.14
1	E	165	GLU	N-CA-C	-5.50	105.44	111.82
1	A	320	GLN	N-CA-C	5.46	122.44	110.80
1	A	368	THR	N-CA-C	5.40	119.03	112.23
1	D	224	GLN	N-CA-C	5.39	117.57	111.11
1	D	316	ILE	N-CA-C	5.38	116.10	107.98
1	F	318	LEU	N-CA-C	5.36	122.22	110.80
1	D	119	ARG	N-CA-C	5.36	119.30	112.34
1	A	87	THR	N-CA-C	-5.35	106.53	113.16
1	F	54	PHE	N-CA-C	5.25	118.79	112.38
1	F	92	LYS	N-CA-C	-5.24	107.02	113.41
1	C	344	GLU	N-CA-C	-5.24	105.57	111.28
1	C	103	PHE	N-CA-C	-5.22	105.71	111.71
1	E	113	GLU	N-CA-C	5.16	118.71	112.93
1	B	348	ARG	N-CA-C	5.16	119.44	113.20
1	E	318	LEU	N-CA-C	5.15	121.77	110.80
1	B	290	ALA	N-CA-C	5.14	116.96	111.36
1	A	64	GLU	N-CA-C	5.12	117.60	111.71
1	F	45	LYS	N-CA-C	-5.11	105.79	111.36
1	B	174	TYR	N-CA-C	5.11	116.85	111.28
1	E	368	THR	N-CA-C	5.08	119.19	112.89
1	F	132	PHE	N-CA-C	-5.04	105.70	111.14
1	C	323	THR	N-CA-C	-5.02	106.31	114.09
1	E	198	GLY	N-CA-C	-5.01	108.45	114.66
1	B	93	LYS	N-CA-C	5.00	116.81	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	70	0
1	B	2696	0	2676	89	0
1	C	2696	0	2676	85	0
1	D	2696	0	2676	94	0
1	E	2696	0	2676	71	0
1	F	2696	0	2676	107	0
2	A	52	0	56	3	0
2	B	35	0	47	7	0
2	C	26	0	30	4	0
2	D	26	0	30	2	0
2	E	26	0	30	4	0
2	F	52	0	60	5	0
3	A	99	0	0	4	0
3	B	89	0	0	2	0
3	C	86	0	0	2	0
3	D	62	0	0	2	0
3	E	50	0	0	3	0
3	F	41	0	0	3	0
All	All	16820	0	16309	493	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 15.

All (493) 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:E:164:SER:HA	1:E:234:GLU:HG3	1.22	1.08
1:E:316:ILE:HD12	1:E:316:ILE:H	1.19	1.03
1:D:320:GLN:HG2	1:D:321:ALA:H	1.28	0.98
1:B:316:ILE:HD12	1:B:316:ILE:H	1.29	0.96
1:A:316:ILE:HD12	1:A:316:ILE:H	1.29	0.94
1:B:233:GLN:HA	1:B:236:TRP:NE1	1.82	0.94
1:F:320:GLN:HG2	1:F:321:ALA:H	1.33	0.93
1:A:183:LEU:HD21	2:A:402:8PH:H30	1.48	0.92
1:B:320:GLN:HG2	1:B:321:ALA:H	1.34	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:235:VAL:HG11	1:F:262:LEU:HD21	1.51	0.91
1:A:320:GLN:HG2	1:A:321:ALA:H	1.34	0.90
1:B:342:MET:HE2	1:B:342:MET:HA	1.53	0.88
1:F:249:PRO:HA	1:F:252:ILE:HD13	1.55	0.88
1:F:316:ILE:HD12	1:F:316:ILE:H	1.36	0.87
1:E:321:ALA:HA	1:E:324:LEU:HB2	1.54	0.87
1:E:320:GLN:HG2	1:E:321:ALA:H	1.40	0.86
1:E:325:MET:O	1:F:327:ASP:HB2	1.74	0.86
1:B:322:VAL:HG21	1:C:53:SER:HB2	1.56	0.86
1:A:327:ASP:HB2	1:C:325:MET:O	1.76	0.85
1:A:326:MET:HE3	1:B:291:ILE:HD11	1.58	0.85
1:D:249:PRO:HA	1:D:252:ILE:HD13	1.59	0.85
1:E:316:ILE:H	1:E:316:ILE:CD1	1.88	0.84
1:C:211:LEU:HD23	2:C:401:8PH:H19	1.60	0.83
1:B:316:ILE:HD12	1:B:316:ILE:N	1.96	0.81
1:A:320:GLN:HG2	1:A:321:ALA:N	1.95	0.80
1:D:322:VAL:HG22	1:E:295:MET:HE1	1.65	0.79
1:E:260:ASN:HD22	1:E:353:ASN:ND2	1.80	0.79
1:B:211:LEU:HD22	2:B:401:8PH:H20	1.63	0.78
1:F:320:GLN:O	1:F:323:THR:HG22	1.84	0.77
1:C:320:GLN:HG2	1:C:321:ALA:H	1.50	0.77
1:F:150:MET:HG3	1:F:174:TYR:O	1.83	0.77
1:D:116:GLU:O	1:D:119:ARG:HD3	1.86	0.76
1:B:118:ASP:O	1:B:121:VAL:HG22	1.85	0.76
1:B:320:GLN:HG2	1:B:321:ALA:N	2.00	0.76
1:B:320:GLN:O	1:B:323:THR:HG22	1.86	0.75
1:B:325:MET:HG3	1:C:291:ILE:HG23	1.69	0.75
1:D:211:LEU:HD23	2:D:401:8PH:H18	1.68	0.74
1:B:325:MET:O	1:C:327:ASP:HB2	1.88	0.74
1:B:326:MET:HE3	1:C:291:ILE:HD11	1.69	0.73
1:D:321:ALA:HA	1:D:324:LEU:HG	1.71	0.73
1:A:325:MET:O	1:B:327:ASP:HB2	1.89	0.73
1:D:353:ASN:ND2	1:D:355:SER:H	1.86	0.73
1:C:316:ILE:HD12	1:C:316:ILE:H	1.54	0.73
1:C:321:ALA:O	1:C:325:MET:HG2	1.88	0.72
1:E:164:SER:CA	1:E:234:GLU:HG3	2.10	0.72
3:A:586:HOH:O	1:C:325:MET:HE1	1.90	0.72
1:E:320:GLN:HG2	1:E:321:ALA:N	2.05	0.72
1:A:211:LEU:HD13	1:A:293:GLN:NE2	2.04	0.71
1:E:115:LYS:HE2	1:E:115:LYS:HA	1.72	0.71
1:A:320:GLN:CG	1:A:321:ALA:H	2.02	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:320:GLN:HG2	1:D:321:ALA:N	2.04	0.71
1:D:319:GLY:HA3	1:D:323:THR:HG21	1.72	0.71
1:C:210:PHE:CE2	1:C:297:ILE:HD13	2.26	0.70
1:F:320:GLN:HG2	1:F:321:ALA:N	2.06	0.70
1:E:228:ARG:HG3	1:E:228:ARG:HH11	1.56	0.70
1:C:320:GLN:O	1:C:324:LEU:HG	1.91	0.70
1:D:150:MET:O	1:D:154:MET:HG3	1.92	0.70
1:A:353:ASN:HD22	1:A:355:SER:H	1.37	0.70
2:E:401:8PH:H6	3:E:522:HOH:O	1.92	0.70
1:A:353:ASN:ND2	1:A:355:SER:H	1.89	0.69
1:B:316:ILE:H	1:B:316:ILE:CD1	2.00	0.69
1:E:297:ILE:HD13	1:E:338:ILE:HG12	1.73	0.68
1:B:183:LEU:HD12	2:B:401:8PH:H26	1.75	0.68
1:B:343:GLU:OE1	1:C:48:ASN:HB3	1.93	0.68
1:F:342:MET:HE2	1:F:342:MET:HA	1.76	0.68
1:B:260:ASN:HD22	1:B:353:ASN:ND2	1.92	0.67
1:F:297:ILE:HD12	1:F:338:ILE:HG12	1.75	0.67
1:E:150:MET:O	1:E:154:MET:HG3	1.95	0.67
1:D:325:MET:O	1:E:327:ASP:HB2	1.95	0.66
1:F:255:ALA:HB1	1:F:310:PHE:CE2	2.30	0.66
1:A:306:ASN:OD1	1:A:308:GLN:HB3	1.96	0.66
1:F:100:PHE:HA	1:F:103:PHE:CD2	2.30	0.66
1:F:233:GLN:HA	1:F:236:TRP:NE1	2.11	0.66
1:F:350:PRO:HG2	1:F:353:ASN:HB2	1.77	0.66
1:F:94:VAL:HB	1:F:95:PRO:HD3	1.79	0.66
1:F:235:VAL:HG11	1:F:262:LEU:CD2	2.24	0.65
1:B:331:MET:HE3	1:B:335:LYS:HG3	1.78	0.65
1:F:65:MET:HE2	1:F:192:PHE:CD2	2.31	0.65
1:A:190:SER:O	1:A:191:GLU:HB2	1.95	0.65
1:C:316:ILE:HD12	1:C:316:ILE:N	2.12	0.65
1:B:73:TYR:HD2	2:B:402:8PH:H28	1.60	0.64
1:E:297:ILE:CD1	1:E:338:ILE:HG12	2.27	0.64
1:E:51:SER:OG	1:E:54:PHE:HB2	1.97	0.64
1:F:353:ASN:ND2	1:F:355:SER:H	1.95	0.64
1:B:233:GLN:HA	1:B:236:TRP:CD1	2.32	0.63
1:D:87:THR:O	1:D:87:THR:HG22	1.97	0.63
1:D:326:MET:HE3	1:E:291:ILE:CD1	2.29	0.63
1:F:150:MET:O	1:F:154:MET:HG3	1.99	0.63
1:A:316:ILE:HD12	1:A:316:ILE:N	2.09	0.62
1:B:297:ILE:HD12	1:B:338:ILE:HG12	1.81	0.62
1:C:320:GLN:HG2	1:C:321:ALA:N	2.15	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:320:GLN:O	1:D:323:THR:HG22	2.00	0.62
1:E:83:GLU:HG3	1:E:154:MET:HB3	1.82	0.62
1:F:54:PHE:CD2	2:F:402:8PH:H17	2.34	0.62
1:D:233:GLN:HA	1:D:236:TRP:NE1	2.15	0.62
1:C:131:GLU:HA	1:C:134:ASN:ND2	2.14	0.61
1:D:85:ASP:OD2	1:D:87:THR:HB	2.00	0.61
1:C:317:ARG:HD2	1:C:318:LEU:HD12	1.82	0.61
1:D:263:ILE:O	1:D:267:LEU:HG	2.01	0.61
1:E:238:ARG:HH22	1:E:261:GLU:CD	2.07	0.61
1:D:211:LEU:HD13	1:D:293:GLN:NE2	2.15	0.61
1:B:331:MET:HB3	1:B:332:PRO:HD3	1.82	0.61
1:D:52:ARG:HG3	3:D:559:HOH:O	2.00	0.61
1:A:200:ASP:OD2	1:A:203:ARG:HB2	2.00	0.60
1:E:325:MET:C	1:F:327:ASP:HB2	2.26	0.60
1:C:211:LEU:CD2	2:C:401:8PH:H19	2.29	0.60
1:B:350:PRO:HG2	1:B:353:ASN:HB2	1.84	0.60
1:B:110:ARG:HH11	1:B:112:MET:HE2	1.66	0.60
1:D:365:THR:O	1:D:369:GLN:HG3	2.02	0.60
1:E:119:ARG:O	1:E:123:GLU:HG3	2.02	0.60
1:F:200:ASP:OD2	1:F:203:ARG:HB2	2.02	0.60
1:C:325:MET:HA	1:C:325:MET:HE2	1.83	0.59
1:A:293:GLN:O	1:A:297:ILE:HG12	2.02	0.59
1:B:368:THR:HG22	1:B:368:THR:O	2.01	0.59
1:E:260:ASN:HD22	1:E:353:ASN:HD21	1.46	0.59
1:C:211:LEU:HD13	1:C:293:GLN:NE2	2.16	0.59
1:F:245:ASP:O	1:F:248:LEU:HG	2.02	0.59
1:E:94:VAL:HB	1:E:95:PRO:HD3	1.84	0.59
1:A:211:LEU:HD13	1:A:293:GLN:HE22	1.67	0.59
1:D:326:MET:HE3	1:E:291:ILE:HD11	1.83	0.59
1:A:325:MET:HA	1:A:325:MET:HE2	1.85	0.58
1:A:233:GLN:HA	1:A:236:TRP:NE1	2.18	0.58
1:D:159:ASP:OD1	1:D:160:LYS:N	2.37	0.58
1:C:353:ASN:ND2	1:C:355:SER:H	2.01	0.58
1:A:319:GLY:O	1:A:320:GLN:O	2.20	0.58
1:F:119:ARG:O	1:F:123:GLU:HG3	2.02	0.58
1:C:100:PHE:HA	1:C:103:PHE:CD2	2.38	0.58
1:C:212:GLN:OE1	2:C:401:8PH:H12	2.04	0.58
1:D:291:ILE:HB	1:D:292:PRO:HD3	1.85	0.58
1:D:322:VAL:O	1:D:322:VAL:HG12	2.04	0.58
1:F:183:LEU:CD1	2:F:401:8PH:H26	2.33	0.58
1:A:220:TYR:HB2	1:A:231:TRP:CZ2	2.38	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:343:GLU:OE1	1:B:367:ARG:NH2	2.37	0.57
1:D:94:VAL:HB	1:D:95:PRO:HD3	1.86	0.57
1:E:180:GLY:HA2	2:E:401:8PH:H27	1.86	0.57
1:A:83:GLU:OE1	1:A:154:MET:HE2	2.03	0.57
1:A:215:ASN:HD21	2:A:403:8PH:CAS	2.16	0.57
1:C:348:ARG:O	1:C:350:PRO:HD3	2.05	0.57
1:D:242:LYS:HB2	1:D:242:LYS:NZ	2.19	0.57
1:B:66:ARG:HD2	3:B:565:HOH:O	2.03	0.57
1:D:255:ALA:HB1	1:D:310:PHE:CE1	2.38	0.57
1:F:297:ILE:CD1	1:F:338:ILE:HG12	2.34	0.57
1:E:325:MET:HA	1:E:325:MET:HE3	1.86	0.57
1:D:211:LEU:HD13	1:D:293:GLN:HE22	1.70	0.57
1:D:291:ILE:HD11	1:F:326:MET:HE3	1.87	0.57
1:A:291:ILE:HB	1:A:292:PRO:HD3	1.87	0.56
1:A:291:ILE:HD11	1:C:326:MET:CE	2.35	0.56
1:B:112:MET:SD	1:B:123:GLU:HG2	2.45	0.56
1:D:346:TYR:HA	1:D:349:ILE:CD1	2.35	0.56
1:F:79:LEU:HD12	1:F:79:LEU:O	2.05	0.56
1:D:325:MET:HE3	1:D:325:MET:HA	1.86	0.56
1:B:320:GLN:CG	1:B:321:ALA:H	2.15	0.56
1:D:159:ASP:CG	1:D:160:LYS:N	2.64	0.56
1:F:120:GLN:HA	1:F:120:GLN:HE21	1.70	0.56
1:E:97:LEU:HD12	1:E:158:LEU:HD11	1.88	0.56
1:B:190:SER:O	1:B:191:GLU:HB2	2.05	0.56
1:D:220:TYR:HB2	1:D:231:TRP:CZ2	2.42	0.55
1:E:316:ILE:HD12	1:E:316:ILE:N	2.04	0.55
1:A:331:MET:HB3	1:A:332:PRO:HD3	1.87	0.55
1:F:324:LEU:O	1:F:325:MET:HE2	2.06	0.55
1:C:342:MET:HE2	1:C:342:MET:HA	1.88	0.55
1:E:263:ILE:HG12	1:E:300:LEU:HD22	1.87	0.55
1:A:61:LEU:HA	1:A:284:SER:HB2	1.88	0.55
1:B:94:VAL:HB	1:B:95:PRO:HD3	1.87	0.55
1:F:291:ILE:HB	1:F:292:PRO:HD3	1.87	0.55
1:C:211:LEU:HD13	1:C:293:GLN:HE22	1.72	0.55
1:C:277:LEU:HA	1:C:280:LEU:HD22	1.87	0.55
1:D:163:THR:OG1	1:D:164:SER:N	2.36	0.55
1:A:327:ASP:HB2	1:C:325:MET:C	2.31	0.55
1:C:338:ILE:O	1:C:342:MET:HG2	2.07	0.55
1:B:211:LEU:CD1	1:B:293:GLN:NE2	2.70	0.55
1:A:316:ILE:H	1:A:316:ILE:CD1	2.09	0.55
1:E:77:ARG:O	1:E:77:ARG:HD3	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:255:ALA:HB1	1:F:310:PHE:CZ	2.41	0.54
1:A:112:MET:SD	1:A:123:GLU:HG2	2.47	0.54
1:A:210:PHE:CZ	1:A:297:ILE:HD13	2.43	0.54
1:B:91:GLU:CD	1:B:91:GLU:H	2.15	0.54
1:A:94:VAL:HB	1:A:95:PRO:HD3	1.89	0.54
1:B:362:ILE:O	1:B:366:ILE:HG13	2.07	0.54
1:E:320:GLN:CG	1:E:321:ALA:H	2.18	0.54
1:D:343:GLU:CD	1:D:367:ARG:HH21	2.15	0.54
1:B:297:ILE:HD13	1:B:338:ILE:HG23	1.88	0.54
1:E:209:LEU:HD23	1:E:269:HIS:CE1	2.43	0.54
1:F:83:GLU:OE2	1:F:154:MET:HE2	2.08	0.54
1:F:233:GLN:HA	1:F:236:TRP:CD1	2.43	0.54
1:C:321:ALA:HA	1:C:324:LEU:HG	1.90	0.54
1:B:36:LEU:HB2	1:D:137:GLU:OE1	2.08	0.54
1:B:338:ILE:O	1:B:342:MET:HG2	2.08	0.54
1:C:263:ILE:O	1:C:267:LEU:HG	2.08	0.54
1:E:85:ASP:HB3	1:E:88:ILE:HD12	1.89	0.54
1:A:48:ASN:ND2	1:C:367:ARG:HD3	2.24	0.53
1:F:281:ARG:CG	1:F:281:ARG:HH21	2.21	0.53
1:F:320:GLN:CG	1:F:321:ALA:H	2.14	0.53
1:B:229:GLU:HG2	1:B:243:LEU:HD23	1.91	0.53
1:A:108:ASP:HB2	3:A:551:HOH:O	2.08	0.53
1:E:331:MET:HB3	1:E:332:PRO:HD3	1.90	0.53
1:B:252:ILE:CG2	1:B:253:ASP:N	2.71	0.53
1:F:336:ALA:O	1:F:340:GLN:HG2	2.09	0.53
1:D:65:MET:HE2	1:D:192:PHE:CD2	2.44	0.53
1:F:207:MET:HE2	1:F:276:TYR:CD1	2.44	0.53
1:F:349:ILE:HD13	1:F:359:THR:HB	1.92	0.52
1:B:239:TYR:O	1:B:254:LEU:HD13	2.10	0.52
1:B:317:ARG:HG3	1:B:318:LEU:N	2.22	0.52
1:A:291:ILE:HD11	1:C:326:MET:HE3	1.90	0.52
1:C:105:TYR:O	1:C:107:PRO:HD3	2.09	0.52
1:D:245:ASP:O	1:D:247:ALA:N	2.43	0.52
1:E:211:LEU:HD13	1:E:293:GLN:NE2	2.25	0.52
1:F:211:LEU:HD23	2:F:401:8PH:H18	1.90	0.52
1:F:353:ASN:HD22	1:F:355:SER:H	1.57	0.52
1:C:220:TYR:CZ	1:C:246:PHE:HB2	2.43	0.52
1:F:362:ILE:O	1:F:366:ILE:HG13	2.10	0.52
1:F:45:LYS:HG2	1:F:49:GLN:HE21	1.73	0.52
1:A:57:VAL:HG11	1:A:288:PHE:HA	1.91	0.51
1:B:175:VAL:O	2:B:401:8PH:H10	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:344:GLU:O	1:F:348:ARG:HD3	2.10	0.51
1:A:150:MET:HG3	1:A:174:TYR:O	2.10	0.51
1:D:65:MET:O	1:D:69:VAL:HG23	2.10	0.51
1:E:321:ALA:CA	1:E:324:LEU:HB2	2.34	0.51
1:C:350:PRO:HG2	1:C:353:ASN:HB2	1.92	0.51
1:D:55:ALA:O	1:D:59:GLN:HG3	2.10	0.51
1:E:261:GLU:HG3	3:E:543:HOH:O	2.10	0.51
1:F:39:SER:HB2	1:F:127:THR:HG23	1.93	0.51
1:F:55:ALA:O	1:F:59:GLN:HG3	2.10	0.51
1:E:51:SER:HB2	1:E:73:TYR:CZ	2.45	0.51
1:F:271:PRO:HB3	1:F:369:GLN:HE22	1.76	0.51
1:A:321:ALA:HA	1:A:324:LEU:CD1	2.41	0.51
1:C:255:ALA:HB1	1:C:310:PHE:CZ	2.46	0.51
1:F:77:ARG:HE	1:F:77:ARG:HA	1.75	0.51
1:B:193:GLU:OE2	1:B:282:ASN:HB3	2.11	0.51
1:D:45:LYS:O	1:D:49:GLN:HG3	2.10	0.51
1:B:211:LEU:HD13	1:B:293:GLN:HE22	1.76	0.51
1:F:353:ASN:ND2	1:F:355:SER:HB2	2.26	0.51
1:D:91:GLU:HG3	1:D:92:LYS:N	2.25	0.51
1:D:225:GLN:H	1:D:225:GLN:HE21	1.57	0.51
1:A:321:ALA:HA	1:A:324:LEU:HG	1.93	0.50
1:C:131:GLU:HA	1:C:134:ASN:HD22	1.73	0.50
1:B:233:GLN:HA	1:B:236:TRP:HE1	1.72	0.50
1:C:291:ILE:HB	1:C:292:PRO:HD3	1.91	0.50
1:C:339:TYR:O	1:C:342:MET:HB2	2.11	0.50
1:D:163:THR:HA	1:D:233:GLN:HB3	1.93	0.50
1:F:357:SER:O	1:F:361:GLN:HG3	2.11	0.50
1:D:85:ASP:OD1	1:D:114:SER:HA	2.12	0.50
1:E:356:SER:O	1:E:360:ARG:HG3	2.12	0.50
1:B:71:ILE:O	1:B:75:VAL:HG13	2.11	0.50
1:B:322:VAL:HG12	1:B:322:VAL:O	2.11	0.50
1:C:293:GLN:O	1:C:297:ILE:HG12	2.11	0.50
1:E:193:GLU:HA	1:E:281:ARG:HD3	1.93	0.50
1:B:291:ILE:HB	1:B:292:PRO:HD3	1.94	0.50
1:D:259:LEU:HD21	1:D:309:VAL:HG21	1.92	0.50
1:E:183:LEU:HD12	2:E:401:8PH:H26	1.92	0.50
1:E:118:ASP:O	1:E:121:VAL:HG13	2.12	0.50
1:E:149:ARG:NH2	1:E:185:ARG:HH12	2.10	0.50
1:B:256:VAL:HG21	1:B:307:GLN:HG2	1.94	0.49
1:B:322:VAL:CG2	1:C:53:SER:HB2	2.33	0.49
1:E:321:ALA:HA	1:E:324:LEU:HD12	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:281:ARG:HD2	3:F:523:HOH:O	2.12	0.49
1:F:348:ARG:O	1:F:350:PRO:HD3	2.12	0.49
1:D:87:THR:O	1:D:87:THR:CG2	2.60	0.49
1:D:125:PHE:N	1:D:126:PRO:CD	2.76	0.49
1:A:149:ARG:NH2	1:A:185:ARG:HH22	2.11	0.49
1:C:314:VAL:CG1	1:C:315:LEU:N	2.74	0.49
1:F:322:VAL:HG12	1:F:322:VAL:O	2.12	0.49
1:A:100:PHE:HA	1:A:103:PHE:CD2	2.47	0.49
1:B:51:SER:O	1:B:52:ARG:C	2.55	0.49
1:F:190:SER:O	1:F:191:GLU:HB2	2.13	0.49
1:C:157:PHE:CE2	1:C:160:LYS:HE3	2.48	0.49
1:C:320:GLN:O	1:C:323:THR:HG22	2.13	0.49
1:D:41:LYS:HD3	1:F:364:SER:OG	2.13	0.49
1:D:254:LEU:HA	1:D:257:GLN:NE2	2.27	0.49
1:A:125:PHE:N	1:A:126:PRO:CD	2.76	0.49
1:A:210:PHE:CE2	1:A:297:ILE:HD13	2.48	0.49
1:B:342:MET:HA	1:B:342:MET:CE	2.35	0.49
1:C:195:PRO:O	1:C:199:GLU:HG3	2.12	0.49
1:C:317:ARG:HD2	1:C:318:LEU:CD1	2.42	0.49
1:D:110:ARG:HD2	1:D:112:MET:SD	2.53	0.49
1:F:240:VAL:HG21	1:F:245:ASP:HB2	1.95	0.49
1:F:331:MET:O	1:F:335:LYS:HG3	2.13	0.49
1:F:248:LEU:HD12	1:F:251:ASN:HD21	1.78	0.48
1:D:246:PHE:CD1	1:D:255:ALA:HA	2.48	0.48
1:F:232:PRO:HD2	1:F:262:LEU:HD11	1.95	0.48
1:D:153:GLY:HA3	1:D:174:TYR:CD1	2.48	0.48
1:D:345:ILE:O	1:D:349:ILE:HG13	2.12	0.48
1:A:118:ASP:O	1:A:121:VAL:HG22	2.14	0.48
1:A:63:GLY:HA3	3:A:510:HOH:O	2.12	0.48
1:A:327:ASP:OD2	1:B:327:ASP:OD2	2.32	0.48
1:B:150:MET:HG3	1:B:174:TYR:O	2.12	0.48
1:F:368:THR:O	1:F:368:THR:HG22	2.12	0.48
1:A:90:VAL:O	1:A:94:VAL:HG23	2.14	0.48
1:A:356:SER:O	1:A:360:ARG:HG3	2.13	0.48
1:D:211:LEU:CD2	2:D:401:8PH:H18	2.42	0.48
1:F:221:LEU:HD13	1:F:310:PHE:HA	1.96	0.48
1:C:92:LYS:O	1:C:95:PRO:HG2	2.14	0.48
1:C:114:SER:HB3	1:C:119:ARG:HB2	1.95	0.48
1:E:223:ASP:OD1	1:E:228:ARG:NH2	2.47	0.48
1:E:228:ARG:HG3	1:E:228:ARG:NH1	2.26	0.48
1:C:150:MET:O	1:C:154:MET:HG3	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:114:SER:O	1:D:115:LYS:HD2	2.14	0.47
1:D:163:THR:HA	1:D:233:GLN:HE21	1.77	0.47
1:D:238:ARG:HB2	1:D:238:ARG:NH1	2.28	0.47
1:A:320:GLN:CG	1:A:321:ALA:N	2.63	0.47
1:A:248:LEU:HB2	1:A:251:ASN:ND2	2.29	0.47
1:B:297:ILE:CD1	1:B:338:ILE:HG23	2.43	0.47
1:C:125:PHE:N	1:C:126:PRO:CD	2.78	0.47
1:D:255:ALA:HB1	1:D:310:PHE:CZ	2.50	0.47
1:C:42:THR:HG22	1:C:46:TYR:CE2	2.50	0.47
1:D:353:ASN:HD22	1:D:355:SER:H	1.61	0.47
1:E:194:ASP:OD1	1:E:195:PRO:HD2	2.15	0.47
1:F:36:LEU:N	3:F:533:HOH:O	2.48	0.47
1:F:210:PHE:CZ	1:F:297:ILE:HG12	2.50	0.47
1:C:65:MET:HE2	1:C:192:PHE:CD2	2.49	0.47
1:D:159:ASP:CG	1:D:160:LYS:H	2.22	0.47
1:A:153:GLY:HA3	1:A:174:TYR:CG	2.50	0.46
1:C:210:PHE:HE2	1:C:297:ILE:HD13	1.78	0.46
1:C:270:ILE:HD13	1:C:342:MET:SD	2.55	0.46
1:F:146:ILE:HG22	1:F:150:MET:CE	2.44	0.46
1:A:326:MET:HE3	1:B:291:ILE:CD1	2.39	0.46
1:C:322:VAL:HG12	1:C:322:VAL:O	2.15	0.46
1:D:325:MET:HE2	1:E:294:VAL:HG11	1.96	0.46
1:E:126:PRO:HA	1:E:129:SER:OG	2.15	0.46
1:F:47:LEU:HD22	1:F:70:CYS:SG	2.55	0.46
1:B:100:PHE:HA	1:B:103:PHE:CD2	2.50	0.46
1:B:306:ASN:OD1	1:B:306:ASN:C	2.57	0.46
1:E:324:LEU:HD13	3:E:517:HOH:O	2.14	0.46
1:C:320:GLN:CG	1:C:321:ALA:H	2.25	0.46
1:A:161:HIS:HD2	1:A:229:GLU:HB2	1.80	0.46
1:B:262:LEU:O	1:B:265:ASN:HB3	2.15	0.46
1:B:339:TYR:O	1:B:342:MET:HB2	2.16	0.46
1:D:83:GLU:HG3	1:D:154:MET:HB3	1.98	0.46
1:B:252:ILE:HG21	1:B:307:GLN:NE2	2.31	0.46
1:F:65:MET:HE2	1:F:192:PHE:CG	2.51	0.46
1:A:316:ILE:N	1:A:316:ILE:CD1	2.76	0.46
1:E:321:ALA:HA	1:E:324:LEU:CB	2.37	0.46
1:F:178:LEU:HD23	1:F:181:ILE:HD12	1.98	0.46
1:B:125:PHE:N	1:B:126:PRO:CD	2.79	0.46
1:C:190:SER:O	1:C:191:GLU:HB2	2.14	0.46
1:D:318:LEU:H	1:D:318:LEU:HG	1.51	0.46
1:D:359:THR:O	1:D:362:ILE:HG22	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:218:ARG:O	1:F:218:ARG:HG2	2.16	0.46
1:A:338:ILE:O	1:A:342:MET:HG2	2.16	0.46
1:B:105:TYR:O	1:B:107:PRO:HD3	2.16	0.46
1:C:317:ARG:NH1	1:C:348:ARG:NH2	2.64	0.46
1:B:65:MET:HE2	1:B:192:PHE:CD2	2.51	0.46
1:D:297:ILE:CD1	1:D:338:ILE:HG12	2.46	0.46
1:A:242:LYS:NZ	1:A:242:LYS:HB2	2.30	0.45
1:B:52:ARG:HG3	3:B:572:HOH:O	2.16	0.45
1:F:66:ARG:HB3	3:F:532:HOH:O	2.16	0.45
1:F:85:ASP:HB3	1:F:88:ILE:HD12	1.98	0.45
1:A:335:LYS:HB3	1:A:339:TYR:CE2	2.52	0.45
1:D:346:TYR:HA	1:D:349:ILE:HD11	1.97	0.45
1:B:153:GLY:HA3	1:B:174:TYR:CG	2.52	0.45
1:C:220:TYR:OH	1:C:246:PHE:HB2	2.16	0.45
1:D:348:ARG:O	1:D:350:PRO:HD3	2.17	0.45
1:F:243:LEU:HD12	1:F:243:LEU:O	2.16	0.45
1:C:316:ILE:H	1:C:316:ILE:CD1	2.13	0.45
1:F:153:GLY:HA3	1:F:174:TYR:CD1	2.51	0.45
1:F:214:THR:HG23	1:F:300:LEU:CD2	2.46	0.45
1:A:251:ASN:O	1:A:252:ILE:C	2.60	0.45
1:B:183:LEU:CD1	2:B:401:8PH:H26	2.46	0.45
1:D:248:LEU:HB2	1:D:251:ASN:HD22	1.82	0.45
1:D:362:ILE:HG23	1:D:363:ILE:N	2.32	0.45
1:E:45:LYS:HG2	1:E:49:GLN:NE2	2.32	0.45
1:F:42:THR:HG22	1:F:46:TYR:CE2	2.51	0.45
1:F:125:PHE:N	1:F:126:PRO:CD	2.79	0.45
1:A:252:ILE:HD11	1:A:307:GLN:HB3	1.99	0.44
1:C:321:ALA:O	1:C:325:MET:CG	2.60	0.44
1:D:238:ARG:HB2	1:D:238:ARG:HH11	1.81	0.44
1:F:233:GLN:C	1:F:235:VAL:N	2.74	0.44
1:F:245:ASP:C	1:F:247:ALA:H	2.26	0.44
1:B:325:MET:HE2	1:C:295:MET:HG3	1.99	0.44
1:D:245:ASP:C	1:D:247:ALA:N	2.75	0.44
1:F:233:GLN:C	1:F:235:VAL:H	2.25	0.44
1:F:233:GLN:O	1:F:237:SER:HB3	2.18	0.44
1:F:248:LEU:HA	1:F:249:PRO:HD3	1.88	0.44
1:A:322:VAL:O	1:A:322:VAL:CG1	2.64	0.44
1:C:331:MET:HB3	1:C:332:PRO:HD3	2.00	0.44
1:D:100:PHE:HA	1:D:103:PHE:CD2	2.53	0.44
1:E:155:ALA:HA	1:E:158:LEU:CD1	2.48	0.44
1:C:54:PHE:O	1:C:57:VAL:HG22	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:321:ALA:O	1:D:325:MET:HG2	2.17	0.44
2:C:401:8PH:H27	3:C:543:HOH:O	2.18	0.44
1:D:42:THR:HG22	1:D:46:TYR:CE2	2.53	0.44
1:B:211:LEU:HD11	1:B:293:GLN:NE2	2.32	0.44
1:F:156:GLU:HG2	1:F:156:GLU:O	2.18	0.44
1:C:318:LEU:HB2	1:C:319:GLY:H	1.47	0.44
1:E:50:THR:OG1	1:E:73:TYR:HE1	2.01	0.44
1:B:267:LEU:HD22	1:B:342:MET:HE1	1.99	0.43
1:C:315:LEU:O	1:C:316:ILE:C	2.59	0.43
1:F:267:LEU:HD22	1:F:342:MET:HE1	2.00	0.43
1:B:252:ILE:HG23	1:B:307:GLN:HE21	1.83	0.43
1:D:221:LEU:O	1:D:222:GLU:C	2.60	0.43
1:D:269:HIS:O	1:D:273:VAL:HG23	2.18	0.43
1:E:100:PHE:HA	1:E:103:PHE:CD2	2.53	0.43
1:F:50:THR:HG22	1:F:118:ASP:OD1	2.18	0.43
1:F:228:ARG:HG3	1:F:228:ARG:HH11	1.82	0.43
1:B:211:LEU:CD1	1:B:293:GLN:HE22	2.31	0.43
1:E:92:LYS:O	1:E:95:PRO:HD2	2.18	0.43
1:A:193:GLU:OE2	1:A:282:ASN:HB3	2.18	0.43
1:D:150:MET:HG3	1:D:174:TYR:O	2.19	0.43
1:A:291:ILE:HG21	1:C:322:VAL:HG13	2.01	0.43
1:C:51:SER:HB3	1:C:55:ALA:HB2	2.00	0.43
1:C:163:THR:HA	1:C:233:GLN:HB3	2.00	0.43
1:E:125:PHE:N	1:E:126:PRO:CD	2.81	0.43
1:E:133:ARG:HA	1:E:140:GLN:NE2	2.33	0.43
1:E:325:MET:C	1:E:326:MET:HG3	2.44	0.43
1:F:91:GLU:H	1:F:91:GLU:CD	2.25	0.43
1:C:324:LEU:HD12	3:C:535:HOH:O	2.18	0.43
1:D:334:VAL:O	1:D:338:ILE:HG13	2.19	0.43
1:F:213:LYS:O	1:F:217:ILE:HG13	2.19	0.43
1:F:355:SER:O	1:F:356:SER:C	2.62	0.43
1:B:150:MET:O	1:B:154:MET:HG3	2.18	0.43
1:B:359:THR:O	1:B:362:ILE:HG22	2.18	0.43
1:D:190:SER:O	1:D:191:GLU:HB2	2.19	0.43
1:B:236:TRP:CZ3	1:B:242:LYS:HA	2.54	0.43
1:C:317:ARG:NH1	1:C:348:ARG:HH21	2.17	0.43
1:D:153:GLY:HA3	1:D:174:TYR:CG	2.53	0.43
1:B:111:PHE:HB3	1:B:122:LEU:HB3	2.01	0.43
1:B:194:ASP:OD1	1:B:195:PRO:HD2	2.18	0.43
1:F:262:LEU:O	1:F:265:ASN:HB3	2.19	0.43
1:A:238:ARG:HB2	1:A:238:ARG:HH11	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:221:LEU:O	1:D:224:GLN:N	2.52	0.43
2:E:401:8PH:H10	2:E:401:8PH:H17	2.01	0.43
1:F:338:ILE:O	1:F:342:MET:HG2	2.19	0.43
1:A:252:ILE:HG23	1:A:253:ASP:N	2.33	0.42
1:E:150:MET:HG3	1:E:174:TYR:O	2.19	0.42
1:E:109:TRP:CH2	1:E:111:PHE:HB2	2.54	0.42
1:F:79:LEU:HB2	1:F:100:PHE:CE2	2.55	0.42
1:F:281:ARG:CG	1:F:281:ARG:NH2	2.81	0.42
1:A:238:ARG:HB2	1:A:238:ARG:NH1	2.35	0.42
1:B:263:ILE:O	1:B:267:LEU:HG	2.20	0.42
1:B:220:TYR:CZ	1:B:224:GLN:HG3	2.54	0.42
1:F:233:GLN:HE21	1:F:233:GLN:HB3	1.51	0.42
3:A:578:HOH:O	1:B:66:ARG:HD3	2.19	0.42
1:D:344:GLU:HG2	3:D:540:HOH:O	2.19	0.42
1:F:45:LYS:HG2	1:F:49:GLN:NE2	2.35	0.42
1:F:90:VAL:O	1:F:94:VAL:HG23	2.18	0.42
1:F:345:ILE:O	1:F:346:TYR:C	2.62	0.42
1:D:342:MET:HE2	1:D:342:MET:HA	2.01	0.42
1:E:150:MET:HE3	1:E:150:MET:HB2	1.90	0.42
1:F:146:ILE:HG22	1:F:150:MET:HE3	2.00	0.42
1:A:170:LYS:O	1:A:173:HIS:HB3	2.20	0.42
1:A:353:ASN:ND2	1:A:355:SER:HB2	2.35	0.42
1:B:210:PHE:CZ	1:B:297:ILE:HG12	2.55	0.42
1:C:356:SER:O	1:C:360:ARG:HG3	2.20	0.42
1:E:190:SER:O	1:E:191:GLU:HB2	2.20	0.42
1:F:54:PHE:HE2	1:F:288:PHE:CE1	2.38	0.42
1:B:252:ILE:CG2	1:B:307:GLN:NE2	2.82	0.42
1:C:76:LEU:HD22	1:C:150:MET:SD	2.60	0.42
1:C:355:SER:O	1:C:356:SER:C	2.63	0.42
1:D:116:GLU:O	1:D:119:ARG:CD	2.63	0.42
1:D:220:TYR:OH	1:D:246:PHE:HB2	2.20	0.42
1:E:334:VAL:O	1:E:338:ILE:HG13	2.20	0.42
1:B:306:ASN:OD1	1:B:308:GLN:N	2.50	0.41
1:C:269:HIS:O	1:C:273:VAL:HG23	2.20	0.41
1:F:242:LYS:O	1:F:244:GLY:N	2.53	0.41
1:A:179:VAL:HB	2:A:401:8PH:H19	2.02	0.41
1:E:211:LEU:HD13	1:E:293:GLN:HE22	1.85	0.41
2:B:401:8PH:H21	2:B:402:8PH:H22	2.03	0.41
1:F:100:PHE:HA	1:F:103:PHE:CE2	2.55	0.41
1:F:229:GLU:HB3	1:F:243:LEU:HD23	2.03	0.41
1:C:194:ASP:OD2	1:C:195:PRO:HD2	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:233:GLN:HA	1:E:236:TRP:CE2	2.56	0.41
1:E:291:ILE:HB	1:E:292:PRO:HD3	2.02	0.41
1:F:80:ASP:OD1	2:F:401:8PH:H7	2.19	0.41
1:B:255:ALA:HB1	1:B:310:PHE:CZ	2.56	0.41
1:C:252:ILE:HG23	1:C:253:ASP:N	2.36	0.41
1:D:57:VAL:HG11	1:D:288:PHE:HA	2.03	0.41
1:D:248:LEU:HA	1:D:249:PRO:HD2	1.90	0.41
1:D:349:ILE:HA	1:D:350:PRO:HD3	1.91	0.41
1:E:73:TYR:CD1	1:E:73:TYR:C	2.99	0.41
1:F:153:GLY:HA3	1:F:174:TYR:CG	2.56	0.41
1:F:251:ASN:O	1:F:252:ILE:C	2.63	0.41
1:C:77:ARG:HA	1:C:77:ARG:HD2	1.92	0.41
1:E:149:ARG:HH21	1:E:185:ARG:HH12	1.69	0.41
1:E:221:LEU:HG	1:E:225:GLN:OE1	2.21	0.41
1:B:220:TYR:HB2	1:B:231:TRP:CZ2	2.56	0.40
1:C:51:SER:OG	1:C:54:PHE:HB2	2.22	0.40
1:C:78:ALA:O	1:C:79:LEU:C	2.64	0.40
1:D:124:ASP:C	1:D:126:PRO:HD2	2.46	0.40
1:D:218:ARG:O	1:D:218:ARG:HG2	2.20	0.40
1:F:183:LEU:HD13	2:F:401:8PH:H26	2.02	0.40
1:C:233:GLN:HA	1:C:236:TRP:NE1	2.37	0.40
1:D:77:ARG:HA	1:D:77:ARG:HD3	1.97	0.40
1:D:111:PHE:HB3	1:D:122:LEU:HB3	2.03	0.40
1:F:252:ILE:HG22	1:F:253:ASP:OD2	2.21	0.40
1:A:105:TYR:O	1:A:107:PRO:HD3	2.21	0.40
1:B:163:THR:HA	1:B:233:GLN:HB3	2.04	0.40
1:D:163:THR:CA	1:D:233:GLN:HE21	2.35	0.40
1:B:180:GLY:HA2	2:B:401:8PH:H27	2.03	0.40
1:C:262:LEU:O	1:C:265:ASN:HB3	2.21	0.40
1:F:97:LEU:HD12	1:F:158:LEU:HD11	2.04	0.40
1:A:42:THR:HG22	1:A:46:TYR:CE2	2.57	0.40
1:C:353:ASN:HD22	1:C:355:SER:H	1.70	0.40
1:F:132:PHE:C	1:F:134:ASN:H	2.29	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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%)	314 (95%)	15 (4%)	3 (1%)	14	27
1	B	332/360 (92%)	312 (94%)	18 (5%)	2 (1%)	21	38
1	C	332/360 (92%)	317 (96%)	12 (4%)	3 (1%)	14	27
1	D	332/360 (92%)	311 (94%)	17 (5%)	4 (1%)	10	20
1	E	332/360 (92%)	318 (96%)	10 (3%)	4 (1%)	10	20
1	F	332/360 (92%)	305 (92%)	21 (6%)	6 (2%)	6	12
All	All	1992/2160 (92%)	1877 (94%)	93 (5%)	22 (1%)	11	22

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	318	LEU
1	A	320	GLN
1	B	318	LEU
1	C	318	LEU
1	D	318	LEU
1	F	318	LEU
1	D	85	ASP
1	D	246	PHE
1	E	200	ASP
1	F	243	LEU
1	E	54	PHE
1	E	107	PRO
1	E	328	ALA
1	F	252	ILE
1	A	252	ILE
1	B	249	PRO
1	D	252	ILE
1	F	308	GLN
1	C	54	PHE
1	F	119	ARG

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Mol	Chain	Res	Type
1	F	107	PRO
1	C	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%)	280 (94%)	17 (6%)	18	39
1	B	297/320 (93%)	287 (97%)	10 (3%)	32	60
1	C	297/320 (93%)	284 (96%)	13 (4%)	25	50
1	D	297/320 (93%)	277 (93%)	20 (7%)	15	31
1	E	297/320 (93%)	282 (95%)	15 (5%)	21	43
1	F	297/320 (93%)	279 (94%)	18 (6%)	17	35
All	All	1782/1920 (93%)	1689 (95%)	93 (5%)	21	42

All (93) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	41	LYS
1	A	115	LYS
1	A	120	GLN
1	A	203	ARG
1	A	211	LEU
1	A	234	GLU
1	A	241	LYS
1	A	242	LYS
1	A	281	ARG
1	A	295	MET
1	A	316	ILE
1	A	317	ARG
1	A	322	VAL
1	A	323	THR
1	A	344	GLU
1	A	351	ASP

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Mol	Chain	Res	Type
1	A	357	SER
1	B	54	PHE
1	B	77	ARG
1	B	137	GLU
1	B	196	LEU
1	B	233	GLN
1	B	238	ARG
1	B	249	PRO
1	B	315	LEU
1	B	316	ILE
1	B	352	SER
1	C	52	ARG
1	C	59	GLN
1	C	110	ARG
1	C	119	ARG
1	C	129	SER
1	C	211	LEU
1	C	280	LEU
1	C	295	MET
1	C	297	ILE
1	C	315	LEU
1	C	316	ILE
1	C	318	LEU
1	C	326	MET
1	D	52	ARG
1	D	54	PHE
1	D	83	GLU
1	D	119	ARG
1	D	123	GLU
1	D	211	LEU
1	D	221	LEU
1	D	225	GLN
1	D	233	GLN
1	D	242	LYS
1	D	252	ILE
1	D	256	VAL
1	D	314	VAL
1	D	315	LEU
1	D	316	ILE
1	D	318	LEU
1	D	326	MET
1	D	351	ASP

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Mol	Chain	Res	Type
1	D	352	SER
1	D	361	GLN
1	E	42	THR
1	E	52	ARG
1	E	80	ASP
1	E	83	GLU
1	E	86	MET
1	E	121	VAL
1	E	233	GLN
1	E	234	GLU
1	E	281	ARG
1	E	295	MET
1	E	316	ILE
1	E	324	LEU
1	E	325	MET
1	E	340	GLN
1	E	344	GLU
1	F	75	VAL
1	F	77	ARG
1	F	83	GLU
1	F	90	VAL
1	F	110	ARG
1	F	120	GLN
1	F	203	ARG
1	F	221	LEU
1	F	233	GLN
1	F	248	LEU
1	F	252	ILE
1	F	262	LEU
1	F	281	ARG
1	F	295	MET
1	F	316	ILE
1	F	347	HIS
1	F	352	SER
1	F	364	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (73) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	48	ASN
1	A	59	GLN
1	A	140	GLN

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Mol	Chain	Res	Type
1	A	161	HIS
1	A	215	ASN
1	A	224	GLN
1	A	225	GLN
1	A	233	GLN
1	A	268	HIS
1	A	283	GLN
1	A	293	GLN
1	A	307	GLN
1	B	48	ASN
1	B	59	GLN
1	B	67	ASN
1	B	101	HIS
1	B	224	GLN
1	B	225	GLN
1	B	233	GLN
1	B	251	ASN
1	B	257	GLN
1	B	283	GLN
1	B	293	GLN
1	B	307	GLN
1	B	340	GLN
1	B	353	ASN
1	C	48	ASN
1	C	59	GLN
1	C	120	GLN
1	C	134	ASN
1	C	225	GLN
1	C	233	GLN
1	C	251	ASN
1	C	257	GLN
1	C	293	GLN
1	C	320	GLN
1	C	369	GLN
1	D	48	ASN
1	D	59	GLN
1	D	98	HIS
1	D	101	HIS
1	D	224	GLN
1	D	225	GLN
1	D	233	GLN
1	D	251	ASN

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Mol	Chain	Res	Type
1	D	283	GLN
1	D	293	GLN
1	D	353	ASN
1	E	48	ASN
1	E	49	GLN
1	E	140	GLN
1	E	212	GLN
1	E	215	ASN
1	E	224	GLN
1	E	225	GLN
1	E	233	GLN
1	E	257	GLN
1	E	293	GLN
1	E	320	GLN
1	E	353	ASN
1	F	48	ASN
1	F	49	GLN
1	F	59	GLN
1	F	120	GLN
1	F	225	GLN
1	F	233	GLN
1	F	251	ASN
1	F	257	GLN
1	F	260	ASN
1	F	305	ASN
1	F	340	GLN
1	F	353	ASN
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 ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

10 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	8PH	A	403	-	18,18,26	2.32	7 (38%)	26,29,37	1.41	5 (19%)
2	8PH	F	401	-	26,26,26	1.97	7 (26%)	34,37,37	1.30	6 (17%)
2	8PH	A	402	-	7,7,26	0.36	0	6,6,37	0.75	0
2	8PH	C	401	-	26,26,26	1.93	7 (26%)	34,37,37	1.30	6 (17%)
2	8PH	B	401	-	26,26,26	1.95	7 (26%)	34,37,37	1.33	6 (17%)
2	8PH	B	402	-	8,8,26	0.35	0	7,7,37	0.83	0
2	8PH	D	401	-	26,26,26	1.95	7 (26%)	34,37,37	1.24	4 (11%)
2	8PH	F	402	-	26,26,26	1.99	8 (30%)	34,37,37	1.33	5 (14%)
2	8PH	E	401	-	26,26,26	1.96	7 (26%)	34,37,37	1.31	5 (14%)
2	8PH	A	401	-	26,26,26	1.97	7 (26%)	34,37,37	1.25	5 (14%)

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	8PH	A	403	-	-	3/25/25/33	0/1/1/1
2	8PH	F	401	-	-	6/33/33/33	0/1/1/1
2	8PH	A	402	-	-	1/5/5/33	-
2	8PH	C	401	-	-	3/33/33/33	0/1/1/1
2	8PH	B	401	-	-	13/33/33/33	0/1/1/1
2	8PH	B	402	-	-	1/6/6/33	-
2	8PH	D	401	-	-	23/33/33/33	0/1/1/1
2	8PH	F	402	-	-	17/33/33/33	0/1/1/1
2	8PH	E	401	-	-	15/33/33/33	0/1/1/1
2	8PH	A	401	-	-	7/33/33/33	0/1/1/1

All (57) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	402	8PH	PAZ-OAH	6.86	1.61	1.50
2	B	401	8PH	PAZ-OAH	6.82	1.61	1.50
2	A	401	8PH	PAZ-OAH	6.81	1.61	1.50
2	C	401	8PH	PAZ-OAH	6.81	1.61	1.50
2	F	401	8PH	PAZ-OAH	6.79	1.61	1.50
2	D	401	8PH	PAZ-OAH	6.77	1.61	1.50
2	E	401	8PH	PAZ-OAH	6.76	1.61	1.50
2	A	403	8PH	PAZ-OAH	6.75	1.61	1.50
2	A	401	8PH	PAZ-CAX	3.10	1.87	1.85
2	F	402	8PH	PAZ-CAX	2.98	1.87	1.85
2	F	401	8PH	PAZ-CAX	2.95	1.87	1.85
2	E	401	8PH	PAZ-CAX	2.91	1.87	1.85
2	F	402	8PH	PAY-CAX	2.88	1.87	1.85
2	A	403	8PH	PAY-CAX	2.77	1.87	1.85
2	D	401	8PH	PAZ-CAX	2.77	1.87	1.85
2	E	401	8PH	PAY-CAX	2.74	1.87	1.85
2	D	401	8PH	PAY-CAX	2.74	1.87	1.85
2	B	401	8PH	PAZ-CAX	2.73	1.87	1.85
2	A	403	8PH	PAZ-CAX	2.70	1.87	1.85
2	B	401	8PH	PAY-CAX	2.70	1.87	1.85
2	F	401	8PH	PAY-CAX	2.67	1.87	1.85
2	C	401	8PH	PAZ-CAX	2.66	1.87	1.85
2	C	401	8PH	PAY-CAX	2.59	1.87	1.85
2	F	402	8PH	PAZ-OAC	-2.56	1.48	1.54
2	A	401	8PH	PAZ-OAC	-2.56	1.48	1.54
2	F	401	8PH	PAZ-OAC	-2.54	1.48	1.54
2	A	403	8PH	PAZ-OAG	2.53	1.61	1.54
2	A	403	8PH	PAZ-OAC	-2.52	1.48	1.54
2	D	401	8PH	PAY-OAF	2.52	1.61	1.54
2	E	401	8PH	PAY-OAE	2.52	1.61	1.54
2	A	401	8PH	PAY-CAX	2.51	1.86	1.85
2	D	401	8PH	PAY-OAE	2.51	1.61	1.54
2	F	401	8PH	PAY-OAF	2.51	1.61	1.54
2	F	402	8PH	PAZ-OAG	2.51	1.61	1.54
2	E	401	8PH	PAY-OAF	2.51	1.61	1.54
2	A	401	8PH	PAZ-OAG	2.51	1.61	1.54
2	C	401	8PH	PAY-OAF	2.51	1.61	1.54
2	E	401	8PH	PAZ-OAC	-2.51	1.48	1.54
2	E	401	8PH	PAZ-OAG	2.51	1.61	1.54
2	B	401	8PH	PAZ-OAC	-2.51	1.48	1.54
2	F	402	8PH	PAY-OAF	2.50	1.61	1.54
2	C	401	8PH	PAY-OAE	2.50	1.61	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	403	8PH	PAY-OAF	2.50	1.61	1.54
2	F	401	8PH	PAY-OAE	2.50	1.61	1.54
2	F	401	8PH	PAZ-OAG	2.50	1.61	1.54
2	A	401	8PH	PAY-OAE	2.49	1.61	1.54
2	F	402	8PH	PAY-OAE	2.49	1.61	1.54
2	D	401	8PH	PAZ-OAC	-2.49	1.48	1.54
2	D	401	8PH	PAZ-OAG	2.49	1.61	1.54
2	B	401	8PH	PAZ-OAG	2.48	1.61	1.54
2	A	403	8PH	PAY-OAE	2.48	1.61	1.54
2	B	401	8PH	PAY-OAE	2.48	1.61	1.54
2	C	401	8PH	PAZ-OAG	2.48	1.61	1.54
2	A	401	8PH	PAY-OAF	2.48	1.61	1.54
2	C	401	8PH	PAZ-OAC	-2.47	1.48	1.54
2	B	401	8PH	PAY-OAF	2.45	1.61	1.54
2	F	402	8PH	CAI-NAV	-2.03	1.33	1.37

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	402	8PH	OAC-PAZ-CAX	3.07	112.95	106.19
2	F	401	8PH	CAX-CAU-NAW	-3.01	109.84	113.67
2	B	401	8PH	OAC-PAZ-CAX	3.00	112.78	106.19
2	C	401	8PH	OAC-PAZ-CAX	2.99	112.76	106.19
2	A	401	8PH	OAC-PAZ-CAX	2.98	112.74	106.19
2	A	403	8PH	OAC-PAZ-CAX	2.95	112.69	106.19
2	A	403	8PH	CAX-CAU-NAW	-2.93	109.95	113.67
2	D	401	8PH	OAC-PAZ-CAX	2.93	112.62	106.19
2	A	401	8PH	OAG-PAZ-OAH	-2.87	105.92	112.96
2	E	401	8PH	CAX-CAU-NAW	-2.87	110.02	113.67
2	E	401	8PH	OAC-PAZ-CAX	2.84	112.44	106.19
2	F	402	8PH	OAG-PAZ-OAH	-2.84	106.00	112.96
2	F	401	8PH	PAY-CAX-OAD	2.83	113.90	107.42
2	D	401	8PH	OAG-PAZ-OAH	-2.82	106.05	112.96
2	F	401	8PH	OAC-PAZ-CAX	2.82	112.39	106.19
2	A	403	8PH	OAG-PAZ-OAH	-2.80	106.11	112.96
2	B	401	8PH	OAG-PAZ-OAH	-2.76	106.20	112.96
2	F	402	8PH	PAY-CAX-OAD	2.75	113.72	107.42
2	E	401	8PH	OAG-PAZ-OAH	-2.74	106.24	112.96
2	C	401	8PH	OAG-PAZ-OAH	-2.74	106.26	112.96
2	F	401	8PH	OAG-PAZ-OAH	-2.74	106.26	112.96
2	A	401	8PH	PAZ-CAX-PAY	-2.72	107.84	112.72
2	B	401	8PH	PAY-CAX-OAD	2.72	113.65	107.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	401	8PH	PAY-CAX-OAD	2.71	113.62	107.42
2	F	402	8PH	PAZ-CAX-OAD	2.67	113.54	107.42
2	C	401	8PH	PAZ-CAX-OAD	2.62	113.43	107.42
2	A	401	8PH	PAY-CAX-OAD	2.61	113.40	107.42
2	A	403	8PH	PAZ-CAX-OAD	2.61	113.39	107.42
2	F	401	8PH	PAZ-CAX-OAD	2.61	113.39	107.42
2	A	403	8PH	PAY-CAX-OAD	2.60	113.37	107.42
2	B	401	8PH	PAZ-CAX-OAD	2.59	113.34	107.42
2	A	401	8PH	PAZ-CAX-OAD	2.53	113.21	107.42
2	D	401	8PH	PAY-CAX-OAD	2.49	113.12	107.42
2	E	401	8PH	PAY-CAX-OAD	2.45	113.03	107.42
2	D	401	8PH	PAZ-CAX-OAD	2.40	112.91	107.42
2	C	401	8PH	CAX-CAU-NAW	-2.30	110.74	113.67
2	F	401	8PH	PAZ-CAX-PAY	-2.30	108.59	112.72
2	C	401	8PH	PAZ-CAX-PAY	-2.29	108.62	112.72
2	B	401	8PH	PAZ-CAX-PAY	-2.21	108.77	112.72
2	E	401	8PH	PAZ-CAX-OAD	2.20	112.45	107.42
2	B	401	8PH	CAU-NAW-CAJ	-2.16	123.00	127.02
2	F	402	8PH	PAZ-CAX-PAY	-2.08	109.00	112.72

There are no chirality outliers.

All (89) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	403	8PH	NAW-CAU-CAX-PAY
2	A	403	8PH	NAW-CAU-CAX-PAZ
2	A	403	8PH	NAW-CAU-CAX-OAD
2	B	401	8PH	PAY-CAX-PAZ-OAH
2	B	401	8PH	OAD-CAX-PAZ-OAH
2	B	401	8PH	OAD-CAX-PAZ-OAC
2	B	401	8PH	CAU-CAX-PAZ-OAH
2	B	401	8PH	CAX-CAU-NAW-CAJ
2	D	401	8PH	OAD-CAX-PAY-OAE
2	D	401	8PH	CAU-CAX-PAY-OAE
2	D	401	8PH	OAD-CAX-PAY-OAB
2	D	401	8PH	CAU-CAX-PAY-OAB
2	D	401	8PH	PAY-CAX-PAZ-OAG
2	D	401	8PH	NAW-CAU-CAX-PAY
2	D	401	8PH	NAW-CAU-CAX-PAZ
2	E	401	8PH	PAZ-CAX-PAY-OAE
2	E	401	8PH	OAD-CAX-PAY-OAE
2	E	401	8PH	CAU-CAX-PAY-OAE

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Mol	Chain	Res	Type	Atoms
2	E	401	8PH	PAZ-CAX-PAY-OAF
2	E	401	8PH	OAD-CAX-PAY-OAF
2	E	401	8PH	CAU-CAX-PAY-OAF
2	E	401	8PH	PAZ-CAX-PAY-OAB
2	E	401	8PH	OAD-CAX-PAY-OAB
2	E	401	8PH	CAU-CAX-PAY-OAB
2	E	401	8PH	OAD-CAX-PAZ-OAH
2	F	402	8PH	PAY-CAX-PAZ-OAH
2	F	402	8PH	PAY-CAX-PAZ-OAC
2	F	402	8PH	OAD-CAX-PAZ-OAG
2	F	402	8PH	OAD-CAX-PAZ-OAH
2	F	402	8PH	OAD-CAX-PAZ-OAC
2	F	402	8PH	CAU-CAX-PAZ-OAG
2	F	402	8PH	CAU-CAX-PAZ-OAH
2	F	402	8PH	CAU-CAX-PAZ-OAC
2	D	401	8PH	CAP-CAQ-CAR-CAS
2	C	401	8PH	CAP-CAQ-CAR-CAS
2	F	402	8PH	CAP-CAQ-CAR-CAS
2	B	401	8PH	CAM-CAN-CAO-CAP
2	D	401	8PH	PAY-CAX-PAZ-OAC
2	F	402	8PH	PAY-CAX-PAZ-OAG
2	A	401	8PH	CAP-CAQ-CAR-CAS
2	E	401	8PH	CAQ-CAR-CAS-CAT
2	A	401	8PH	CAX-CAU-NAW-CAJ
2	F	401	8PH	CAS-CAT-NAV-CAI
2	B	401	8PH	CAP-CAQ-CAR-CAS
2	B	401	8PH	CAS-CAT-NAV-CAI
2	E	401	8PH	CAS-CAT-NAV-CAI
2	B	401	8PH	CAS-CAT-NAV-CAK
2	E	401	8PH	CAS-CAT-NAV-CAK
2	A	401	8PH	OAD-CAX-PAY-OAE
2	D	401	8PH	OAD-CAX-PAY-OAF
2	D	401	8PH	OAD-CAX-PAZ-OAG
2	E	401	8PH	NAW-CAU-CAX-PAY
2	F	402	8PH	CAS-CAT-NAV-CAK
2	B	401	8PH	CAU-CAX-PAZ-OAC
2	D	401	8PH	CAU-CAX-PAZ-OAG
2	E	401	8PH	CAU-CAX-PAZ-OAH
2	C	401	8PH	CAS-CAT-NAV-CAI
2	A	402	8PH	CAO-CAP-CAQ-CAR
2	C	401	8PH	CAS-CAT-NAV-CAK
2	B	402	8PH	CAO-CAP-CAQ-CAR

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Mol	Chain	Res	Type	Atoms
2	F	402	8PH	CAQ-CAR-CAS-CAT
2	F	401	8PH	CAP-CAQ-CAR-CAS
2	F	401	8PH	CAX-CAU-NAW-CAK
2	F	402	8PH	CAO-CAP-CAQ-CAR
2	D	401	8PH	CAO-CAP-CAQ-CAR
2	B	401	8PH	PAY-CAX-PAZ-OAC
2	B	401	8PH	OAD-CAX-PAZ-OAG
2	D	401	8PH	PAZ-CAX-PAY-OAE
2	D	401	8PH	CAU-CAX-PAZ-OAC
2	D	401	8PH	NAW-CAU-CAX-OAD
2	F	402	8PH	CAN-CAO-CAP-CAQ
2	D	401	8PH	PAZ-CAX-PAY-OAB
2	D	401	8PH	PAY-CAX-PAZ-OAH
2	D	401	8PH	OAD-CAX-PAZ-OAH
2	F	401	8PH	OAD-CAX-PAZ-OAH
2	F	402	8PH	NAW-CAU-CAX-PAY
2	F	402	8PH	NAW-CAU-CAX-PAZ
2	F	402	8PH	CAS-CAT-NAV-CAI
2	A	401	8PH	CAQ-CAR-CAS-CAT
2	A	401	8PH	CAU-CAX-PAY-OAE
2	A	401	8PH	CAU-CAX-PAY-OAB
2	D	401	8PH	CAU-CAX-PAY-OAF
2	D	401	8PH	CAU-CAX-PAZ-OAH
2	F	401	8PH	CAS-CAT-NAV-CAK
2	A	401	8PH	CAX-CAU-NAW-CAK
2	B	401	8PH	CAX-CAU-NAW-CAK
2	D	401	8PH	CAX-CAU-NAW-CAK
2	D	401	8PH	CAX-CAU-NAW-CAJ
2	F	401	8PH	CAX-CAU-NAW-CAJ

There are no ring outliers.

10 monomers are involved in 25 short contacts:

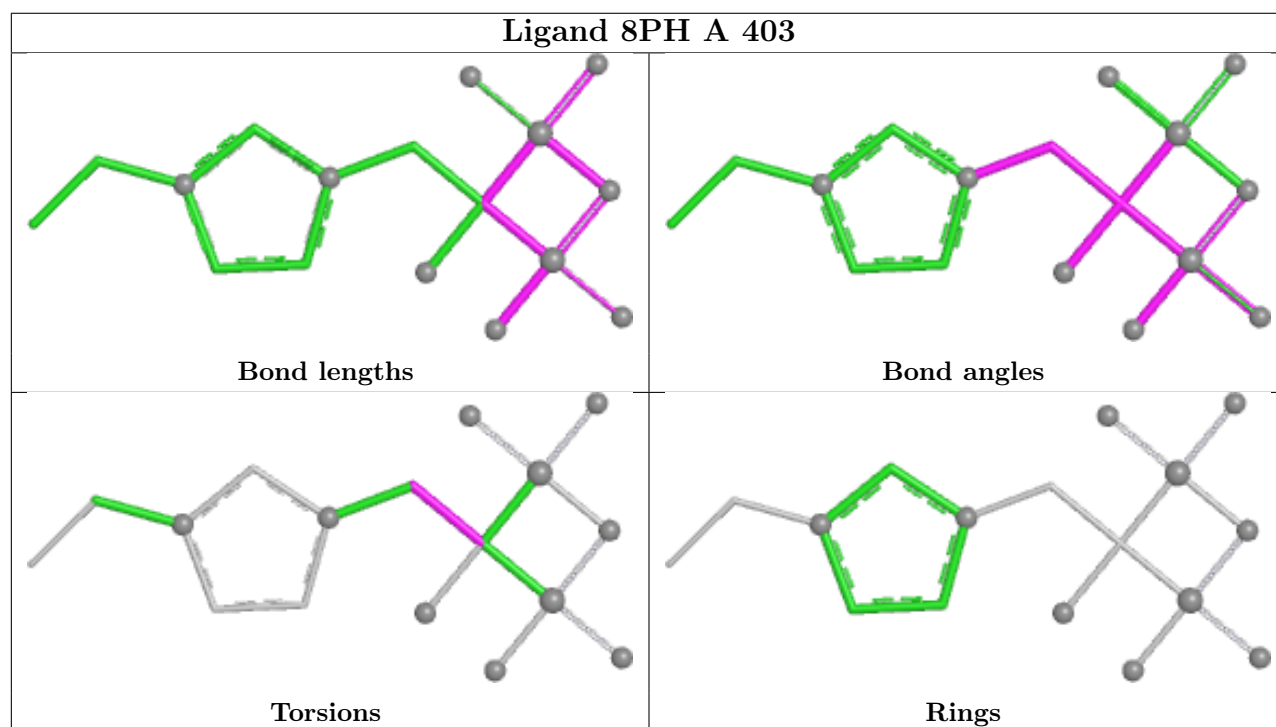
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	403	8PH	1	0
2	F	401	8PH	4	0
2	A	402	8PH	1	0
2	C	401	8PH	4	0
2	B	401	8PH	6	0
2	B	402	8PH	2	0
2	D	401	8PH	2	0
2	F	402	8PH	1	0

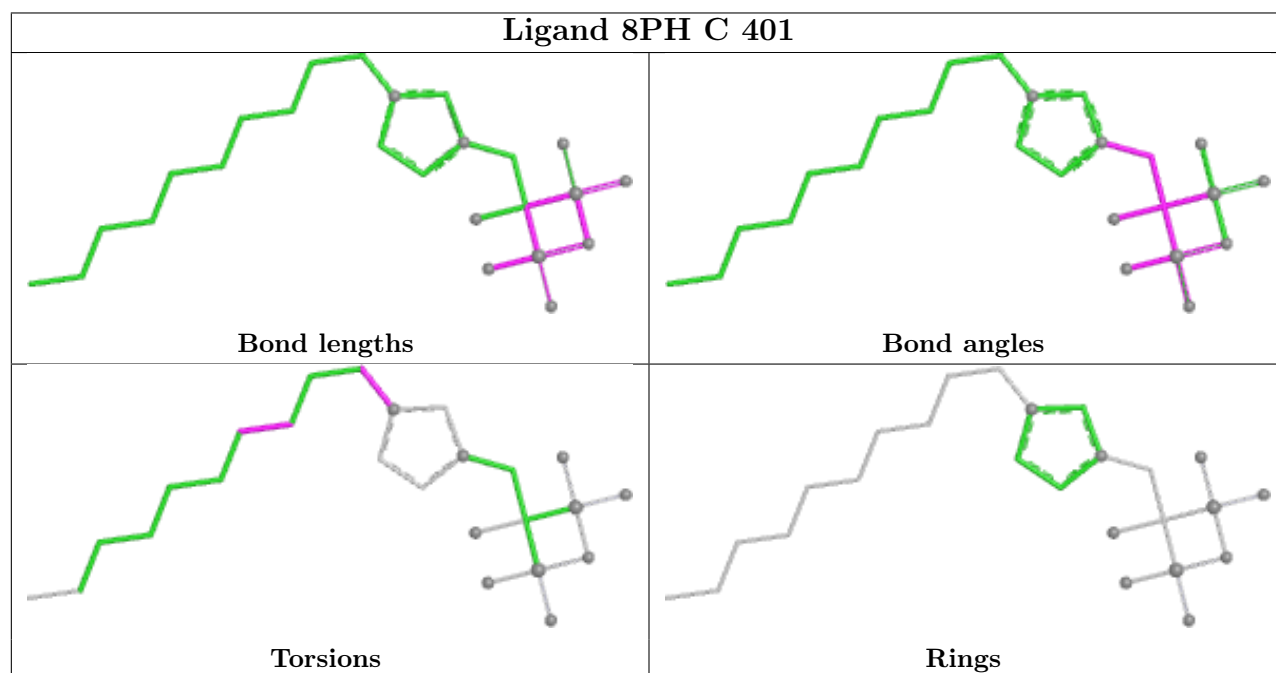
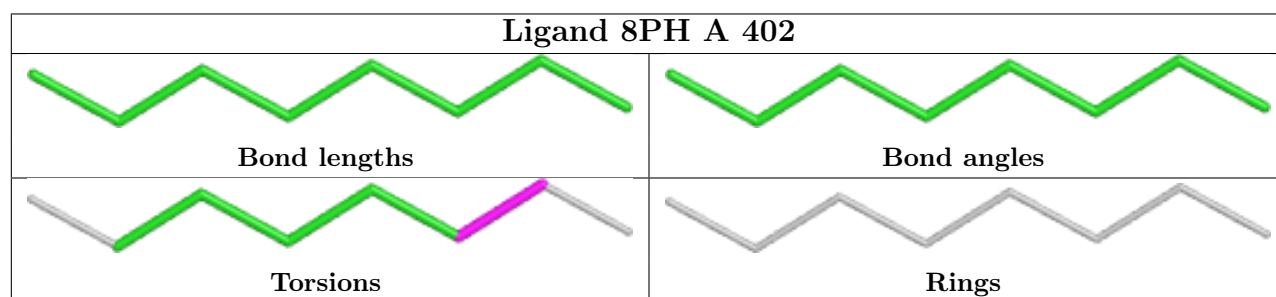
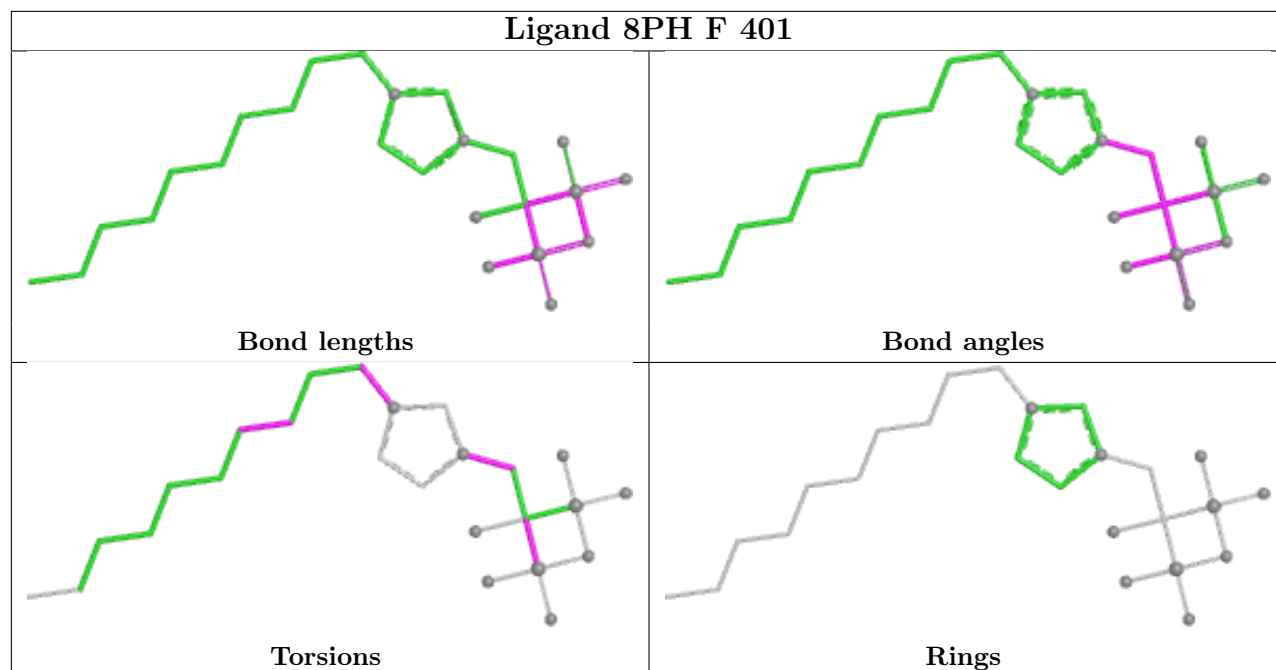
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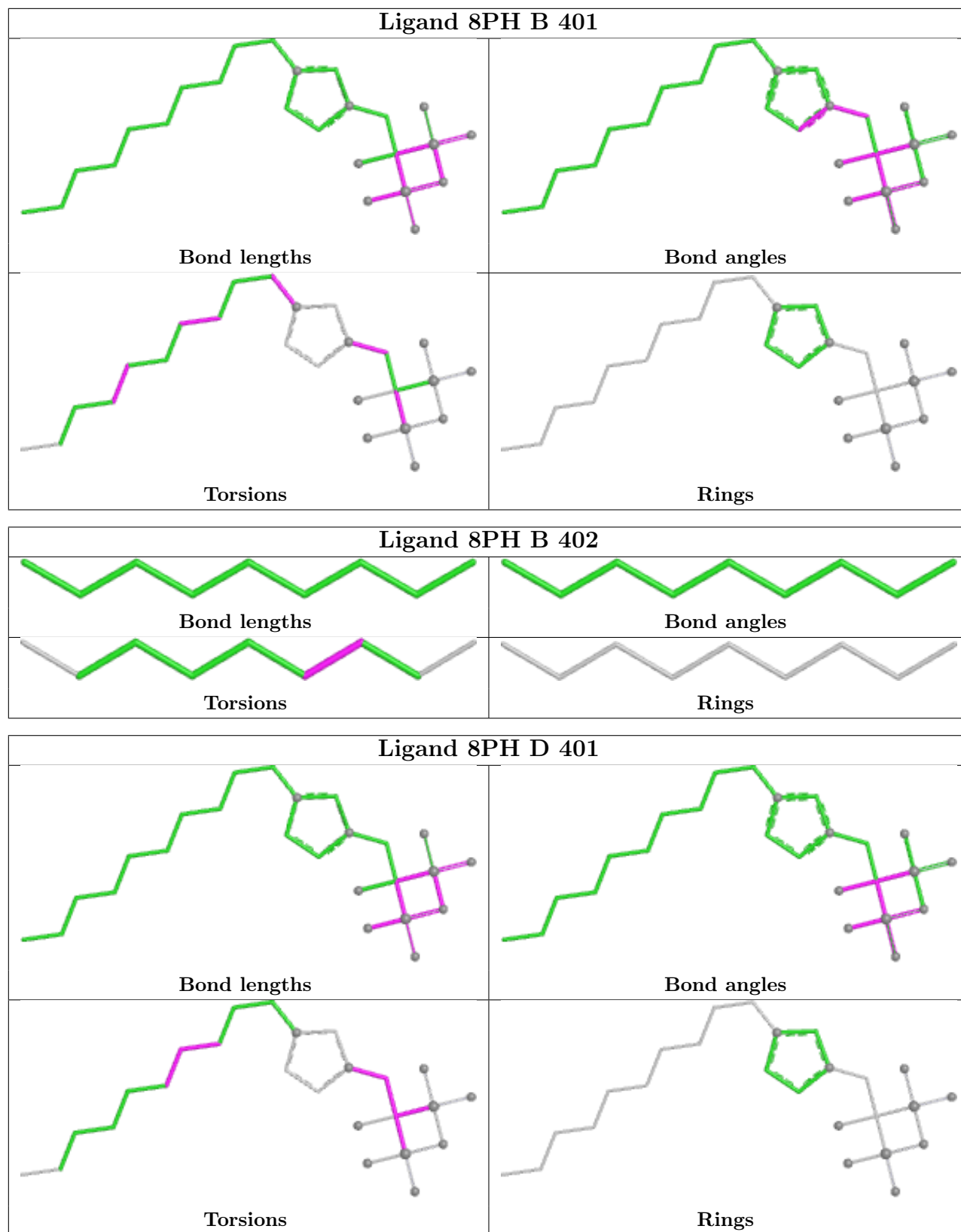
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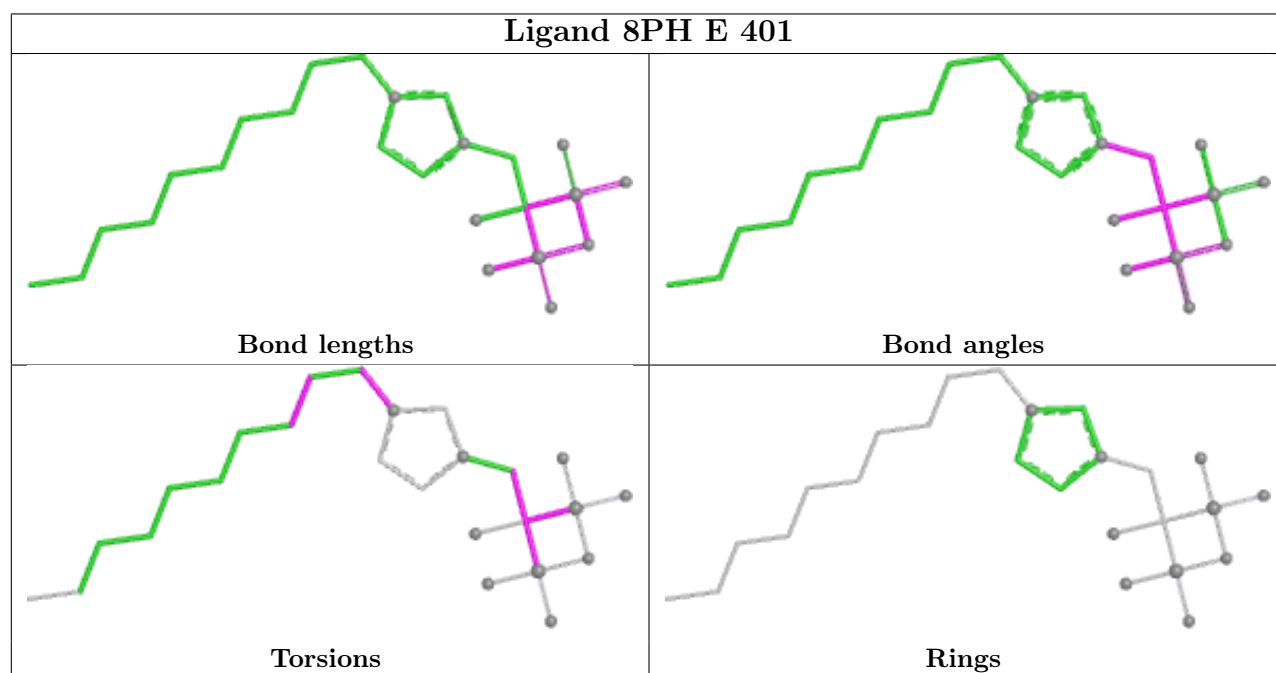
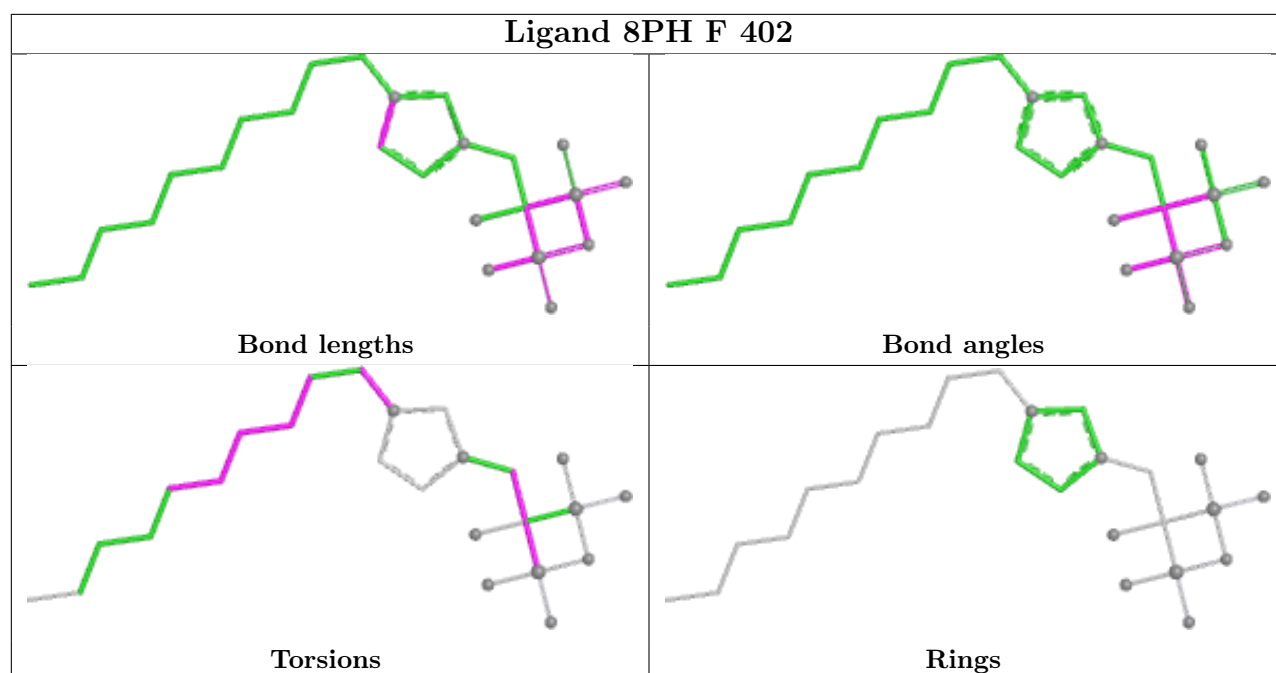
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	401	8PH	4	0
2	A	401	8PH	1	0

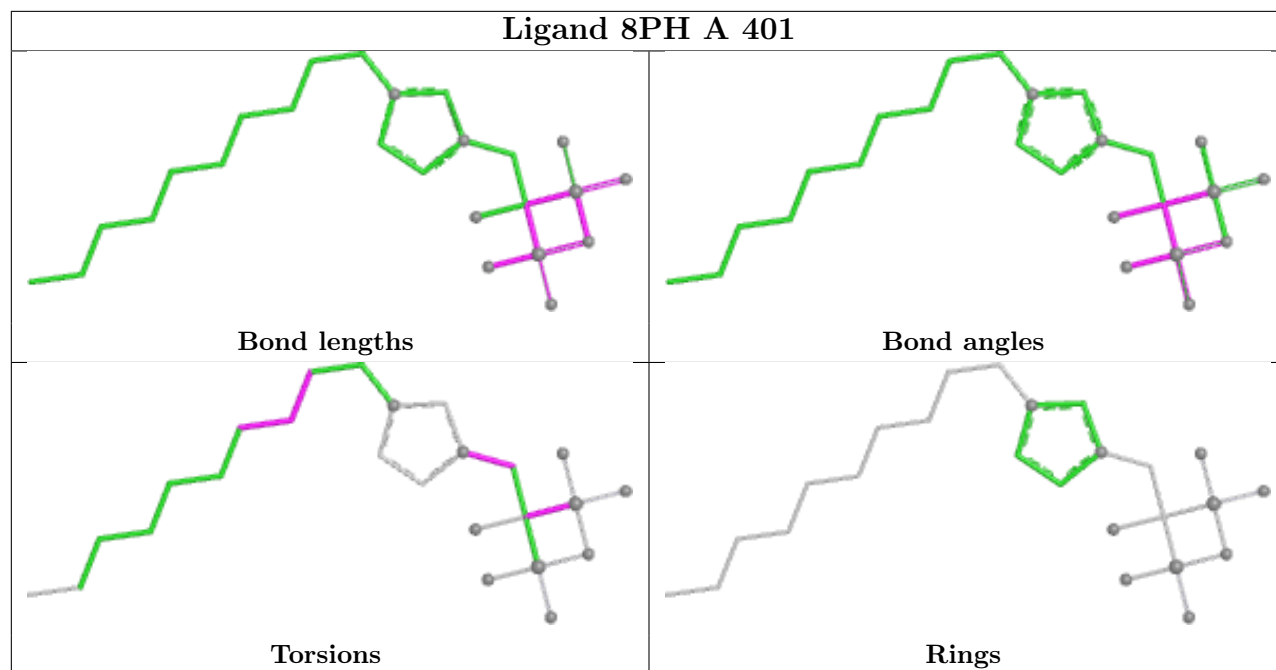
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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











5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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%)	-0.32	7 (2%) 63 59	17, 32, 68, 114	0
1	B	334/360 (92%)	-0.14	6 (1%) 67 64	20, 36, 74, 119	0
1	C	334/360 (92%)	-0.30	5 (1%) 72 68	19, 35, 63, 107	0
1	D	334/360 (92%)	0.13	10 (2%) 52 48	22, 48, 86, 123	0
1	E	334/360 (92%)	0.07	15 (4%) 38 33	28, 48, 85, 123	0
1	F	334/360 (92%)	0.21	13 (3%) 43 38	28, 56, 95, 137	0
All	All	2004/2160 (92%)	-0.06	56 (2%) 55 50	17, 41, 86, 137	0

All (56) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	321	ALA	5.3
1	E	322	VAL	4.9
1	D	321	ALA	4.5
1	E	323	THR	4.4
1	F	316	ILE	4.4
1	F	320	GLN	3.9
1	E	324	LEU	3.8
1	B	321	ALA	3.8
1	F	318	LEU	3.7
1	F	322	VAL	3.7
1	B	318	LEU	3.6
1	E	321	ALA	3.6
1	E	318	LEU	3.6
1	A	322	VAL	3.5
1	D	322	VAL	3.5
1	F	323	THR	3.5
1	F	314	VAL	3.4
1	F	324	LEU	3.4
1	E	315	LEU	3.3

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Mol	Chain	Res	Type	RSRZ
1	E	320	GLN	3.2
1	B	322	VAL	3.2
1	D	318	LEU	3.2
1	D	315	LEU	3.0
1	A	321	ALA	2.9
1	A	36	LEU	2.9
1	C	321	ALA	2.8
1	A	318	LEU	2.8
1	F	369	GLN	2.8
1	D	324	LEU	2.7
1	E	326	MET	2.7
1	F	315	LEU	2.7
1	A	320	GLN	2.6
1	D	36	LEU	2.6
1	C	322	VAL	2.6
1	F	313	ALA	2.6
1	C	315	LEU	2.6
1	F	36	LEU	2.5
1	A	315	LEU	2.5
1	D	319	GLY	2.5
1	F	319	GLY	2.5
1	E	36	LEU	2.4
1	B	324	LEU	2.4
1	E	316	ILE	2.3
1	B	319	GLY	2.3
1	E	53	SER	2.3
1	E	319	GLY	2.2
1	E	325	MET	2.2
1	C	317	ARG	2.2
1	D	316	ILE	2.2
1	D	320	GLN	2.2
1	C	36	LEU	2.2
1	D	347	HIS	2.1
1	E	369	GLN	2.1
1	E	327	ASP	2.1
1	A	314	VAL	2.0
1	B	36	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

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](#)

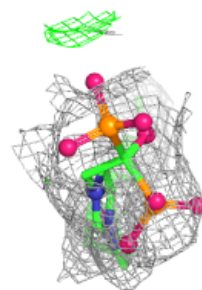
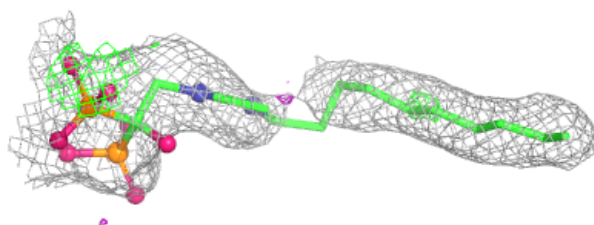
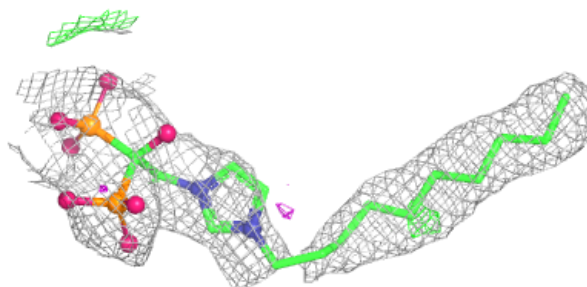
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	8PH	E	401	26/26	0.71	0.17	47,93,106,106	0
2	8PH	A	403	18/26	0.75	0.15	101,106,107,107	0
2	8PH	F	402	26/26	0.78	0.15	62,101,110,110	0
2	8PH	D	401	26/26	0.81	0.14	52,103,114,115	0
2	8PH	F	401	26/26	0.83	0.13	50,92,102,102	0
2	8PH	C	401	26/26	0.84	0.13	33,86,96,97	0
2	8PH	B	401	26/26	0.87	0.12	31,81,93,94	0
2	8PH	A	401	26/26	0.88	0.11	27,54,69,70	0
2	8PH	B	402	9/26	0.90	0.11	38,38,40,42	0
2	8PH	A	402	8/26	0.94	0.11	36,38,40,42	0

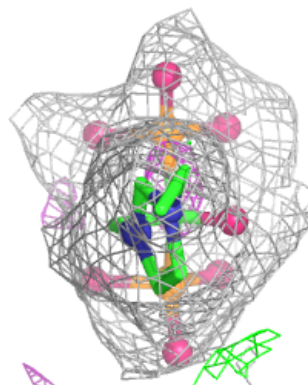
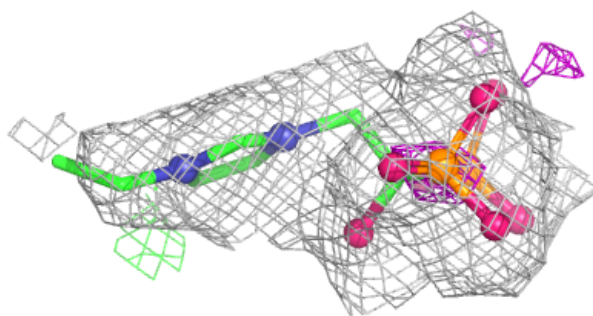
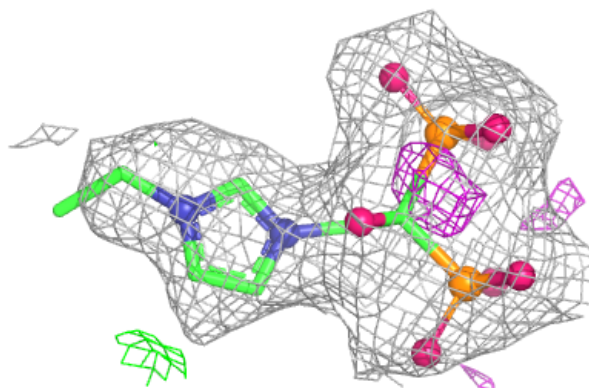
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around 8PH E 401:

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

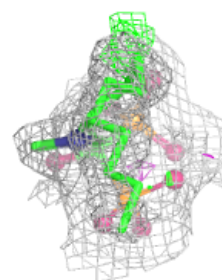
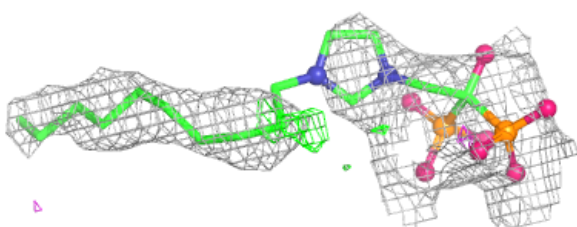
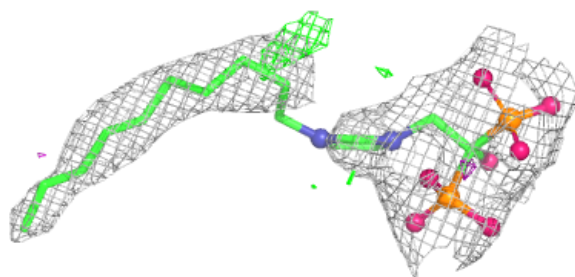
**Electron density around 8PH A 403:**

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

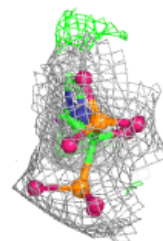
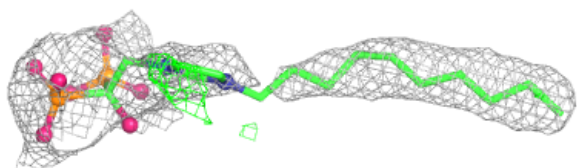
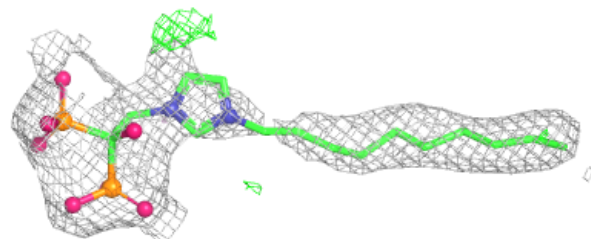


Electron density around 8PH F 402:

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

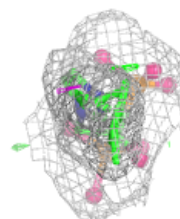
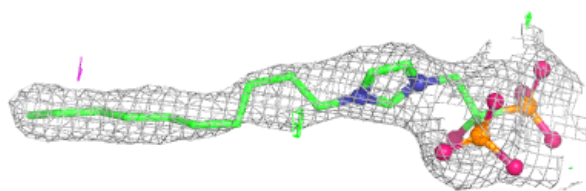
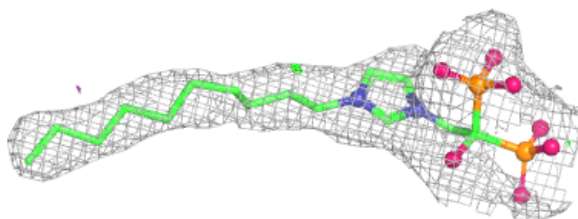
**Electron density around 8PH D 401:**

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

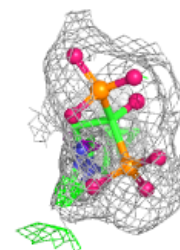
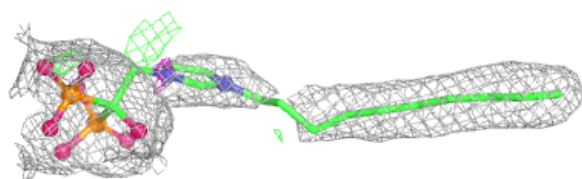
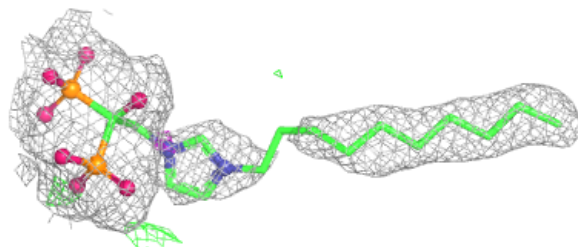


Electron density around 8PH F 401:

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

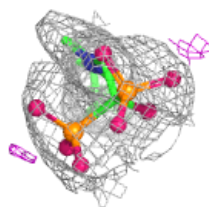
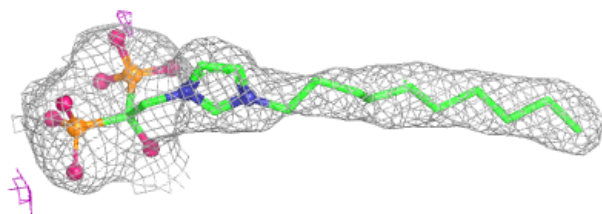
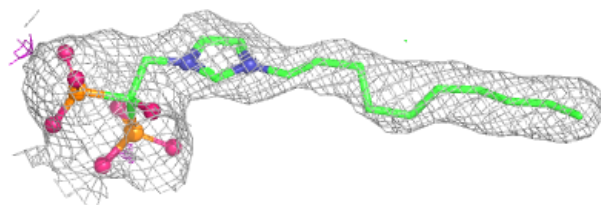
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and green (positive)

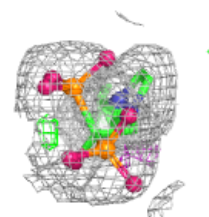
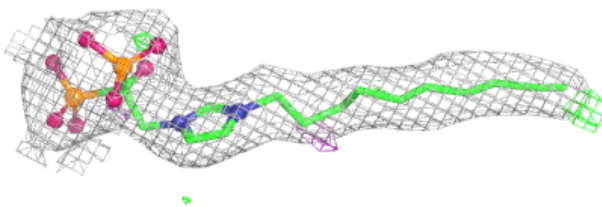
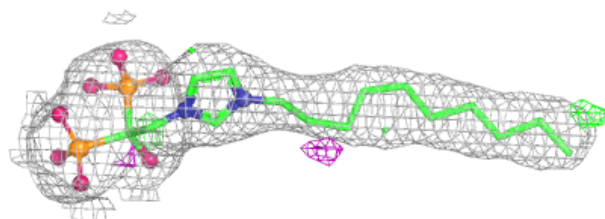


Electron density around 8PH B 401:

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

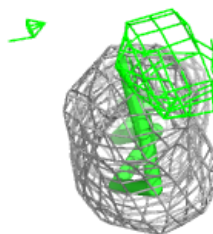
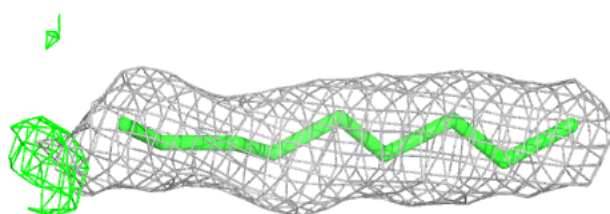
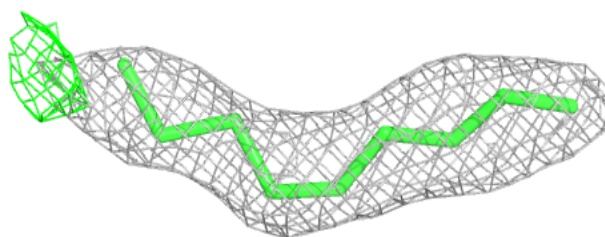
**Electron density around 8PH A 401:**

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

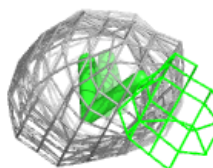
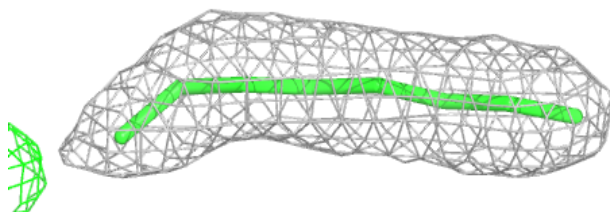
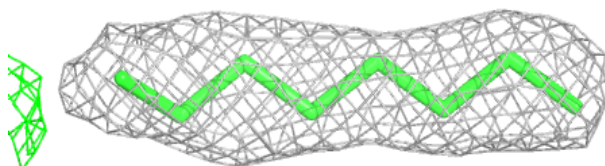


Electron density around 8PH B 402:

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

**Electron density around 8PH A 402:**

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



6.5 Other polymers [i](#)

There are no such residues in this entry.