



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 05:56 PM EDT

PDB ID : 6UNZ / pdb_00006unz
Title : Crystal structure of cytosolic fumarate hydratase from Leishmania major
Authors : Feliciano, P.R.; Drennan, C.L.
Deposited on : 2019-10-14
Resolution : 3.19 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

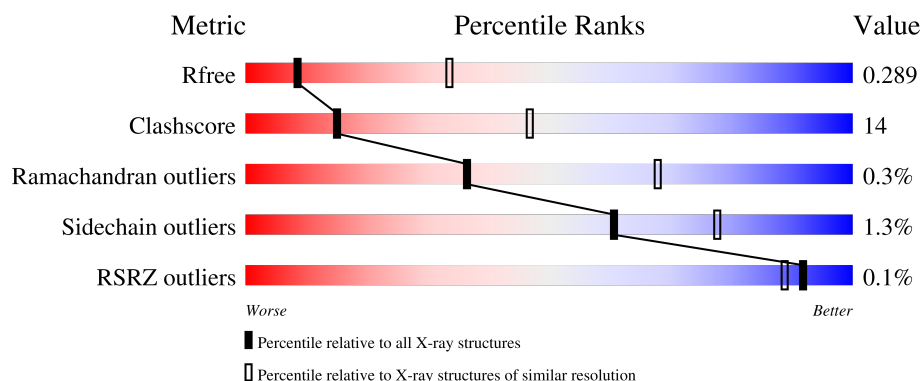
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.19 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	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	604	68% 21% 11%
1	B	604	65% 22% 12%
1	C	604	61% 26% 12%
1	D	604	66% 22% 12%
1	E	604	64% 23% 12%

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Mol	Chain	Length	Quality of chain
1	F	604	
1	G	604	
1	H	604	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	SF4	B	601	-	-	X	-
2	SF4	C	601	-	-	X	-
2	SF4	F	601	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 31246 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 fumarate hydratase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	539	Total	C	N	O	S	0	0	0
			3933	2469	688	758	18			
1	B	534	Total	C	N	O	S	0	0	0
			3933	2480	681	755	17			
1	C	532	Total	C	N	O	S	0	0	0
			3822	2387	678	737	20			
1	D	530	Total	C	N	O	S	0	0	0
			3874	2440	677	738	19			
1	E	531	Total	C	N	O	S	0	0	0
			3863	2429	677	736	21			
1	F	531	Total	C	N	O	S	0	1	0
			3981	2512	690	759	20			
1	G	532	Total	C	N	O	S	0	0	0
			3873	2430	684	742	17			
1	H	531	Total	C	N	O	S	0	0	0
			3888	2448	673	749	18			

There are 288 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-35	MET	-	expression tag	UNP E9AE57
A	-34	GLY	-	expression tag	UNP E9AE57
A	-33	SER	-	expression tag	UNP E9AE57
A	-32	SER	-	expression tag	UNP E9AE57
A	-31	HIS	-	expression tag	UNP E9AE57
A	-30	HIS	-	expression tag	UNP E9AE57
A	-29	HIS	-	expression tag	UNP E9AE57
A	-28	HIS	-	expression tag	UNP E9AE57
A	-27	HIS	-	expression tag	UNP E9AE57
A	-26	HIS	-	expression tag	UNP E9AE57
A	-25	SER	-	expression tag	UNP E9AE57
A	-24	SER	-	expression tag	UNP E9AE57
A	-23	GLY	-	expression tag	UNP E9AE57

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-22	LEU	-	expression tag	UNP E9AE57
A	-21	VAL	-	expression tag	UNP E9AE57
A	-20	PRO	-	expression tag	UNP E9AE57
A	-19	ARG	-	expression tag	UNP E9AE57
A	-18	GLY	-	expression tag	UNP E9AE57
A	-17	SER	-	expression tag	UNP E9AE57
A	-16	HIS	-	expression tag	UNP E9AE57
A	-15	MET	-	expression tag	UNP E9AE57
A	-14	ALA	-	expression tag	UNP E9AE57
A	-13	SER	-	expression tag	UNP E9AE57
A	-12	MET	-	expression tag	UNP E9AE57
A	-11	THR	-	expression tag	UNP E9AE57
A	-10	GLY	-	expression tag	UNP E9AE57
A	-9	GLY	-	expression tag	UNP E9AE57
A	-8	GLN	-	expression tag	UNP E9AE57
A	-7	GLN	-	expression tag	UNP E9AE57
A	-6	MET	-	expression tag	UNP E9AE57
A	-5	GLY	-	expression tag	UNP E9AE57
A	-4	ARG	-	expression tag	UNP E9AE57
A	-3	GLY	-	expression tag	UNP E9AE57
A	-2	SER	-	expression tag	UNP E9AE57
A	-1	GLU	-	expression tag	UNP E9AE57
A	0	PHE	-	expression tag	UNP E9AE57
B	-35	MET	-	expression tag	UNP E9AE57
B	-34	GLY	-	expression tag	UNP E9AE57
B	-33	SER	-	expression tag	UNP E9AE57
B	-32	SER	-	expression tag	UNP E9AE57
B	-31	HIS	-	expression tag	UNP E9AE57
B	-30	HIS	-	expression tag	UNP E9AE57
B	-29	HIS	-	expression tag	UNP E9AE57
B	-28	HIS	-	expression tag	UNP E9AE57
B	-27	HIS	-	expression tag	UNP E9AE57
B	-26	HIS	-	expression tag	UNP E9AE57
B	-25	SER	-	expression tag	UNP E9AE57
B	-24	SER	-	expression tag	UNP E9AE57
B	-23	GLY	-	expression tag	UNP E9AE57
B	-22	LEU	-	expression tag	UNP E9AE57
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B	-20	PRO	-	expression tag	UNP E9AE57
B	-19	ARG	-	expression tag	UNP E9AE57
B	-18	GLY	-	expression tag	UNP E9AE57
B	-17	SER	-	expression tag	UNP E9AE57

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-16	HIS	-	expression tag	UNP E9AE57
B	-15	MET	-	expression tag	UNP E9AE57
B	-14	ALA	-	expression tag	UNP E9AE57
B	-13	SER	-	expression tag	UNP E9AE57
B	-12	MET	-	expression tag	UNP E9AE57
B	-11	THR	-	expression tag	UNP E9AE57
B	-10	GLY	-	expression tag	UNP E9AE57
B	-9	GLY	-	expression tag	UNP E9AE57
B	-8	GLN	-	expression tag	UNP E9AE57
B	-7	GLN	-	expression tag	UNP E9AE57
B	-6	MET	-	expression tag	UNP E9AE57
B	-5	GLY	-	expression tag	UNP E9AE57
B	-4	ARG	-	expression tag	UNP E9AE57
B	-3	GLY	-	expression tag	UNP E9AE57
B	-2	SER	-	expression tag	UNP E9AE57
B	-1	GLU	-	expression tag	UNP E9AE57
B	0	PHE	-	expression tag	UNP E9AE57
C	-35	MET	-	expression tag	UNP E9AE57
C	-34	GLY	-	expression tag	UNP E9AE57
C	-33	SER	-	expression tag	UNP E9AE57
C	-32	SER	-	expression tag	UNP E9AE57
C	-31	HIS	-	expression tag	UNP E9AE57
C	-30	HIS	-	expression tag	UNP E9AE57
C	-29	HIS	-	expression tag	UNP E9AE57
C	-28	HIS	-	expression tag	UNP E9AE57
C	-27	HIS	-	expression tag	UNP E9AE57
C	-26	HIS	-	expression tag	UNP E9AE57
C	-25	SER	-	expression tag	UNP E9AE57
C	-24	SER	-	expression tag	UNP E9AE57
C	-23	GLY	-	expression tag	UNP E9AE57
C	-22	LEU	-	expression tag	UNP E9AE57
C	-21	VAL	-	expression tag	UNP E9AE57
C	-20	PRO	-	expression tag	UNP E9AE57
C	-19	ARG	-	expression tag	UNP E9AE57
C	-18	GLY	-	expression tag	UNP E9AE57
C	-17	SER	-	expression tag	UNP E9AE57
C	-16	HIS	-	expression tag	UNP E9AE57
C	-15	MET	-	expression tag	UNP E9AE57
C	-14	ALA	-	expression tag	UNP E9AE57
C	-13	SER	-	expression tag	UNP E9AE57
C	-12	MET	-	expression tag	UNP E9AE57
C	-11	THR	-	expression tag	UNP E9AE57

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-10	GLY	-	expression tag	UNP E9AE57
C	-9	GLY	-	expression tag	UNP E9AE57
C	-8	GLN	-	expression tag	UNP E9AE57
C	-7	GLN	-	expression tag	UNP E9AE57
C	-6	MET	-	expression tag	UNP E9AE57
C	-5	GLY	-	expression tag	UNP E9AE57
C	-4	ARG	-	expression tag	UNP E9AE57
C	-3	GLY	-	expression tag	UNP E9AE57
C	-2	SER	-	expression tag	UNP E9AE57
C	-1	GLU	-	expression tag	UNP E9AE57
C	0	PHE	-	expression tag	UNP E9AE57
D	-35	MET	-	expression tag	UNP E9AE57
D	-34	GLY	-	expression tag	UNP E9AE57
D	-33	SER	-	expression tag	UNP E9AE57
D	-32	SER	-	expression tag	UNP E9AE57
D	-31	HIS	-	expression tag	UNP E9AE57
D	-30	HIS	-	expression tag	UNP E9AE57
D	-29	HIS	-	expression tag	UNP E9AE57
D	-28	HIS	-	expression tag	UNP E9AE57
D	-27	HIS	-	expression tag	UNP E9AE57
D	-26	HIS	-	expression tag	UNP E9AE57
D	-25	SER	-	expression tag	UNP E9AE57
D	-24	SER	-	expression tag	UNP E9AE57
D	-23	GLY	-	expression tag	UNP E9AE57
D	-22	LEU	-	expression tag	UNP E9AE57
D	-21	VAL	-	expression tag	UNP E9AE57
D	-20	PRO	-	expression tag	UNP E9AE57
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D	-17	SER	-	expression tag	UNP E9AE57
D	-16	HIS	-	expression tag	UNP E9AE57
D	-15	MET	-	expression tag	UNP E9AE57
D	-14	ALA	-	expression tag	UNP E9AE57
D	-13	SER	-	expression tag	UNP E9AE57
D	-12	MET	-	expression tag	UNP E9AE57
D	-11	THR	-	expression tag	UNP E9AE57
D	-10	GLY	-	expression tag	UNP E9AE57
D	-9	GLY	-	expression tag	UNP E9AE57
D	-8	GLN	-	expression tag	UNP E9AE57
D	-7	GLN	-	expression tag	UNP E9AE57
D	-6	MET	-	expression tag	UNP E9AE57
D	-5	GLY	-	expression tag	UNP E9AE57

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-4	ARG	-	expression tag	UNP E9AE57
D	-3	GLY	-	expression tag	UNP E9AE57
D	-2	SER	-	expression tag	UNP E9AE57
D	-1	GLU	-	expression tag	UNP E9AE57
D	0	PHE	-	expression tag	UNP E9AE57
E	-35	MET	-	expression tag	UNP E9AE57
E	-34	GLY	-	expression tag	UNP E9AE57
E	-33	SER	-	expression tag	UNP E9AE57
E	-32	SER	-	expression tag	UNP E9AE57
E	-31	HIS	-	expression tag	UNP E9AE57
E	-30	HIS	-	expression tag	UNP E9AE57
E	-29	HIS	-	expression tag	UNP E9AE57
E	-28	HIS	-	expression tag	UNP E9AE57
E	-27	HIS	-	expression tag	UNP E9AE57
E	-26	HIS	-	expression tag	UNP E9AE57
E	-25	SER	-	expression tag	UNP E9AE57
E	-24	SER	-	expression tag	UNP E9AE57
E	-23	GLY	-	expression tag	UNP E9AE57
E	-22	LEU	-	expression tag	UNP E9AE57
E	-21	VAL	-	expression tag	UNP E9AE57
E	-20	PRO	-	expression tag	UNP E9AE57
E	-19	ARG	-	expression tag	UNP E9AE57
E	-18	GLY	-	expression tag	UNP E9AE57
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E	-8	GLN	-	expression tag	UNP E9AE57
E	-7	GLN	-	expression tag	UNP E9AE57
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E	-5	GLY	-	expression tag	UNP E9AE57
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E	-3	GLY	-	expression tag	UNP E9AE57
E	-2	SER	-	expression tag	UNP E9AE57
E	-1	GLU	-	expression tag	UNP E9AE57
E	0	PHE	-	expression tag	UNP E9AE57
F	-35	MET	-	expression tag	UNP E9AE57

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Chain	Residue	Modelled	Actual	Comment	Reference
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F	-33	SER	-	expression tag	UNP E9AE57
F	-32	SER	-	expression tag	UNP E9AE57
F	-31	HIS	-	expression tag	UNP E9AE57
F	-30	HIS	-	expression tag	UNP E9AE57
F	-29	HIS	-	expression tag	UNP E9AE57
F	-28	HIS	-	expression tag	UNP E9AE57
F	-27	HIS	-	expression tag	UNP E9AE57
F	-26	HIS	-	expression tag	UNP E9AE57
F	-25	SER	-	expression tag	UNP E9AE57
F	-24	SER	-	expression tag	UNP E9AE57
F	-23	GLY	-	expression tag	UNP E9AE57
F	-22	LEU	-	expression tag	UNP E9AE57
F	-21	VAL	-	expression tag	UNP E9AE57
F	-20	PRO	-	expression tag	UNP E9AE57
F	-19	ARG	-	expression tag	UNP E9AE57
F	-18	GLY	-	expression tag	UNP E9AE57
F	-17	SER	-	expression tag	UNP E9AE57
F	-16	HIS	-	expression tag	UNP E9AE57
F	-15	MET	-	expression tag	UNP E9AE57
F	-14	ALA	-	expression tag	UNP E9AE57
F	-13	SER	-	expression tag	UNP E9AE57
F	-12	MET	-	expression tag	UNP E9AE57
F	-11	THR	-	expression tag	UNP E9AE57
F	-10	GLY	-	expression tag	UNP E9AE57
F	-9	GLY	-	expression tag	UNP E9AE57
F	-8	GLN	-	expression tag	UNP E9AE57
F	-7	GLN	-	expression tag	UNP E9AE57
F	-6	MET	-	expression tag	UNP E9AE57
F	-5	GLY	-	expression tag	UNP E9AE57
F	-4	ARG	-	expression tag	UNP E9AE57
F	-3	GLY	-	expression tag	UNP E9AE57
F	-2	SER	-	expression tag	UNP E9AE57
F	-1	GLU	-	expression tag	UNP E9AE57
F	0	PHE	-	expression tag	UNP E9AE57
G	-35	MET	-	expression tag	UNP E9AE57
G	-34	GLY	-	expression tag	UNP E9AE57
G	-33	SER	-	expression tag	UNP E9AE57
G	-32	SER	-	expression tag	UNP E9AE57
G	-31	HIS	-	expression tag	UNP E9AE57
G	-30	HIS	-	expression tag	UNP E9AE57
G	-29	HIS	-	expression tag	UNP E9AE57

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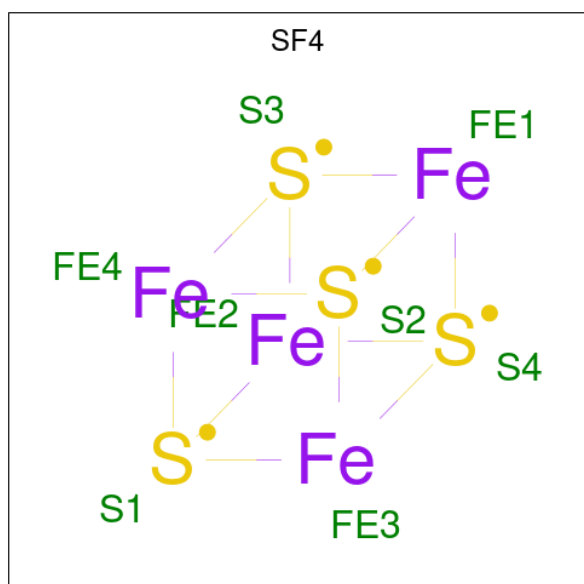
Chain	Residue	Modelled	Actual	Comment	Reference
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G	-27	HIS	-	expression tag	UNP E9AE57
G	-26	HIS	-	expression tag	UNP E9AE57
G	-25	SER	-	expression tag	UNP E9AE57
G	-24	SER	-	expression tag	UNP E9AE57
G	-23	GLY	-	expression tag	UNP E9AE57
G	-22	LEU	-	expression tag	UNP E9AE57
G	-21	VAL	-	expression tag	UNP E9AE57
G	-20	PRO	-	expression tag	UNP E9AE57
G	-19	ARG	-	expression tag	UNP E9AE57
G	-18	GLY	-	expression tag	UNP E9AE57
G	-17	SER	-	expression tag	UNP E9AE57
G	-16	HIS	-	expression tag	UNP E9AE57
G	-15	MET	-	expression tag	UNP E9AE57
G	-14	ALA	-	expression tag	UNP E9AE57
G	-13	SER	-	expression tag	UNP E9AE57
G	-12	MET	-	expression tag	UNP E9AE57
G	-11	THR	-	expression tag	UNP E9AE57
G	-10	GLY	-	expression tag	UNP E9AE57
G	-9	GLY	-	expression tag	UNP E9AE57
G	-8	GLN	-	expression tag	UNP E9AE57
G	-7	GLN	-	expression tag	UNP E9AE57
G	-6	MET	-	expression tag	UNP E9AE57
G	-5	GLY	-	expression tag	UNP E9AE57
G	-4	ARG	-	expression tag	UNP E9AE57
G	-3	GLY	-	expression tag	UNP E9AE57
G	-2	SER	-	expression tag	UNP E9AE57
G	-1	GLU	-	expression tag	UNP E9AE57
G	0	PHE	-	expression tag	UNP E9AE57
H	-35	MET	-	expression tag	UNP E9AE57
H	-34	GLY	-	expression tag	UNP E9AE57
H	-33	SER	-	expression tag	UNP E9AE57
H	-32	SER	-	expression tag	UNP E9AE57
H	-31	HIS	-	expression tag	UNP E9AE57
H	-30	HIS	-	expression tag	UNP E9AE57
H	-29	HIS	-	expression tag	UNP E9AE57
H	-28	HIS	-	expression tag	UNP E9AE57
H	-27	HIS	-	expression tag	UNP E9AE57
H	-26	HIS	-	expression tag	UNP E9AE57
H	-25	SER	-	expression tag	UNP E9AE57
H	-24	SER	-	expression tag	UNP E9AE57
H	-23	GLY	-	expression tag	UNP E9AE57

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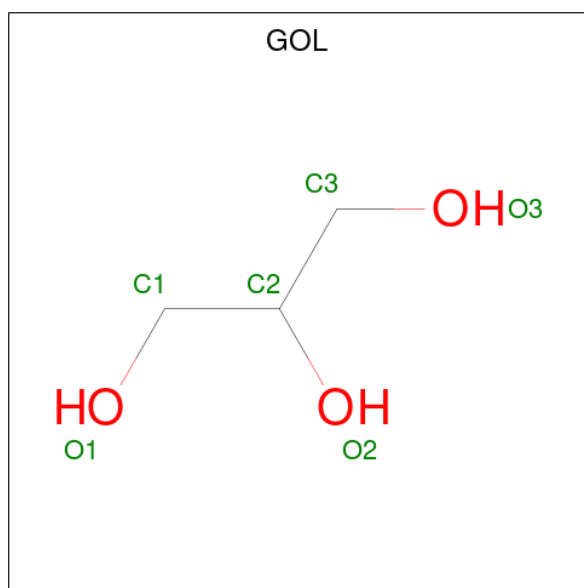
Chain	Residue	Modelled	Actual	Comment	Reference
H	-22	LEU	-	expression tag	UNP E9AE57
H	-21	VAL	-	expression tag	UNP E9AE57
H	-20	PRO	-	expression tag	UNP E9AE57
H	-19	ARG	-	expression tag	UNP E9AE57
H	-18	GLY	-	expression tag	UNP E9AE57
H	-17	SER	-	expression tag	UNP E9AE57
H	-16	HIS	-	expression tag	UNP E9AE57
H	-15	MET	-	expression tag	UNP E9AE57
H	-14	ALA	-	expression tag	UNP E9AE57
H	-13	SER	-	expression tag	UNP E9AE57
H	-12	MET	-	expression tag	UNP E9AE57
H	-11	THR	-	expression tag	UNP E9AE57
H	-10	GLY	-	expression tag	UNP E9AE57
H	-9	GLY	-	expression tag	UNP E9AE57
H	-8	GLN	-	expression tag	UNP E9AE57
H	-7	GLN	-	expression tag	UNP E9AE57
H	-6	MET	-	expression tag	UNP E9AE57
H	-5	GLY	-	expression tag	UNP E9AE57
H	-4	ARG	-	expression tag	UNP E9AE57
H	-3	GLY	-	expression tag	UNP E9AE57
H	-2	SER	-	expression tag	UNP E9AE57
H	-1	GLU	-	expression tag	UNP E9AE57
H	0	PHE	-	expression tag	UNP E9AE57

- Molecule 2 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe_4S_4) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	Fe	S	0	0
			8	4	4		
2	B	1	Total	Fe	S	0	0
			8	4	4		
2	C	1	Total	Fe	S	0	0
			8	4	4		
2	D	1	Total	Fe	S	0	0
			8	4	4		
2	E	1	Total	Fe	S	0	0
			8	4	4		
2	F	1	Total	Fe	S	0	0
			8	4	4		
2	G	1	Total	Fe	S	0	0
			8	4	4		
2	H	1	Total	Fe	S	0	0
			8	4	4		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			6	3	3		
3	E	1	Total	C	O	0	0
			6	3	3		

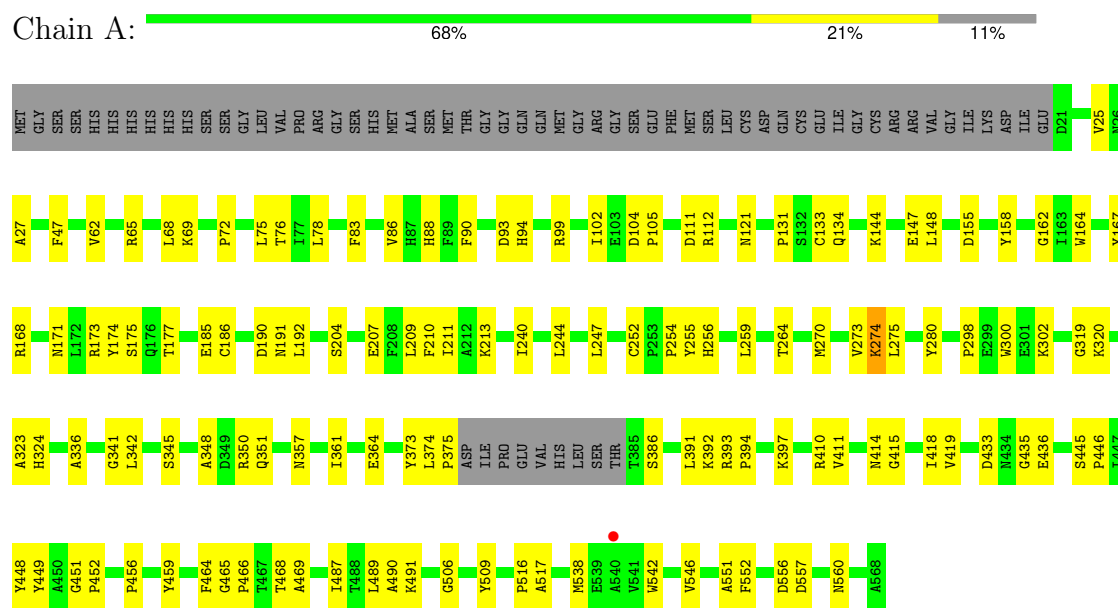
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total 1	O 1	0	0
4	D	1	Total 1	O 1	0	0
4	F	1	Total 1	O 1	0	0

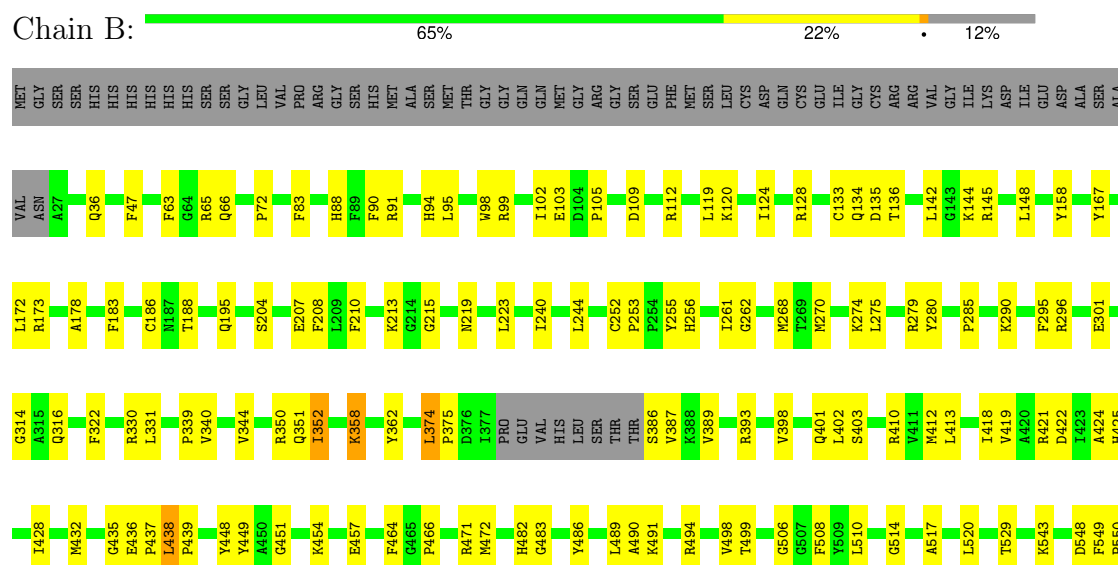
3 Residue-property plots

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: fumarate hydratase 2



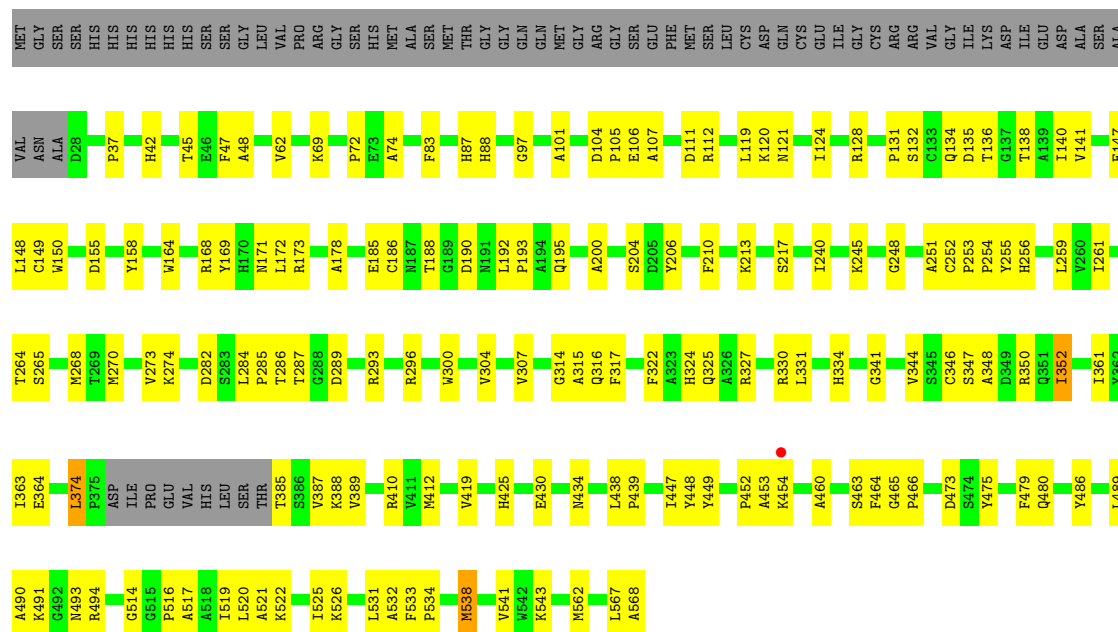
• Molecule 1: fumarate hydratase 2





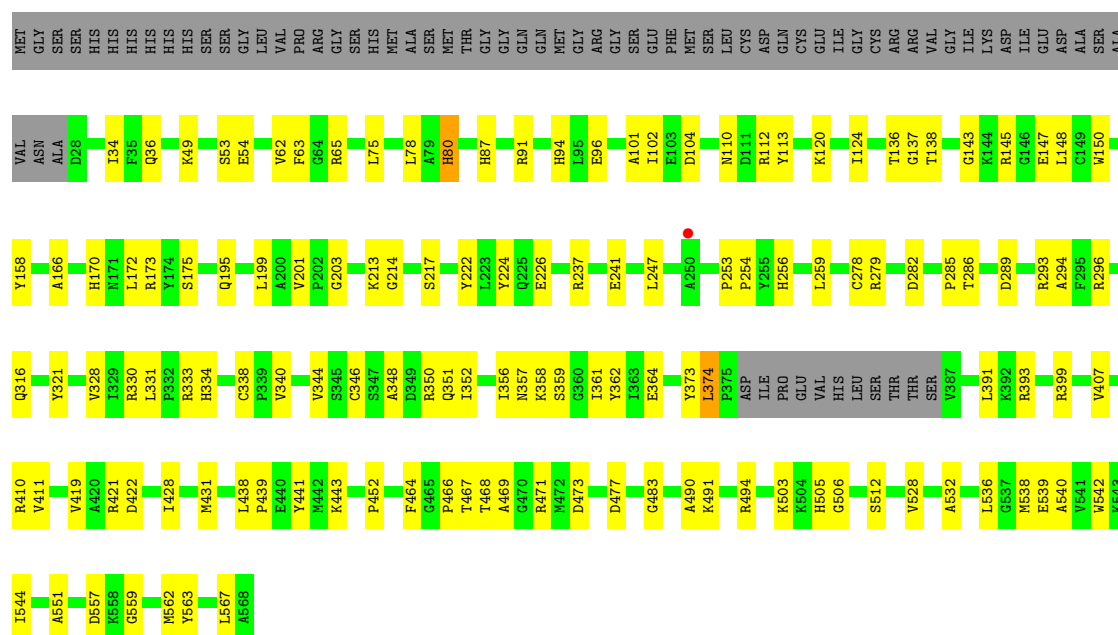
- Molecule 1: fumarate hydratase 2

Chain C: 61% 26% 12%



- Molecule 1: fumarate hydratase 2

Chain D: 66% 22% 12%



- Molecule 1: fumarate hydratase 2

Chain E:



K543	M412	A523	E185		VAL	MET
I544	L413	H524	E185		ASN	GLY
E545	N414		P193		ALA	SER
V546	G415	R327			ASP	HIS
F547					F29	HIS
D548	I418	R330	Y206			HIS
	V419	L331			Q36	HIS
F552	A420	P332	L209			HIS
I553	R421	R333			D39	HIS
	D422	H334	N219			HIS
D557		G335	K220		H42	SER
	P437	A336			H43	SER
T566	L438		Q225			GLY
L567	P439	P339			V62	LEU
A568		V340	L230		VAL	VAL
	S445	G341			R65	PRO
	P446		P233		O66	ARG
	I447	V344			A67	GLY
	Y448	S345	L236			SER
	Y449	C346			L75	HIS
		S347	I240			MET
	A453				H88	ALA
		Q351	K243			SER
	S463	I352	L244		R91	MET
			K245			THR
	T467	I361	T246		H94	GLY
	T468	V362				GLY
	A469	I363	A250		V98	GLN
	G470	E364				GLN
	R471		P253		I102	MET
		P375	ASP		E103	GLY
	Y475	ILE	Y255		D104	ARG
		PRO	H256		P105	GLY
	F479	GLU				SER
			L259		D111	GLU
	T488	VAL			R112	PHE
	L489	HIS	G262		Y113	MET
	A490	LEU				SER
		SER	T272		K120	LEU
	R494	THR			N121	CYS
			T280			ASP
	V498	S386	Y281		I124	GLN
		V387			A126	CYS
	G506		P285			GLU
						ILE
	P516	X392	T286		R128	GLY
	A517	R393				CYS
	A518	P394	R293		D136	ARG
		I395			T136	ARG
		D396	W300			VAL
	V528	K397			R144	GLY
					L145	ILE
	F533	R399	G312			LYS
	P534	Q400	T313		L148	ASP
		Q401	G314		C149	ILE
					W150	GLU
	M638		G318			ASP
	E539	G406	G319		D155	ALA
	A540	T409	K320			GLU
	V541	R410				ASP
			X321		L172	SER

- Molecule 1: fumarate hydratase 2

Chain F:



L531	A532	L536	G537	M538	E539	A540	V541	W542	V546	E547	D548	F549	P550	V554	D561	A568	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K491	R494	G507	S512	A517	THR	S386	L391	K392	R410	M412	L413	N414	G415	T416	L417	I418	V419	H425	Y441	S445	Y449	P452	P456	Y459	S463	F464	G465	T467	T468	R471	M472	D473	Y475	F479	Q480	S485	A490	K4
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- Molecule 1: fumarate hydratase 2

Chain G:



MET	GLY	SER	SER	SER	HIS	HIS	HIS	HIS	HIS	SER	SER	GLY	LEU	VAL	PRO	ARG	GLY	SER	HIS	SER	MET	MET	ALA	ALA	MET	THR	GLY	GLY	GLN	GLN	MET	GLY	ARG	ARG	SER	SER	GLY	GLU	PHE	GLU	MET	MET	SER	SER	LEU	LEU	CYS	ASP	GLN	GLN	CYS	GLY	CYS	ILE	ILE	GLY	CYS	LYS	ILE	GLU	ILE	GLY	ASP	ALA	SER	...
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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	159.09Å 66.07Å 238.09Å 90.00° 90.02° 90.00°	Depositor
Resolution (Å)	49.71 – 3.19 49.71 – 3.19	Depositor EDS
% Data completeness (in resolution range)	90.2 (49.71-3.19) 90.2 (49.71-3.19)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	714.55 (at 3.19Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
R, R_{free}	0.245 , 0.289 0.245 , 0.289	Depositor DCC
R_{free} test set	3753 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å ²)	65.6	Xtriage
Anisotropy	0.543	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 136.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.22$	Xtriage
Estimated twinning fraction	0.398 for h,-k,-l	Xtriage
Reported twinning fraction	0.500 for h,-k,-l	Depositor
Outliers	0 of 75107 reflections	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	31246	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 17.75% 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: SF4, GOL

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	0.30	0/4024	0.74	2/5476 (0.0%)
1	B	0.33	0/4029	0.78	0/5489
1	C	0.27	0/3911	0.72	0/5329
1	D	0.26	0/3966	0.69	0/5399
1	E	0.31	1/3952 (0.0%)	0.70	0/5380
1	F	0.29	0/4076	0.68	0/5542
1	G	0.29	0/3965	0.71	0/5401
1	H	0.30	0/3984	0.73	0/5432
All	All	0.29	1/31907 (0.0%)	0.72	2/43448 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	144	LYS	C-O	-5.29	1.17	1.23

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	433	ASP	CA-C-N	-5.03	114.31	122.65
1	A	433	ASP	C-N-CA	-5.03	114.31	122.65

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	3933	0	3563	92	0
1	B	3933	0	3571	100	0
1	C	3822	0	3400	125	0
1	D	3874	0	3531	96	0
1	E	3863	0	3537	109	0
1	F	3981	0	3721	117	0
1	G	3873	0	3489	111	0
1	H	3888	0	3506	126	0
2	A	8	0	0	1	0
2	B	8	0	0	2	0
2	C	8	0	0	5	0
2	D	8	0	0	0	0
2	E	8	0	0	1	0
2	F	8	0	0	3	0
2	G	8	0	0	0	0
2	H	8	0	0	0	0
3	B	6	0	8	2	0
3	E	6	0	8	2	0
4	A	1	0	0	0	0
4	D	1	0	0	0	0
4	F	1	0	0	0	0
All	All	31246	0	28334	815	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 14.

All (815) 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:G:494:ARG:HD2	1:G:509:TYR:CD1	2.00	0.96
1:B:421:ARG:HH11	1:B:451:GLY:HA3	1.32	0.95
1:G:494:ARG:HD2	1:G:509:TYR:CE1	2.07	0.89
1:F:412:MET:HE3	1:F:548:ASP:HB3	1.53	0.89
1:E:399:ARG:HD2	1:E:506:GLY:O	1.73	0.89
1:C:134:GLN:NE2	2:C:601:SF4:S4	2.48	0.87
1:F:72:PRO:HB2	1:F:158:TYR:HD2	1.42	0.85
1:D:102:ILE:HA	1:D:112:ARG:HG2	1.58	0.83
1:F:217:SER:HB3	1:F:346:CYS:HB3	1.61	0.83
1:B:65:ARG:HE	1:B:148:LEU:HD11	1.42	0.82
1:H:412:MET:HE3	1:H:548:ASP:HB3	1.60	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:473:ASP:HA	1:C:494:ARG:HH21	1.45	0.81
1:B:136:THR:HG21	1:B:172:LEU:HB3	1.61	0.80
1:H:72:PRO:HB2	1:H:158:TYR:HD2	1.46	0.80
1:F:532:ALA:HB3	1:F:541:VAL:HB	1.65	0.78
1:F:418:ILE:HG12	1:F:445:SER:HB2	1.65	0.77
1:B:506:GLY:HA2	1:B:557:ASP:HB2	1.66	0.77
1:A:336:ALA:HB2	3:B:602:GOL:H31	1.67	0.76
1:E:399:ARG:CD	1:E:506:GLY:O	2.33	0.76
1:C:245:LYS:HD2	1:C:567:LEU:HA	1.68	0.76
1:E:136:THR:HG21	1:E:172:LEU:HB3	1.67	0.75
1:A:104:ASP:O	1:A:112:ARG:NH2	2.20	0.75
1:G:219:ASN:HD21	1:H:337:SER:HA	1.51	0.75
1:G:93:ASP:OD2	1:G:279:ARG:NE	2.20	0.75
1:C:140:ILE:HA	1:C:195:GLN:HB3	1.69	0.74
1:H:88:HIS:ND1	1:H:125:ALA:O	2.19	0.74
1:H:275:LEU:HD22	1:H:280:TYR:HD2	1.53	0.74
1:B:148:LEU:HB2	1:B:204:SER:HB3	1.70	0.74
1:A:72:PRO:HB2	1:A:158:TYR:HD2	1.53	0.74
1:B:102:ILE:O	1:B:112:ARG:NH2	2.22	0.73
1:D:358:LYS:H	1:D:358:LYS:HD3	1.52	0.73
1:B:393:ARG:NH2	1:B:401:GLN:OE1	2.21	0.72
1:F:109:ASP:OD1	1:F:112:ARG:NH1	2.22	0.72
1:A:469:ALA:HB2	1:A:490:ALA:HB3	1.71	0.72
1:G:282:ASP:O	1:G:296:ARG:NH1	2.23	0.72
1:A:336:ALA:O	1:B:219:ASN:ND2	2.23	0.72
1:F:143:GLY:O	1:F:199:LEU:N	2.20	0.72
1:H:72:PRO:HB2	1:H:158:TYR:CD2	2.24	0.72
1:D:49:LYS:NZ	1:D:359:SER:O	2.22	0.71
1:D:138:THR:N	1:D:213:LYS:O	2.20	0.71
1:E:102:ILE:O	1:E:112:ARG:NH1	2.24	0.71
1:C:47:PHE:HA	1:C:364:GLU:HA	1.72	0.70
1:F:62:VAL:HG11	1:F:150:TRP:HH2	1.55	0.70
1:D:399:ARG:NH2	1:D:506:GLY:O	2.23	0.70
1:D:330:ARG:NH1	1:D:331:LEU:O	2.23	0.70
1:B:213:LYS:NZ	1:B:344:VAL:O	2.25	0.70
1:B:489:LEU:HA	1:B:510:LEU:O	1.92	0.69
1:D:473:ASP:OD1	1:D:494:ARG:NH2	2.26	0.69
1:G:494:ARG:HH11	1:G:509:TYR:HD1	1.40	0.69
1:B:99:ARG:HG3	1:B:375:PRO:HG3	1.74	0.69
1:H:425:HIS:HD2	1:H:479:PHE:HE2	1.41	0.69
1:D:136:THR:HG21	1:D:172:LEU:HB3	1.74	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:395:ILE:O	1:E:397:LYS:N	2.21	0.69
1:B:173:ARG:NH2	3:B:602:GOL:O3	2.26	0.69
1:F:72:PRO:HB2	1:F:158:TYR:CD2	2.28	0.69
1:H:330:ARG:NH1	1:H:331:LEU:O	2.26	0.68
1:D:148:LEU:N	1:D:203:GLY:O	2.24	0.68
1:A:435:GLY:O	1:A:436:GLU:HG3	1.93	0.68
1:H:107:ALA:O	1:H:112:ARG:NH2	2.27	0.68
1:D:422:ASP:OD2	1:D:471:ARG:NH1	2.26	0.68
1:F:467:THR:OG1	1:F:471:ARG:NH1	2.27	0.68
1:C:48:ALA:N	1:C:363:ILE:O	2.18	0.67
1:E:65:ARG:HH22	1:F:152:GLY:HA2	1.60	0.67
1:A:252:CYS:SG	1:A:491:LYS:NZ	2.63	0.67
1:C:164:TRP:HB2	1:C:192:LEU:HD11	1.77	0.66
1:G:134:GLN:OE1	1:G:173:ARG:NH1	2.29	0.66
1:G:135:ASP:OD1	1:H:334:HIS:NE2	2.27	0.66
1:B:428:ILE:HG12	1:B:439:PRO:HG2	1.77	0.66
1:A:72:PRO:HB2	1:A:158:TYR:CD2	2.30	0.66
1:A:134:GLN:NE2	2:A:601:SF4:S1	2.68	0.66
1:E:421:ARG:HG3	1:E:539:GLU:HG2	1.77	0.66
1:A:167:TYR:HB3	1:A:174:TYR:HE1	1.59	0.65
1:B:253:PRO:HG3	1:B:255:TYR:CZ	2.31	0.65
1:E:135:ASP:OD1	1:F:334:HIS:NE2	2.27	0.65
1:F:134:GLN:NE2	2:F:601:SF4:S2	2.68	0.65
1:E:287:THR:OG1	1:E:293:ARG:NH2	2.30	0.65
1:A:133:CYS:HB3	1:A:350:ARG:HH22	1.62	0.65
1:G:88:HIS:ND1	1:G:125:ALA:O	2.30	0.64
1:G:247:LEU:HD11	1:G:342:LEU:HD21	1.77	0.64
1:H:178:ALA:N	1:H:186:CYS:O	2.28	0.64
1:H:217:SER:HB3	1:H:346:CYS:HB3	1.79	0.64
1:F:97:GLY:HA3	1:F:324:HIS:HD2	1.63	0.64
1:A:175:SER:N	1:A:191:ASN:OD1	2.24	0.64
1:D:407:VAL:HB	1:D:562:MET:HB3	1.79	0.64
1:A:336:ALA:HB3	1:B:215:GLY:HA3	1.80	0.64
1:E:494:ARG:HD3	1:E:498:VAL:HG11	1.80	0.64
1:C:425:HIS:HB3	1:C:475:TYR:CD1	2.33	0.64
1:D:147:GLU:N	1:D:201:VAL:O	2.31	0.64
1:A:102:ILE:HA	1:A:112:ARG:HG2	1.80	0.64
1:C:270:MET:HA	1:C:273:VAL:HG12	1.79	0.64
1:G:111:ASP:OD2	1:G:320:LYS:N	2.21	0.63
1:D:213:LYS:NZ	1:D:344:VAL:O	2.31	0.63
1:F:333:ARG:NH1	1:F:337:SER:O	2.29	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:254:PRO:HB2	1:A:324:HIS:CD2	2.33	0.63
1:A:65:ARG:NH1	1:A:147:GLU:OE2	2.32	0.63
1:C:330:ARG:NH1	1:C:331:LEU:O	2.31	0.63
1:F:245:LYS:HZ1	1:F:568:ALA:N	1.97	0.63
1:F:142:LEU:HD11	1:F:199:LEU:HD11	1.81	0.63
1:G:511:GLY:N	1:G:552:PHE:O	2.31	0.63
1:H:277:SER:O	1:H:279:ARG:NH1	2.31	0.63
1:C:111:ASP:HA	1:C:315:ALA:HB3	1.81	0.62
1:D:75:LEU:HB2	1:D:158:TYR:HB3	1.82	0.62
1:F:413:LEU:O	1:F:548:ASP:N	2.32	0.62
1:G:469:ALA:CB	1:G:509:TYR:CE1	2.81	0.62
1:A:394:PRO:HD2	1:A:397:LYS:HD2	1.81	0.62
1:C:287:THR:OG1	1:C:293:ARG:NH2	2.31	0.62
1:B:386:SER:OG	1:B:410:ARG:NH2	2.32	0.62
1:B:529:THR:O	1:B:543:LYS:N	2.22	0.62
1:E:422:ASP:OD2	1:E:471:ARG:NH1	2.31	0.62
1:E:150:TRP:HB3	1:F:65:ARG:HH12	1.64	0.62
1:G:486:TYR:O	1:G:508:PHE:N	2.32	0.62
1:A:164:TRP:HB2	1:A:192:LEU:HD11	1.81	0.62
1:F:473:ASP:OD1	1:F:494:ARG:NE	2.31	0.62
1:F:449:TYR:HB2	1:F:490:ALA:HB2	1.81	0.61
1:H:418:ILE:HG12	1:H:445:SER:HB2	1.81	0.61
1:G:256:HIS:O	1:G:344:VAL:HA	1.99	0.61
1:H:480:GLN:NE2	1:H:485:SER:O	2.32	0.61
1:A:144:LYS:HE3	1:A:207:GLU:HB2	1.82	0.61
1:H:171:ASN:HA	1:H:454:LYS:HB2	1.80	0.61
1:A:456:PRO:HB2	1:A:459:TYR:CD2	2.35	0.61
1:E:385:THR:O	1:E:387:VAL:N	2.33	0.61
1:D:411:VAL:N	1:D:551:ALA:O	2.28	0.61
1:E:488:THR:OG1	1:E:494:ARG:NH1	2.34	0.61
1:A:174:TYR:HD1	1:A:191:ASN:HB2	1.65	0.60
1:A:391:LEU:N	1:A:414:ASN:O	2.22	0.60
1:C:74:ALA:HB1	1:C:361:ILE:HD11	1.83	0.60
1:E:422:ASP:HB3	1:F:228:LYS:HB3	1.81	0.60
1:F:173:ARG:HD2	1:F:538:MET:HB3	1.83	0.60
1:D:490:ALA:O	1:D:512:SER:N	2.34	0.60
1:G:274:LYS:NZ	1:H:267:GLU:OE2	2.34	0.60
1:G:399:ARG:HD2	1:G:486:TYR:HB3	1.83	0.60
1:D:282:ASP:O	1:D:296:ARG:NH1	2.27	0.60
1:H:415:GLY:N	1:H:546:VAL:O	2.28	0.60
1:C:425:HIS:HB3	1:C:475:TYR:CE1	2.37	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:243:LYS:O	1:E:246:THR:OG1	2.18	0.60
1:C:193:PRO:HB2	1:D:34:ILE:HD11	1.83	0.60
1:G:114:VAL:HG21	1:G:315:ALA:O	2.01	0.60
1:H:254:PRO:HG3	1:H:315:ALA:HB1	1.82	0.60
1:D:137:GLY:HA2	1:D:214:GLY:HA2	1.83	0.59
1:H:257:ILE:O	1:H:326:ALA:HA	2.02	0.59
1:C:87:HIS:O	1:C:132:SER:N	2.26	0.59
1:F:148:LEU:HB2	1:F:204:SER:HB3	1.84	0.59
1:H:449:TYR:HB3	1:H:467:THR:HG22	1.84	0.59
1:H:176:GLN:HG2	1:H:191:ASN:HA	1.85	0.59
1:D:419:VAL:HB	1:D:542:TRP:HB2	1.83	0.59
1:F:100:ARG:HG2	1:F:321:TYR:OH	2.03	0.59
1:G:396:ASP:HA	1:G:399:ARG:HD3	1.83	0.59
1:F:134:GLN:HB3	1:F:465:GLY:H	1.67	0.59
1:C:72:PRO:HB2	1:C:158:TYR:HD2	1.68	0.59
1:C:136:THR:HG21	1:C:172:LEU:HB3	1.85	0.59
1:D:91:ARG:HD2	1:D:279:ARG:HH21	1.68	0.58
1:G:256:HIS:CD2	1:G:345:SER:HB3	2.38	0.58
1:H:416:THR:HG21	1:H:543:LYS:HE2	1.84	0.58
1:E:135:ASP:HB2	3:E:602:GOL:H2	1.84	0.58
1:F:254:PRO:HB2	1:F:324:HIS:CE1	2.38	0.58
1:B:483:GLY:HA2	1:B:486:TYR:OH	2.03	0.58
1:E:336:ALA:O	1:F:219:ASN:ND2	2.28	0.58
1:H:506:GLY:HA2	1:H:557:ASP:HB2	1.86	0.58
1:E:220:LYS:HE2	1:E:250:ALA:HB3	1.85	0.58
1:C:248:GLY:HA3	1:C:493:ASN:HD21	1.68	0.58
1:E:104:ASP:O	1:E:112:ARG:NH2	2.36	0.57
1:E:314:GLY:HA3	1:E:318:GLY:C	2.29	0.57
1:E:393:ARG:NH2	1:E:401:GLN:OE1	2.37	0.57
1:C:388:LYS:HA	1:C:412:MET:HB2	1.85	0.57
1:A:452:PRO:HB3	1:A:464:PHE:CD2	2.40	0.57
1:G:488:THR:O	1:G:510:LEU:N	2.31	0.57
1:B:438:LEU:HD21	1:B:482:HIS:O	2.05	0.57
1:E:98:TRP:CD1	1:E:324:HIS:HE2	2.21	0.57
1:E:245:LYS:NZ	1:E:568:ALA:HB2	2.20	0.57
1:G:254:PRO:HB2	1:G:324:HIS:CE1	2.40	0.57
1:D:253:PRO:HD2	1:D:316:GLN:HG2	1.85	0.57
1:E:67:ALA:HA	1:E:148:LEU:HB3	1.87	0.57
1:F:490:ALA:O	1:F:512:SER:N	2.37	0.57
1:G:480:GLN:NE2	1:G:507:GLY:HA3	2.19	0.57
1:D:431:MET:HE1	1:D:439:PRO:HG3	1.85	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:268:MET:HE1	1:C:331:LEU:HD13	1.87	0.56
1:D:506:GLY:HA2	1:D:557:ASP:HB2	1.85	0.56
1:B:358:LYS:H	1:B:358:LYS:HD3	1.70	0.56
1:C:119:LEU:HB3	1:C:374:LEU:HD22	1.87	0.56
1:C:449:TYR:HB2	1:C:490:ALA:HB2	1.88	0.56
1:F:87:HIS:HE1	1:F:136:THR:HB	1.71	0.56
1:F:135[B]:ASP:O	1:F:350:ARG:NH1	2.34	0.56
1:G:134:GLN:HB3	1:G:517:ALA:HB1	1.86	0.56
1:G:442:MET:HE2	1:G:479:PHE:CE2	2.40	0.56
1:H:88:HIS:CE1	1:H:128:ARG:HA	2.41	0.56
1:C:452:PRO:HB3	1:C:464:PHE:CG	2.39	0.56
1:F:135[A]:ASP:O	1:F:350:ARG:NH1	2.34	0.56
1:G:111:ASP:OD1	1:G:319:GLY:HA3	2.06	0.56
1:G:145:ARG:CZ	1:H:200:ALA:HB1	2.35	0.56
1:H:83:PHE:HE1	1:H:210:PHE:HB3	1.69	0.56
1:E:185:GLU:OE2	1:F:330:ARG:NH2	2.39	0.56
1:C:251:ALA:HA	2:C:601:SF4:S1	2.45	0.56
1:F:155:ASP:OD1	1:F:155:ASP:N	2.38	0.56
1:H:369:ASN:O	1:H:372:GLN:HG2	2.06	0.56
1:B:253:PRO:HD2	1:B:316:GLN:HG2	1.86	0.56
1:D:49:LYS:NZ	1:D:54:GLU:OE1	2.37	0.56
1:D:503:LYS:HD2	1:D:559:GLY:HA3	1.88	0.56
1:E:120:LYS:O	1:E:124:ILE:HG12	2.06	0.56
1:F:78:LEU:C	1:F:80:HIS:H	2.14	0.56
1:G:385:THR:O	1:G:410:ARG:NH1	2.39	0.56
1:G:399:ARG:HD2	1:G:486:TYR:CB	2.36	0.56
1:A:213:LYS:HE3	1:A:274:LYS:HD2	1.87	0.55
1:B:144:LYS:HE3	1:B:207:GLU:HB2	1.87	0.55
1:A:490:ALA:H	1:A:509:TYR:HE1	1.54	0.55
1:B:120:LYS:O	1:B:124:ILE:HG12	2.06	0.55
1:C:147:GLU:HA	1:D:145:ARG:NH2	2.21	0.55
1:D:54:GLU:HB3	1:D:359:SER:O	2.06	0.55
1:B:413:LEU:HB2	1:B:549:PHE:HB3	1.88	0.55
1:C:178:ALA:N	1:C:186:CYS:O	2.37	0.55
1:A:134:GLN:HB3	1:A:465:GLY:H	1.72	0.55
1:C:213:LYS:HE2	1:C:274:LYS:HE2	1.88	0.55
1:E:397:LYS:O	1:E:401:GLN:HG3	2.06	0.55
1:G:142:LEU:O	1:G:208:PHE:HA	2.06	0.55
1:B:314:GLY:O	1:B:322:PHE:HD1	1.89	0.55
1:C:97:GLY:HA3	1:C:324:HIS:ND1	2.21	0.55
1:C:134:GLN:HB3	1:C:465:GLY:H	1.71	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:528:VAL:HG22	1:D:544:ILE:HG22	1.87	0.55
1:B:167:TYR:O	1:B:454:LYS:NZ	2.31	0.55
1:B:428:ILE:HG12	1:B:439:PRO:CG	2.37	0.55
1:F:177:THR:HG22	1:F:185:GLU:HB3	1.89	0.55
1:C:453:ALA:N	1:C:463:SER:O	2.33	0.55
1:G:33:ALA:O	1:G:36:GLN:NE2	2.28	0.55
1:D:62:VAL:HG22	1:D:63:PHE:CD2	2.43	0.55
1:E:395:ILE:C	1:E:397:LYS:H	2.13	0.55
1:C:135:ASP:O	1:C:350:ARG:NH2	2.36	0.54
1:E:475:TYR:HB3	1:E:479:PHE:CE2	2.43	0.54
1:G:253:PRO:HG3	1:G:255:TYR:CZ	2.41	0.54
1:A:411:VAL:N	1:A:551:ALA:O	2.31	0.54
1:C:217:SER:HB3	1:C:346:CYS:HB3	1.89	0.54
1:C:521:ALA:HA	1:C:525:ILE:HB	1.90	0.54
1:A:68:LEU:HD11	1:A:357:ASN:C	2.32	0.54
1:E:272:THR:HG23	1:E:281:TYR:HE2	1.72	0.54
1:H:62:VAL:HG22	1:H:63:PHE:CD2	2.43	0.54
1:H:508:PHE:HD1	1:H:556:ASP:HA	1.72	0.54
1:B:252:CYS:HB2	2:B:601:SF4:S2	2.48	0.54
1:D:120:LYS:O	1:D:124:ILE:HG12	2.07	0.54
1:C:149:CYS:SG	1:C:206:TYR:HB2	2.47	0.54
1:E:65:ARG:NH2	1:F:152:GLY:HA2	2.23	0.54
1:E:391:LEU:HD22	1:E:446:PRO:HG2	1.90	0.54
1:H:155:ASP:N	1:H:155:ASP:OD1	2.40	0.54
1:H:213:LYS:HE2	1:H:274:LYS:HE2	1.90	0.54
1:H:254:PRO:HB2	1:H:324:HIS:NE2	2.23	0.54
1:D:53:SER:OG	1:D:361:ILE:N	2.32	0.54
1:H:495:SER:OG	1:H:497:GLN:HG2	2.07	0.54
1:D:49:LYS:HB2	1:D:362:TYR:CE2	2.42	0.54
1:E:254:PRO:HB2	1:E:324:HIS:CD2	2.43	0.54
1:H:469:ALA:HB2	1:H:490:ALA:HB3	1.89	0.53
1:C:178:ALA:HB3	1:C:186:CYS:SG	2.48	0.53
1:E:225:GLN:HG3	1:F:222:TYR:HE1	1.73	0.53
1:B:135:ASP:O	1:B:350:ARG:NH1	2.41	0.53
1:G:104:ASP:O	1:G:112:ARG:NH2	2.41	0.53
1:A:556:ASP:OD2	1:A:560:ASN:ND2	2.37	0.53
1:C:135:ASP:HA	1:C:173:ARG:CZ	2.38	0.53
1:D:452:PRO:HB3	1:D:464:PHE:CE2	2.44	0.53
1:A:83:PHE:CE1	1:A:210:PHE:HB3	2.43	0.53
1:D:36:GLN:O	1:D:285:PRO:HD3	2.08	0.53
1:D:469:ALA:O	1:D:494:ARG:NH2	2.42	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:111:ASP:OD2	1:E:320:LYS:N	2.42	0.53
1:E:287:THR:O	1:E:293:ARG:NE	2.37	0.53
1:A:174:TYR:HA	1:A:191:ASN:HB2	1.90	0.53
1:C:171:ASN:HB3	1:C:454:LYS:O	2.09	0.53
1:B:72:PRO:HB2	1:B:158:TYR:CG	2.44	0.53
1:D:217:SER:HB3	1:D:346:CYS:HB3	1.90	0.53
1:D:226:GLU:C	1:D:338:CYS:HB3	2.34	0.53
1:G:147:GLU:OE2	1:H:152:GLY:N	2.35	0.53
1:D:253:PRO:O	1:D:316:GLN:NE2	2.34	0.52
1:G:469:ALA:HB2	1:G:509:TYR:CE1	2.44	0.52
1:D:237:ARG:O	1:D:241:GLU:HG3	2.09	0.52
1:E:121:ASN:ND2	1:E:516:PRO:HA	2.23	0.52
1:F:62:VAL:HG22	1:F:63:PHE:CD2	2.45	0.52
1:F:480:GLN:NE2	1:F:507:GLY:HA3	2.25	0.52
1:C:282:ASP:CG	1:C:327:ARG:HE	2.15	0.52
1:D:101:ALA:HA	1:D:321:TYR:CZ	2.43	0.52
1:E:449:TYR:HB3	1:E:467:THR:HG22	1.91	0.52
1:H:145:ARG:HB2	1:H:206:TYR:CZ	2.43	0.52
1:H:416:THR:OG1	1:H:545:GLU:OE1	2.22	0.52
1:B:418:ILE:O	1:B:448:TYR:N	2.28	0.52
1:C:141:VAL:N	1:C:195:GLN:O	2.29	0.52
1:D:222:TYR:HE2	1:D:247:LEU:HD23	1.75	0.52
1:D:285:PRO:O	1:D:296:ARG:N	2.39	0.52
1:H:413:LEU:O	1:H:548:ASP:N	2.43	0.52
1:C:253:PRO:HD2	1:C:316:GLN:HG2	1.92	0.52
1:C:134:GLN:HG3	1:C:465:GLY:O	2.10	0.52
1:C:452:PRO:HD2	1:C:538:MET:HB2	1.91	0.52
1:H:83:PHE:CE1	1:H:210:PHE:HB3	2.45	0.52
1:A:186:CYS:SG	1:B:290:LYS:HA	2.49	0.52
1:C:452:PRO:HB3	1:C:464:PHE:CD2	2.44	0.52
1:D:469:ALA:HB2	1:D:490:ALA:HB3	1.92	0.52
1:A:105:PRO:HA	1:A:112:ARG:HH22	1.75	0.51
1:C:264:THR:O	1:D:138:THR:HG21	2.10	0.51
1:E:245:LYS:HZ1	1:E:568:ALA:HB2	1.75	0.51
1:G:93:ASP:OD2	1:G:325:GLN:NE2	2.41	0.51
1:H:393:ARG:NH1	1:H:401:GLN:HB2	2.26	0.51
1:B:90:PHE:CD2	1:B:94:HIS:CD2	2.98	0.51
1:B:210:PHE:HB2	1:B:352:ILE:HG22	1.93	0.51
1:B:464:PHE:H	1:B:517:ALA:HB1	1.75	0.51
1:H:295:PHE:CE1	1:H:330:ARG:HB3	2.45	0.51
1:C:347:SER:HB2	2:C:601:SF4:S3	2.49	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:387:VAL:O	1:C:412:MET:N	2.29	0.51
1:B:91:ARG:HD2	1:B:279:ARG:HH21	1.76	0.51
1:C:150:TRP:HB3	1:D:65:ARG:HH12	1.75	0.51
1:G:469:ALA:CB	1:G:509:TYR:HE1	2.24	0.51
1:B:109:ASP:OD2	1:F:531:LEU:HA	2.10	0.51
1:B:135:ASP:O	1:B:350:ARG:NH2	2.44	0.51
1:C:240:ILE:HG22	1:C:307:VAL:HG11	1.92	0.51
1:F:419:VAL:O	1:F:541:VAL:HA	2.09	0.51
1:B:464:PHE:N	1:B:517:ALA:HB1	2.25	0.51
1:E:193:PRO:HD3	1:F:291:TYR:O	2.10	0.51
1:H:145:ARG:NH1	1:H:155:ASP:OD2	2.42	0.51
1:G:442:MET:HE2	1:G:479:PHE:CD2	2.46	0.51
1:G:90:PHE:CD2	1:G:94:HIS:HD2	2.29	0.51
1:G:167:TYR:HB3	1:G:174:TYR:HE1	1.76	0.51
1:H:136:THR:HB	1:H:167:TYR:HE2	1.76	0.51
1:H:411:VAL:O	1:H:550:PRO:HA	2.11	0.51
1:A:88:HIS:CE1	1:A:131:PRO:HA	2.46	0.50
1:E:336:ALA:O	1:F:471:ARG:NH2	2.43	0.50
1:F:101:ALA:HA	1:F:321:TYR:CE2	2.46	0.50
1:H:74:ALA:HB1	1:H:361:ILE:HD11	1.93	0.50
1:H:419:VAL:O	1:H:541:VAL:HA	2.11	0.50
1:A:255:TYR:H	1:A:323:ALA:HA	1.76	0.50
1:C:171:ASN:O	1:C:453:ALA:HB1	2.10	0.50
1:E:414:ASN:OD1	1:E:548:ASP:N	2.32	0.50
1:A:264:THR:OG1	1:B:195:GLN:HB2	2.11	0.50
1:F:253:PRO:HG2	1:F:316:GLN:HG2	1.92	0.50
1:H:275:LEU:HD22	1:H:280:TYR:CD2	2.39	0.50
1:H:285:PRO:O	1:H:296:ARG:N	2.42	0.50
1:A:466:PRO:O	1:A:491:LYS:HE2	2.12	0.50
1:B:393:ARG:NH2	1:B:398:VAL:HA	2.27	0.50
1:A:155:ASP:OD1	1:A:155:ASP:N	2.44	0.50
1:B:213:LYS:HE2	1:B:274:LYS:HE2	1.92	0.50
1:B:275:LEU:HB3	1:B:280:TYR:HB3	1.93	0.50
1:C:300:TRP:O	1:C:304:VAL:HG23	2.12	0.50
1:E:330:ARG:NH1	1:E:331:LEU:O	2.43	0.50
1:F:257:ILE:O	1:F:326:ALA:HA	2.11	0.50
1:G:135:ASP:O	1:G:350:ARG:NH1	2.40	0.50
1:A:133:CYS:HB3	1:A:350:ARG:NH2	2.26	0.50
1:C:121:ASN:OD1	1:C:519:ILE:N	2.31	0.50
1:E:113:TYR:HE2	1:E:552:PHE:HZ	1.60	0.50
1:E:386:SER:HA	1:E:412:MET:HG3	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:134:GLN:OE1	1:F:517:ALA:HB2	2.12	0.50
1:C:107:ALA:HB3	1:C:112:ARG:HE	1.76	0.50
1:D:466:PRO:HG2	1:D:491:LYS:HG2	1.94	0.50
1:E:256:HIS:HB2	1:E:345:SER:HB3	1.92	0.50
1:F:240:ILE:O	1:F:244:LEU:HG	2.12	0.50
1:F:411:VAL:O	1:F:550:PRO:HA	2.12	0.50
1:G:79:ALA:HB3	1:G:162:GLY:HA3	1.94	0.50
1:H:414:ASN:HA	1:H:547:GLU:HA	1.92	0.50
1:B:88:HIS:CD2	1:B:128:ARG:HA	2.47	0.50
1:F:252:CYS:O	1:F:347:SER:OG	2.23	0.50
1:A:419:VAL:HG11	1:A:464:PHE:HZ	1.77	0.49
1:A:419:VAL:HB	1:A:542:TRP:HB2	1.94	0.49
1:C:178:ALA:HB2	1:C:188:THR:HG22	1.93	0.49
1:F:386:SER:O	1:F:386:SER:OG	2.30	0.49
1:G:563:TYR:HB3	1:G:567:LEU:HD12	1.93	0.49
1:F:300:TRP:O	1:F:304:VAL:HG23	2.11	0.49
1:H:448:TYR:HE1	1:H:466:PRO:HB3	1.77	0.49
1:A:211:ILE:HG22	1:A:213:LYS:HG2	1.94	0.49
1:C:520:LEU:O	1:C:525:ILE:N	2.40	0.49
1:D:563:TYR:HB3	1:D:567:LEU:HD12	1.94	0.49
1:E:121:ASN:OD1	1:E:518:ALA:HB3	2.12	0.49
1:F:80:HIS:HE1	1:F:170:HIS:CE1	2.31	0.49
1:G:141:VAL:HG11	1:G:159:LEU:HD13	1.94	0.49
1:B:99:ARG:O	1:B:103:GLU:HG3	2.12	0.49
1:C:466:PRO:HG2	1:C:491:LYS:HG2	1.93	0.49
1:H:449:TYR:O	1:H:467:THR:N	2.43	0.49
1:A:69:LYS:HD2	1:B:63:PHE:CE2	2.47	0.49
1:C:352:ILE:HG23	1:C:364:GLU:HB3	1.94	0.49
1:C:419:VAL:HA	1:C:448:TYR:O	2.13	0.49
1:H:412:MET:HB3	1:H:548:ASP:HA	1.94	0.49
1:A:121:ASN:CG	1:A:516:PRO:HA	2.36	0.49
1:A:164:TRP:O	1:A:168:ARG:N	2.35	0.49
1:C:48:ALA:O	1:C:363:ILE:HG22	2.13	0.49
1:D:87:HIS:HA	1:D:350:ARG:HD2	1.94	0.49
1:D:443:LYS:HG2	1:D:483:GLY:O	2.13	0.49
1:A:247:LEU:HD11	1:A:342:LEU:HD21	1.94	0.49
1:A:456:PRO:HB2	1:A:459:TYR:HD2	1.74	0.49
1:B:295:PHE:CZ	1:B:330:ARG:HB3	2.48	0.49
1:C:101:ALA:O	1:C:112:ARG:HG2	2.13	0.49
1:C:491:LYS:HE2	2:C:601:SF4:S4	2.52	0.49
1:E:65:ARG:HH12	1:F:152:GLY:HA2	1.78	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:399:ARG:HD3	1:E:506:GLY:O	2.12	0.49
1:A:148:LEU:HB2	1:A:204:SER:HB3	1.95	0.48
1:E:39:ASP:OD2	1:E:42:HIS:HA	2.13	0.48
1:G:219:ASN:HA	1:H:225:GLN:CD	2.37	0.48
1:G:469:ALA:HB2	1:G:509:TYR:HE1	1.77	0.48
1:B:554:VAL:O	1:B:562:MET:N	2.32	0.48
1:G:480:GLN:NE2	1:G:502:CYS:SG	2.80	0.48
1:H:140:ILE:HG12	1:H:195:GLN:HB3	1.95	0.48
1:H:167:TYR:HA	1:H:172:LEU:HD12	1.94	0.48
1:H:240:ILE:HG22	1:H:307:VAL:HG11	1.95	0.48
1:B:358:LYS:H	1:B:358:LYS:CD	2.26	0.48
1:D:104:ASP:O	1:D:112:ARG:NH2	2.45	0.48
1:F:250:ALA:O	1:F:491:LYS:NZ	2.43	0.48
1:G:47:PHE:HB3	1:G:362:TYR:HB3	1.94	0.48
1:G:407:VAL:HB	1:G:562:MET:HB3	1.95	0.48
1:G:499:THR:HG22	1:G:559:GLY:HA2	1.96	0.48
1:H:466:PRO:HG2	1:H:491:LYS:HG2	1.94	0.48
1:B:119:LEU:HD13	1:B:374:LEU:HD13	1.95	0.48
1:G:142:LEU:HD12	1:G:197:ASP:HB3	1.95	0.48
1:A:164:TRP:HB2	1:A:192:LEU:CD1	2.44	0.48
1:B:316:GLN:CD	1:B:514:GLY:H	2.20	0.48
1:C:138:THR:OG1	1:D:334:HIS:HB2	2.13	0.48
1:A:415:GLY:N	1:A:546:VAL:O	2.34	0.48
1:C:385:THR:C	1:C:410:ARG:HB3	2.39	0.48
1:C:389:VAL:N	1:C:412:MET:O	2.34	0.48
1:E:105:PRO:HA	1:E:112:ARG:HH22	1.77	0.48
1:E:209:LEU:HD11	1:E:351:GLN:HB2	1.96	0.48
1:G:211:ILE:HD13	1:G:274:LYS:HD2	1.96	0.48
1:C:473:ASP:HA	1:C:494:ARG:NH2	2.20	0.48
1:E:75:LEU:HD13	1:E:155:ASP:HA	1.94	0.48
1:F:145:ARG:HB2	1:F:206:TYR:CZ	2.49	0.48
1:G:528:VAL:HG13	1:G:544:ILE:HG22	1.96	0.48
1:H:99:ARG:O	1:H:103:GLU:HG3	2.14	0.48
1:C:460:ALA:HB1	1:C:526:LYS:O	2.12	0.48
1:G:220:LYS:HE3	1:G:250:ALA:HB3	1.95	0.48
1:H:421:ARG:O	1:H:425:HIS:ND1	2.37	0.48
1:G:225:GLN:NE2	1:H:218:ALA:O	2.46	0.48
1:G:373:TYR:O	1:G:375:PRO:HD3	2.13	0.48
1:B:240:ILE:O	1:B:244:LEU:HG	2.13	0.48
1:F:144:LYS:HB3	1:F:201:VAL:HG11	1.94	0.48
1:G:133:CYS:HB2	1:G:348:ALA:HB2	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:145:ARG:NE	1:H:200:ALA:HB1	2.29	0.48
1:G:345:SER:OG	1:G:349:ASP:OD1	2.21	0.48
1:H:275:LEU:HB3	1:H:280:TYR:HB3	1.95	0.48
1:E:262:GLY:C	1:E:339:PRO:HG2	2.39	0.47
1:G:39:ASP:OD2	1:G:42:HIS:HA	2.13	0.47
1:A:147:GLU:HA	1:B:145:ARG:NH2	2.29	0.47
1:A:270:MET:HA	1:A:273:VAL:HG12	1.96	0.47
1:C:286:THR:HA	1:C:296:ARG:O	2.14	0.47
1:D:286:THR:HA	1:D:296:ARG:O	2.14	0.47
1:E:91:ARG:HH21	1:E:364:GLU:CD	2.22	0.47
1:F:425:HIS:HD2	1:F:475:TYR:CG	2.32	0.47
1:G:334:HIS:NE2	1:H:135:ASP:OD1	2.43	0.47
1:G:120:LYS:O	1:G:124:ILE:HG12	2.14	0.47
1:A:298:PRO:O	1:A:302:LYS:HG3	2.14	0.47
1:B:494:ARG:NE	1:B:498:VAL:HG11	2.29	0.47
1:C:121:ASN:ND2	1:C:516:PRO:HA	2.28	0.47
1:H:250:ALA:O	1:H:491:LYS:NZ	2.47	0.47
1:H:412:MET:SD	1:H:550:PRO:HG3	2.54	0.47
1:H:425:HIS:HD2	1:H:479:PHE:CE2	2.28	0.47
1:D:477:ASP:OD1	1:D:505:HIS:NE2	2.47	0.47
1:E:225:GLN:HG3	1:F:222:TYR:CE1	2.49	0.47
1:G:227:THR:HG21	1:H:471:ARG:CB	2.45	0.47
1:B:178:ALA:HB2	1:B:188:THR:HG22	1.96	0.47
1:H:259:LEU:O	1:H:328:VAL:HA	2.14	0.47
1:D:467:THR:HG22	1:D:468:THR:H	1.80	0.47
1:D:503:LYS:HD2	1:D:559:GLY:CA	2.44	0.47
1:E:467:THR:OG1	1:E:471:ARG:NH1	2.35	0.47
1:F:161:LYS:NZ	1:F:165:ASN:OD1	2.34	0.47
1:F:392:LYS:HZ3	1:F:415:GLY:HA2	1.80	0.47
1:G:86:VAL:HG22	1:G:350:ARG:HB3	1.97	0.47
1:G:93:ASP:OD1	1:G:279:ARG:NH2	2.46	0.47
1:G:411:VAL:HG22	1:G:551:ALA:O	2.15	0.47
1:G:469:ALA:HB1	1:G:509:TYR:CE1	2.50	0.47
1:H:75:LEU:HB3	1:H:159:LEU:CD2	2.45	0.47
1:H:91:ARG:NH1	1:H:351:GLN:HG3	2.30	0.47
1:H:393:ARG:HH12	1:H:401:GLN:HB2	1.79	0.47
1:H:416:THR:HA	1:H:544:ILE:O	2.14	0.47
1:A:448:TYR:HD1	1:A:489:LEU:O	1.98	0.47
1:B:142:LEU:O	1:B:208:PHE:HA	2.15	0.47
1:B:223:LEU:HB2	1:B:270:MET:SD	2.55	0.47
1:C:148:LEU:HB2	1:C:204:SER:HB3	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:252:CYS:SG	1:C:514:GLY:HA3	2.55	0.47
1:E:39:ASP:OD1	1:E:43:HIS:N	2.48	0.47
1:E:111:ASP:OD2	1:E:319:GLY:HA3	2.15	0.47
1:G:219:ASN:O	1:H:225:GLN:HG3	2.14	0.47
1:G:473:ASP:CG	1:G:498:VAL:HG21	2.40	0.47
1:A:99:ARG:HD3	1:A:373:TYR:HA	1.97	0.47
1:C:464:PHE:HB3	1:C:525:ILE:CD1	2.45	0.47
1:C:532:ALA:HB3	1:C:541:VAL:HB	1.95	0.47
1:G:391:LEU:HB2	1:G:415:GLY:HA3	1.97	0.47
1:H:452:PRO:HB3	1:H:464:PHE:CD2	2.50	0.47
1:B:402:LEU:HD23	1:B:402:LEU:HA	1.72	0.46
1:C:105:PRO:O	1:C:112:ARG:NH2	2.48	0.46
1:C:289:ASP:OD2	1:C:293:ARG:HD3	2.15	0.46
1:D:391:LEU:C	1:D:393:ARG:H	2.23	0.46
1:G:486:TYR:O	1:G:507:GLY:HA2	2.15	0.46
1:B:36:GLN:O	1:B:285:PRO:HD3	2.15	0.46
1:E:528:VAL:HG22	1:E:544:ILE:CB	2.45	0.46
1:G:193:PRO:HG3	1:H:292:GLY:O	2.15	0.46
1:C:69:LYS:HD2	1:D:63:PHE:CE1	2.50	0.46
1:C:261:ILE:HB	1:C:330:ARG:CD	2.45	0.46
1:E:313:ILE:HG12	1:E:566:THR:HB	1.97	0.46
1:F:263:GLY:HA3	1:F:269:THR:OG1	2.16	0.46
1:F:347:SER:HB2	2:F:601:SF4:S4	2.55	0.46
1:F:419:VAL:HB	1:F:542:TRP:CD1	2.50	0.46
1:E:395:ILE:C	1:E:397:LYS:N	2.74	0.46
1:E:408:GLY:N	1:E:553:ILE:O	2.30	0.46
1:G:452:PRO:HB3	1:G:464:PHE:CD2	2.51	0.46
1:E:506:GLY:HA2	1:E:557:ASP:HB2	1.97	0.46
1:A:83:PHE:HE1	1:A:210:PHE:CD1	2.34	0.46
1:A:134:GLN:OE1	1:A:517:ALA:HB2	2.15	0.46
1:C:438:LEU:HD12	1:C:439:PRO:HD2	1.97	0.46
1:D:278:CYS:HA	1:D:351:GLN:HE21	1.81	0.46
1:E:144:LYS:CE	1:E:280:TYR:OH	2.64	0.46
1:G:65:ARG:NH1	1:H:150:TRP:HB3	2.31	0.46
1:H:419:VAL:HB	1:H:542:TRP:HD1	1.80	0.46
1:C:425:HIS:HB3	1:C:475:TYR:CG	2.51	0.46
1:F:391:LEU:HB2	1:F:415:GLY:HA3	1.96	0.46
1:G:183:PHE:HD1	1:H:231:LEU:HB3	1.81	0.46
1:G:215:GLY:HA3	1:H:334:HIS:CG	2.51	0.46
1:G:231:LEU:HD13	1:G:330:ARG:NH2	2.31	0.46
1:H:46:GLU:O	1:H:364:GLU:HA	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:66:GLN:H	1:B:66:GLN:CD	2.23	0.46
1:C:348:ALA:N	2:C:601:SF4:S3	2.88	0.46
1:D:136:THR:HG21	1:D:173:ARG:H	1.81	0.46
1:H:207:GLU:HG2	1:H:355:HIS:HB3	1.97	0.46
1:B:419:VAL:HA	1:B:448:TYR:O	2.16	0.45
1:C:155:ASP:N	1:C:155:ASP:OD1	2.49	0.45
1:C:248:GLY:HA3	1:C:493:ASN:ND2	2.31	0.45
1:F:119:LEU:HD13	1:F:374:LEU:HD13	1.99	0.45
1:G:452:PRO:HB3	1:G:464:PHE:CE2	2.51	0.45
1:H:29:PHE:CZ	1:H:31:PHE:HB2	2.51	0.45
1:A:76:THR:HG23	1:A:162:GLY:N	2.32	0.45
1:C:531:LEU:N	1:C:541:VAL:O	2.37	0.45
1:E:259:LEU:N	1:E:327:ARG:O	2.43	0.45
1:F:89:PHE:CZ	1:F:368:GLN:HG2	2.51	0.45
1:F:145:ARG:N	1:F:199:LEU:O	2.46	0.45
1:A:240:ILE:O	1:A:244:LEU:HG	2.17	0.45
1:D:166:ALA:O	1:D:170:HIS:HB2	2.16	0.45
1:D:348:ALA:HB1	1:D:350:ARG:CZ	2.45	0.45
1:E:256:HIS:O	1:E:344:VAL:HA	2.15	0.45
1:E:415:GLY:O	1:E:546:VAL:HG22	2.16	0.45
1:H:490:ALA:O	1:H:512:SER:N	2.48	0.45
1:A:177:THR:HG22	1:A:185:GLU:HB3	1.97	0.45
1:E:145:ARG:HB2	1:E:206:TYR:CZ	2.51	0.45
1:F:414:ASN:HA	1:F:547:GLU:HA	1.98	0.45
1:H:76:THR:HG23	1:H:162:GLY:N	2.31	0.45
1:H:313:ILE:O	1:H:318:GLY:N	2.41	0.45
1:C:265:SER:OG	1:D:195:GLN:OE1	2.34	0.45
1:E:419:VAL:HA	1:E:448:TYR:O	2.17	0.45
1:E:453:ALA:HB3	1:E:463:SER:HB2	1.98	0.45
1:G:145:ARG:HB3	1:H:200:ALA:CB	2.47	0.45
1:A:102:ILE:HG21	1:A:375:PRO:HG2	1.99	0.45
1:A:173:ARG:NH1	1:A:451:GLY:O	2.50	0.45
1:G:55:LYS:HE3	1:G:56:TYR:CZ	2.52	0.45
1:G:556:ASP:OD2	1:G:560:ASN:HB2	2.17	0.45
1:H:88:HIS:HB3	1:H:126:ALA:HA	1.99	0.45
1:A:506:GLY:HA2	1:A:557:ASP:HB2	1.99	0.45
1:B:90:PHE:HD2	1:B:94:HIS:CD2	2.34	0.45
1:B:91:ARG:HD2	1:B:279:ARG:NH2	2.32	0.45
1:B:253:PRO:HG3	1:B:255:TYR:CE2	2.51	0.45
1:B:421:ARG:O	1:B:425:HIS:ND1	2.41	0.45
1:F:261:ILE:HD13	1:F:340:VAL:HG13	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:254:PRO:HG3	1:H:315:ALA:CB	2.47	0.45
1:H:280:TYR:HD1	1:H:280:TYR:O	2.00	0.45
1:H:419:VAL:HB	1:H:542:TRP:CD1	2.52	0.45
1:A:47:PHE:HD1	1:A:364:GLU:HA	1.82	0.45
1:B:403:SER:HA	1:B:508:PHE:HZ	1.82	0.45
1:E:253:PRO:HA	1:E:254:PRO:C	2.42	0.45
1:C:164:TRP:CD1	1:C:190:ASP:HB2	2.52	0.45
1:C:374:LEU:HD12	1:C:374:LEU:HA	1.82	0.45
1:E:469:ALA:HB2	1:E:490:ALA:HB3	1.98	0.45
1:G:135:ASP:OD1	1:G:173:ARG:NH2	2.50	0.45
1:H:348:ALA:HB1	1:H:350:ARG:CZ	2.47	0.45
1:C:438:LEU:HD22	1:C:479:PHE:O	2.17	0.45
1:D:91:ARG:HD2	1:D:279:ARG:NH2	2.31	0.45
1:F:174:TYR:O	1:F:187:ASN:ND2	2.41	0.45
1:H:180:LEU:N	1:H:184:LYS:O	2.50	0.45
1:H:333:ARG:HD2	1:H:337:SER:O	2.17	0.45
1:H:392:LYS:NZ	1:H:415:GLY:HA2	2.32	0.45
1:B:134:GLN:NE2	2:B:601:SF4:S4	2.85	0.44
1:C:69:LYS:HD2	1:D:63:PHE:CZ	2.51	0.44
1:F:136:THR:HG21	1:F:172:LEU:HD22	2.00	0.44
1:G:494:ARG:NH1	1:G:509:TYR:CD1	2.77	0.44
1:A:256:HIS:HB2	1:A:345:SER:HB3	1.99	0.44
1:C:83:PHE:HE1	1:C:210:PHE:CD1	2.35	0.44
1:C:480:GLN:HG2	1:C:486:TYR:CD1	2.52	0.44
1:D:224:TYR:O	1:D:340:VAL:N	2.46	0.44
1:E:334:HIS:NE2	1:F:135[B]:ASP:OD1	2.50	0.44
1:G:65:ARG:CZ	1:G:148:LEU:HD11	2.47	0.44
1:H:76:THR:OG1	1:H:158:TYR:O	2.33	0.44
1:C:172:LEU:O	1:C:454:LYS:N	2.49	0.44
1:E:254:PRO:HA	1:E:322:PHE:O	2.17	0.44
1:E:272:THR:HG23	1:E:281:TYR:CE2	2.52	0.44
1:E:334:HIS:NE2	1:F:135[A]:ASP:OD1	2.50	0.44
1:G:253:PRO:HD2	1:G:316:GLN:HG2	1.99	0.44
1:H:47:PHE:HB3	1:H:362:TYR:HB3	2.00	0.44
1:A:490:ALA:N	1:A:509:TYR:HE1	2.14	0.44
1:D:113:TYR:OH	1:D:410:ARG:HD3	2.17	0.44
1:D:421:ARG:HG3	1:D:539:GLU:OE2	2.18	0.44
1:E:419:VAL:O	1:E:541:VAL:HA	2.18	0.44
1:F:325:GLN:OE1	1:F:327:ARG:NH2	2.49	0.44
1:B:449:TYR:CD2	1:B:490:ALA:HB2	2.53	0.44
1:D:253:PRO:HA	1:D:254:PRO:C	2.42	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:532:ALA:O	1:F:540:ALA:HB1	2.17	0.44
1:G:489:LEU:HA	1:G:510:LEU:HB2	1.99	0.44
1:D:91:ARG:NE	1:D:364:GLU:OE2	2.45	0.44
1:E:150:TRP:HB3	1:F:65:ARG:NH1	2.31	0.44
1:F:214:GLY:HA3	2:F:601:SF4:S3	2.58	0.44
1:B:94:HIS:CE1	1:B:256:HIS:HE2	2.34	0.44
1:B:424:ALA:O	1:B:428:ILE:HG13	2.18	0.44
1:D:143:GLY:O	1:D:199:LEU:N	2.43	0.44
1:E:219:ASN:HA	1:F:225:GLN:CD	2.43	0.44
1:E:240:ILE:O	1:E:244:LEU:HG	2.17	0.44
1:G:255:TYR:O	1:G:323:ALA:HA	2.17	0.44
1:A:209:LEU:HD11	1:A:351:GLN:HB2	1.98	0.44
1:B:210:PHE:O	1:B:351:GLN:HA	2.18	0.44
1:C:173:ARG:HD2	1:C:538:MET:SD	2.58	0.44
1:C:251:ALA:HB3	1:C:255:TYR:CE1	2.52	0.44
1:C:531:LEU:HD22	1:C:543:LYS:HE3	1.98	0.44
1:F:466:PRO:HG2	1:F:491:LYS:HG2	1.99	0.44
1:B:262:GLY:C	1:B:339:PRO:HG2	2.43	0.44
1:C:543:LYS:HE3	1:C:543:LYS:HB2	1.76	0.44
1:E:418:ILE:HG12	1:E:445:SER:HB2	2.00	0.44
1:E:438:LEU:HD12	1:E:439:PRO:HD2	2.00	0.44
1:F:62:VAL:HG11	1:F:150:TRP:CH2	2.44	0.44
1:E:346:CYS:CB	2:E:601:SF4:S4	3.06	0.43
1:C:120:LYS:O	1:C:124:ILE:HG12	2.19	0.43
1:E:193:PRO:HG3	1:F:292:GLY:O	2.17	0.43
1:F:419:VAL:HB	1:F:542:TRP:HB2	2.00	0.43
1:C:131:PRO:HD2	1:C:463:SER:OG	2.18	0.43
1:C:256:HIS:O	1:C:344:VAL:HA	2.18	0.43
1:C:264:THR:OG1	1:D:195:GLN:HB2	2.17	0.43
1:F:192:LEU:HB3	1:F:193:PRO:HA	1.99	0.43
1:F:210:PHE:HE1	1:F:354:ALA:HB3	1.83	0.43
1:C:531:LEU:HD22	1:C:543:LYS:HB2	2.01	0.43
1:D:102:ILE:O	1:D:112:ARG:HD3	2.18	0.43
1:D:469:ALA:HB1	1:D:494:ARG:CZ	2.49	0.43
1:F:75:LEU:HB3	1:F:159:LEU:CD2	2.48	0.43
1:F:177:THR:CG2	1:F:185:GLU:HB3	2.47	0.43
1:F:250:ALA:HB1	1:F:468:THR:HG23	2.00	0.43
1:C:124:ILE:HD12	1:C:522:LYS:CD	2.49	0.43
1:D:80:HIS:HD1	1:D:80:HIS:C	2.27	0.43
1:E:88:HIS:CE1	1:E:128:ARG:HA	2.54	0.43
1:E:219:ASN:HD21	1:F:337:SER:HA	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:437:PRO:O	1:E:438:LEU:C	2.60	0.43
1:H:133:CYS:HB3	1:H:350:ARG:HH22	1.84	0.43
1:H:480:GLN:HE22	1:H:486:TYR:HA	1.83	0.43
1:B:489:LEU:HD21	1:B:549:PHE:CE1	2.54	0.43
1:E:88:HIS:ND1	1:E:125:ALA:O	2.39	0.43
1:G:133:CYS:SG	1:G:134:GLN:N	2.91	0.43
1:H:171:ASN:CA	1:H:454:LYS:HB2	2.47	0.43
1:B:425:HIS:CD2	1:B:472:MET:HG2	2.52	0.43
1:E:233:PRO:HA	1:E:300:TRP:CZ2	2.54	0.43
1:A:93:ASP:OD1	1:A:94:HIS:N	2.52	0.43
1:F:57:VAL:HA	1:F:69:LYS:O	2.18	0.43
1:G:57:VAL:HB	1:G:68:LEU:HD11	2.00	0.43
1:H:88:HIS:CD2	1:H:128:ARG:HA	2.53	0.43
1:F:285:PRO:O	1:F:295:PHE:HB2	2.18	0.43
1:F:473:ASP:HA	1:F:494:ARG:HH21	1.83	0.43
1:G:410:ARG:HB3	1:G:552:PHE:CZ	2.54	0.43
1:G:425:HIS:ND1	1:G:472:MET:HG3	2.34	0.43
1:A:275:LEU:HB3	1:A:280:TYR:HB3	1.99	0.43
1:B:529:THR:N	1:B:543:LYS:O	2.46	0.43
1:D:94:HIS:CE1	1:D:256:HIS:HE2	2.37	0.43
1:F:418:ILE:HG13	1:F:441:TYR:O	2.19	0.43
1:G:446:PRO:HG3	1:G:487:ILE:HD11	2.00	0.43
1:G:480:GLN:OE1	1:G:505:HIS:HB2	2.19	0.43
1:H:98:TRP:CZ2	1:H:118:LEU:HB3	2.54	0.43
1:D:428:ILE:HD11	1:D:441:TYR:HE1	1.84	0.42
1:D:438:LEU:HD12	1:D:439:PRO:HD2	2.00	0.42
1:H:425:HIS:CD2	1:H:479:PHE:HE2	2.28	0.42
1:A:300:TRP:CZ2	1:B:183:PHE:HZ	2.38	0.42
1:F:134:GLN:HA	1:F:463:SER:HB3	2.00	0.42
1:G:279:ARG:HH12	1:G:351:GLN:HE22	1.67	0.42
1:H:268:MET:SD	1:H:331:LEU:HD13	2.59	0.42
1:H:415:GLY:O	1:H:546:VAL:HG22	2.19	0.42
1:B:47:PHE:HB3	1:B:362:TYR:HB3	2.01	0.42
1:C:168:ARG:HG3	1:C:169:TYR:CD1	2.54	0.42
1:E:409:THR:HB	1:E:553:ILE:HD12	2.00	0.42
1:F:452:PRO:HD2	1:F:538:MET:HA	2.00	0.42
1:G:136:THR:OG1	1:G:173:ARG:O	2.33	0.42
1:G:272:THR:HG21	1:G:331:LEU:HD11	2.01	0.42
1:H:158:TYR:HD1	1:H:158:TYR:HA	1.68	0.42
1:H:239:PHE:CE1	1:H:243:LYS:HE2	2.54	0.42
1:H:452:PRO:HD2	1:H:538:MET:HA	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:422:ASP:CG	1:D:471:ARG:HH11	2.25	0.42
1:E:467:THR:O	1:E:490:ALA:HB1	2.19	0.42
1:A:468:THR:HA	1:A:491:LYS:O	2.20	0.42
1:D:259:LEU:O	1:D:328:VAL:HA	2.19	0.42
1:D:289:ASP:OD2	1:D:293:ARG:HD3	2.19	0.42
1:E:313:ILE:O	1:E:318:GLY:N	2.50	0.42
1:G:209:LEU:HD22	1:G:280:TYR:CD1	2.55	0.42
1:A:164:TRP:CD1	1:A:190:ASP:HB2	2.54	0.42
1:A:259:LEU:HA	1:A:341:GLY:O	2.20	0.42
1:B:178:ALA:N	1:B:186:CYS:O	2.51	0.42
1:B:412:MET:HE3	1:B:548:ASP:HB3	2.01	0.42
1:B:432:MET:HA	1:B:436:GLU:O	2.20	0.42
1:C:314:GLY:O	1:C:322:PHE:HD1	2.03	0.42
1:F:419:VAL:HB	1:F:542:TRP:HD1	1.85	0.42
1:G:333:ARG:HE	1:H:185:GLU:CD	2.27	0.42
1:A:111:ASP:OD2	1:A:320:LYS:N	2.52	0.42
1:B:95:LEU:HD23	1:B:98:TRP:CZ3	2.55	0.42
1:B:448:TYR:CZ	1:B:549:PHE:HZ	2.38	0.42
1:B:499:THR:HG21	1:B:561:ASP:HB2	2.01	0.42
1:C:148:LEU:HA	1:C:150:TRP:NE1	2.35	0.42
1:C:150:TRP:HB3	1:D:65:ARG:NH1	2.33	0.42
1:D:34:ILE:O	1:D:294:ALA:N	2.31	0.42
1:E:233:PRO:HA	1:E:300:TRP:HZ2	1.83	0.42
1:F:80:HIS:O	1:F:80:HIS:ND1	2.52	0.42
1:H:309:GLN:HG2	1:H:320:LYS:O	2.20	0.42
1:A:158:TYR:HD1	1:A:158:TYR:HA	1.68	0.42
1:A:392:LYS:HE3	1:A:415:GLY:HA2	2.00	0.42
1:B:435:GLY:O	1:B:437:PRO:HD3	2.19	0.42
1:E:333:ARG:HE	1:F:185:GLU:CD	2.28	0.42
1:F:120:LYS:O	1:F:124:ILE:HG12	2.20	0.42
1:A:464:PHE:N	1:A:517:ALA:HB1	2.35	0.42
1:B:261:ILE:HD13	1:B:340:VAL:HG13	2.02	0.42
1:B:316:GLN:OE1	1:B:514:GLY:N	2.39	0.42
1:B:389:VAL:HG21	1:B:413:LEU:HD23	2.01	0.42
1:C:245:LYS:NZ	1:C:568:ALA:HB3	2.35	0.42
1:C:430:GLU:O	1:C:434:ASN:ND2	2.46	0.42
1:E:220:LYS:CE	1:E:250:ALA:HB3	2.47	0.42
1:E:230:LEU:HD11	1:E:236:LEU:HA	2.00	0.42
1:G:520:LEU:HD23	1:G:520:LEU:HA	1.93	0.42
1:A:254:PRO:HB2	1:A:324:HIS:CG	2.55	0.42
1:A:391:LEU:C	1:A:393:ARG:H	2.28	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:133:CYS:HB3	1:B:350:ARG:NH2	2.35	0.42
1:B:494:ARG:HD3	1:B:498:VAL:HG11	2.01	0.42
1:C:135:ASP:CG	1:C:173:ARG:HH22	2.28	0.42
1:C:317:PHE:HB3	1:C:562:MET:HE1	2.02	0.42
1:E:529:THR:O	1:E:543:LYS:N	2.40	0.42
1:E:533:PHE:N	1:E:534:PRO:HD3	2.34	0.42
1:F:156:GLU:HG3	1:F:198:LEU:HD11	2.02	0.42
1:H:166:ALA:HA	1:H:170:HIS:CE1	2.55	0.42
1:A:211:ILE:CG2	1:A:213:LYS:HG2	2.50	0.41
1:B:520:LEU:HD21	1:B:550:PRO:HD2	2.02	0.41
1:E:94:HIS:NE2	1:E:347:SER:O	2.50	0.41
1:F:134:GLN:CB	1:F:465:GLY:H	2.30	0.41
1:F:456:PRO:HB2	1:F:459:TYR:CD2	2.54	0.41
1:A:171:ASN:CG	1:A:456:PRO:HA	2.45	0.41
1:B:438:LEU:HA	1:B:439:PRO:HD3	1.94	0.41
1:C:88:HIS:CG	1:C:128:ARG:HA	2.55	0.41
1:C:200:ALA:HB1	1:D:145:ARG:CZ	2.50	0.41
1:G:167:TYR:HB3	1:G:174:TYR:CE1	2.55	0.41
1:C:104:ASP:OD2	1:C:106:GLU:HB2	2.20	0.41
1:D:357:ASN:HB2	1:D:358:LYS:HD3	2.02	0.41
1:E:173:ARG:HH12	3:E:602:GOL:HO1	1.62	0.41
1:E:361:ILE:C	1:E:362:TYR:HD1	2.28	0.41
1:G:136:THR:O	1:G:167:TYR:OH	2.35	0.41
1:G:136:THR:HB	1:G:167:TYR:HE2	1.85	0.41
1:G:139:ALA:HB1	1:G:163:ILE:HD13	2.02	0.41
1:G:494:ARG:NH1	1:G:509:TYR:HD1	2.11	0.41
1:E:36:GLN:O	1:E:285:PRO:HD3	2.20	0.41
1:F:87:HIS:CE1	1:F:136:THR:HB	2.54	0.41
1:H:294:ALA:HA	1:H:330:ARG:O	2.21	0.41
1:A:90:PHE:HE1	1:A:348:ALA:HA	1.86	0.41
1:C:425:HIS:HB3	1:C:475:TYR:CZ	2.55	0.41
1:D:452:PRO:HB3	1:D:464:PHE:CD2	2.56	0.41
1:F:417:LEU:HD12	1:F:546:VAL:HG11	2.02	0.41
1:B:296:ARG:HH21	1:B:301:GLU:CD	2.29	0.41
1:C:334:HIS:CE1	1:D:175:SER:HB3	2.55	0.41
1:D:532:ALA:O	1:D:540:ALA:HB1	2.21	0.41
1:F:254:PRO:HB2	1:F:324:HIS:ND1	2.35	0.41
1:H:166:ALA:O	1:H:170:HIS:ND1	2.54	0.41
1:H:508:PHE:CD1	1:H:556:ASP:HA	2.53	0.41
1:H:510:LEU:CD2	1:H:553:ILE:HA	2.50	0.41
1:A:446:PRO:HA	1:A:487:ILE:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:449:TYR:HD2	1:A:469:ALA:HA	1.85	0.41
1:C:42:HIS:HB3	1:C:45:THR:HB	2.02	0.41
1:C:148:LEU:HD22	1:C:150:TRP:HZ2	1.85	0.41
1:D:421:ARG:HG3	1:D:539:GLU:CD	2.46	0.41
1:G:65:ARG:HH22	1:G:147:GLU:CD	2.28	0.41
1:H:300:TRP:O	1:H:304:VAL:HG23	2.21	0.41
1:A:86:VAL:HG22	1:A:350:ARG:HB3	2.03	0.41
1:B:83:PHE:CE2	1:B:210:PHE:HD2	2.38	0.41
1:B:268:MET:HE1	1:B:331:LEU:HD13	2.03	0.41
1:B:466:PRO:HB2	1:B:491:LYS:HG2	2.01	0.41
1:C:37:PRO:HA	1:C:284:LEU:HD23	2.02	0.41
1:H:65:ARG:NE	1:H:148:LEU:HD11	2.35	0.41
1:B:105:PRO:HA	1:B:112:ARG:HH12	1.86	0.41
1:B:223:LEU:HD12	1:B:340:VAL:O	2.20	0.41
1:B:422:ASP:OD1	1:B:471:ARG:NH1	2.47	0.41
1:C:447:ILE:O	1:C:489:LEU:N	2.53	0.41
1:D:78:LEU:HD22	1:D:356:ILE:HD11	2.03	0.41
1:E:422:ASP:CG	1:E:471:ARG:HE	2.28	0.41
1:F:78:LEU:C	1:F:80:HIS:N	2.77	0.41
1:F:97:GLY:HA3	1:F:324:HIS:CD2	2.48	0.41
1:F:467:THR:O	1:F:490:ALA:HB1	2.21	0.41
1:G:193:PRO:HD3	1:H:291:TYR:O	2.21	0.41
1:H:164:TRP:O	1:H:168:ARG:HB3	2.20	0.41
1:D:96:GLU:HG3	1:D:373:TYR:HE1	1.86	0.41
1:H:280:TYR:O	1:H:280:TYR:CD1	2.73	0.41
1:A:111:ASP:OD2	1:A:319:GLY:HA3	2.21	0.40
1:A:192:LEU:HD23	1:A:192:LEU:HA	1.87	0.40
1:C:254:PRO:HA	1:C:322:PHE:O	2.20	0.40
1:C:325:GLN:OE1	1:C:327:ARG:NH2	2.53	0.40
1:E:65:ARG:NH1	1:F:152:GLY:HA2	2.36	0.40
1:E:233:PRO:O	1:E:300:TRP:HH2	2.04	0.40
1:E:259:LEU:HD12	1:E:341:GLY:O	2.21	0.40
1:F:158:TYR:HD1	1:F:158:TYR:HA	1.69	0.40
1:F:412:MET:HB3	1:F:548:ASP:HA	2.02	0.40
1:F:479:PHE:HD2	1:F:485:SER:HB2	1.86	0.40
1:A:75:LEU:CD1	1:A:155:ASP:HA	2.51	0.40
1:B:421:ARG:NH1	1:B:451:GLY:HA3	2.15	0.40
1:C:185:GLU:OE2	1:D:333:ARG:NH2	2.46	0.40
1:C:259:LEU:HA	1:C:341:GLY:O	2.21	0.40
1:C:285:PRO:O	1:C:296:ARG:N	2.36	0.40
1:D:358:LYS:H	1:D:358:LYS:CD	2.26	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:312:GLY:O	1:E:318:GLY:HA2	2.21	0.40
1:F:256:HIS:O	1:F:344:VAL:HA	2.21	0.40
1:G:65:ARG:NH1	1:G:148:LEU:HD21	2.36	0.40
1:A:121:ASN:HB2	1:A:516:PRO:HA	2.02	0.40
1:C:463:SER:HB3	1:C:517:ALA:HB1	2.04	0.40
1:C:520:LEU:O	1:C:525:ILE:HG13	2.22	0.40
1:D:374:LEU:HD12	1:D:374:LEU:HA	1.86	0.40
1:F:554:VAL:O	1:F:561:ASP:HA	2.21	0.40
1:H:164:TRP:CH2	1:H:168:ARG:HG2	2.56	0.40
1:A:78:LEU:HD13	1:A:361:ILE:CG2	2.51	0.40
1:A:418:ILE:HG12	1:A:445:SER:HB2	2.04	0.40
1:B:167:TYR:HA	1:B:172:LEU:HD12	2.03	0.40
1:D:148:LEU:HA	1:D:150:TRP:NE1	2.36	0.40
1:F:391:LEU:N	1:F:414:ASN:O	2.38	0.40
1:G:50:VAL:HB	1:G:53:SER:HB3	2.03	0.40
1:G:179:ALA:HB2	1:H:330:ARG:CZ	2.52	0.40
1:G:218:ALA:O	1:H:225:GLN:NE2	2.55	0.40
1:G:528:VAL:HG22	1:G:544:ILE:HG22	2.03	0.40
1:H:37:PRO:HA	1:H:284:LEU:HD23	2.03	0.40
1:H:452:PRO:HB3	1:H:464:PHE:CG	2.56	0.40
1:A:410:ARG:HA	1:A:552:PHE:HA	2.03	0.40
1:C:533:PHE:N	1:C:534:PRO:HD3	2.36	0.40
1:D:110:ASN:HA	1:D:113:TYR:HB3	2.04	0.40
1:F:266:ALA:O	1:F:270:MET:HG2	2.22	0.40
1:H:248:GLY:HA3	1:H:493:ASN:OD1	2.22	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	535/604 (89%)	513 (96%)	20 (4%)	2 (0%)	30 62

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	530/604 (88%)	511 (96%)	17 (3%)	2 (0%)	30	62
1	C	528/604 (87%)	512 (97%)	16 (3%)	0	100	100
1	D	526/604 (87%)	508 (97%)	17 (3%)	1 (0%)	43	73
1	E	527/604 (87%)	506 (96%)	17 (3%)	4 (1%)	16	50
1	F	528/604 (87%)	510 (97%)	16 (3%)	2 (0%)	30	62
1	G	528/604 (87%)	513 (97%)	15 (3%)	0	100	100
1	H	527/604 (87%)	510 (97%)	16 (3%)	1 (0%)	43	73
All	All	4229/4832 (88%)	4083 (96%)	134 (3%)	12 (0%)	36	68

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	387	VAL
1	B	438	LEU
1	E	386	SER
1	E	395	ILE
1	H	374	LEU
1	D	536	LEU
1	E	396	ASP
1	F	79	ALA
1	A	27	ALA
1	F	536	LEU
1	E	438	LEU
1	A	25	VAL

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	363/492 (74%)	358 (99%)	5 (1%)	59	77
1	B	369/492 (75%)	365 (99%)	4 (1%)	65	79
1	C	347/492 (70%)	343 (99%)	4 (1%)	63	79

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	362/492 (74%)	358 (99%)	4 (1%)	65	79
1	E	363/492 (74%)	357 (98%)	6 (2%)	53	75
1	F	391/492 (80%)	386 (99%)	5 (1%)	61	78
1	G	357/492 (73%)	352 (99%)	5 (1%)	59	77
1	H	365/492 (74%)	361 (99%)	4 (1%)	65	79
All	All	2917/3936 (74%)	2880 (99%)	37 (1%)	61	78

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	62	VAL
1	A	274	LYS
1	A	374	LEU
1	A	386	SER
1	A	538	MET
1	B	352	ILE
1	B	358	LYS
1	B	374	LEU
1	B	457	GLU
1	C	62	VAL
1	C	352	ILE
1	C	374	LEU
1	C	538	MET
1	D	80	HIS
1	D	352	ILE
1	D	374	LEU
1	D	538	MET
1	E	62	VAL
1	E	144	LYS
1	E	352	ILE
1	E	386	SER
1	E	410	ARG
1	E	538	MET
1	F	274	LYS
1	F	352	ILE
1	F	374	LEU
1	F	410	ARG
1	F	538	MET
1	G	62	VAL
1	G	352	ILE

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Mol	Chain	Res	Type
1	G	368	GLN
1	G	467	THR
1	G	538	MET
1	H	352	ILE
1	H	374	LEU
1	H	410	ARG
1	H	467	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	88	HIS
1	A	480	GLN
1	B	30	HIS
1	B	87	HIS
1	B	171	ASN
1	B	325	GLN
1	B	372	GLN
1	C	497	GLN
1	D	87	HIS
1	D	176	GLN
1	E	87	HIS
1	F	30	HIS
1	F	42	HIS
1	F	87	HIS
1	F	170	HIS
1	F	195	GLN
1	F	324	HIS
1	F	355	HIS
1	F	372	GLN
1	F	482	HIS
1	G	482	HIS
1	H	480	GLN
1	H	497	GLN
1	H	560	ASN

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](#)

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	SF4	B	601	1	0,12,12	-	-	-		
2	SF4	C	601	1	0,12,12	-	-	-		
2	SF4	H	601	1	0,12,12	-	-	-		
3	GOL	E	602	-	5,5,5	0.96	0	5,5,5	1.24	1 (20%)
2	SF4	A	601	1	0,12,12	-	-	-		
2	SF4	E	601	1	0,12,12	-	-	-		
2	SF4	D	601	1	0,12,12	-	-	-		
2	SF4	F	601	1	0,12,12	-	-	-		
2	SF4	G	601	1	0,12,12	-	-	-		
3	GOL	B	602	-	5,5,5	1.00	0	5,5,5	1.01	1 (20%)

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	SF4	B	601	1	-	-	0/6/5/5
2	SF4	C	601	1	-	-	0/6/5/5
2	SF4	H	601	1	-	-	0/6/5/5
3	GOL	E	602	-	-	2/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SF4	A	601	1	-	-	0/6/5/5
2	SF4	D	601	1	-	-	0/6/5/5
2	SF4	E	601	1	-	-	0/6/5/5
2	SF4	F	601	1	-	-	0/6/5/5
2	SF4	G	601	1	-	-	0/6/5/5
3	GOL	B	602	-	-	2/4/4/4	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	602	GOL	C3-C2-C1	-2.15	103.91	111.80
3	B	602	GOL	C3-C2-C1	-2.05	104.29	111.80

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	E	602	GOL	C1-C2-C3-O3
3	B	602	GOL	O1-C1-C2-O2
3	B	602	GOL	O1-C1-C2-C3
3	E	602	GOL	O2-C2-C3-O3

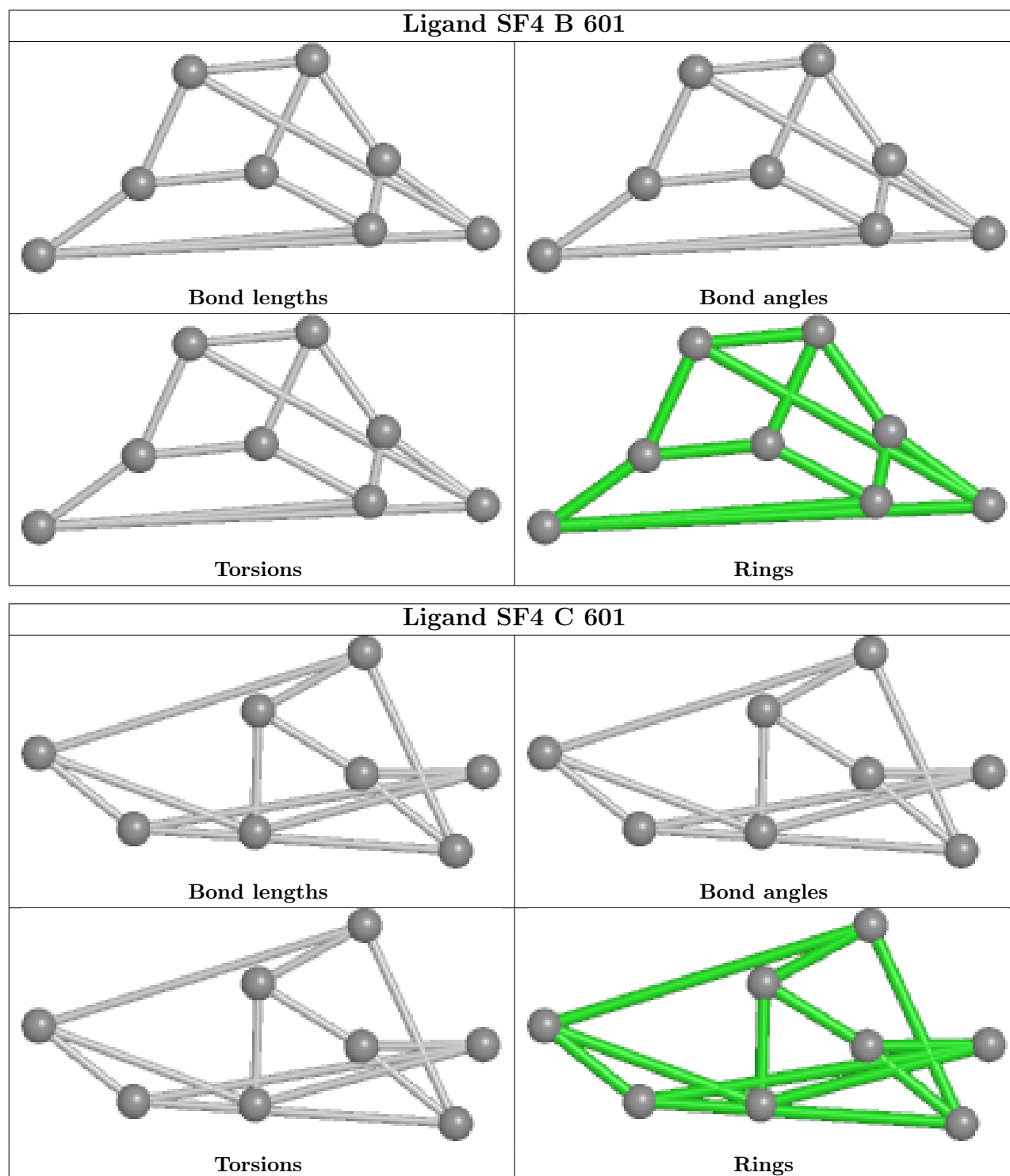
There are no ring outliers.

7 monomers are involved in 16 short contacts:

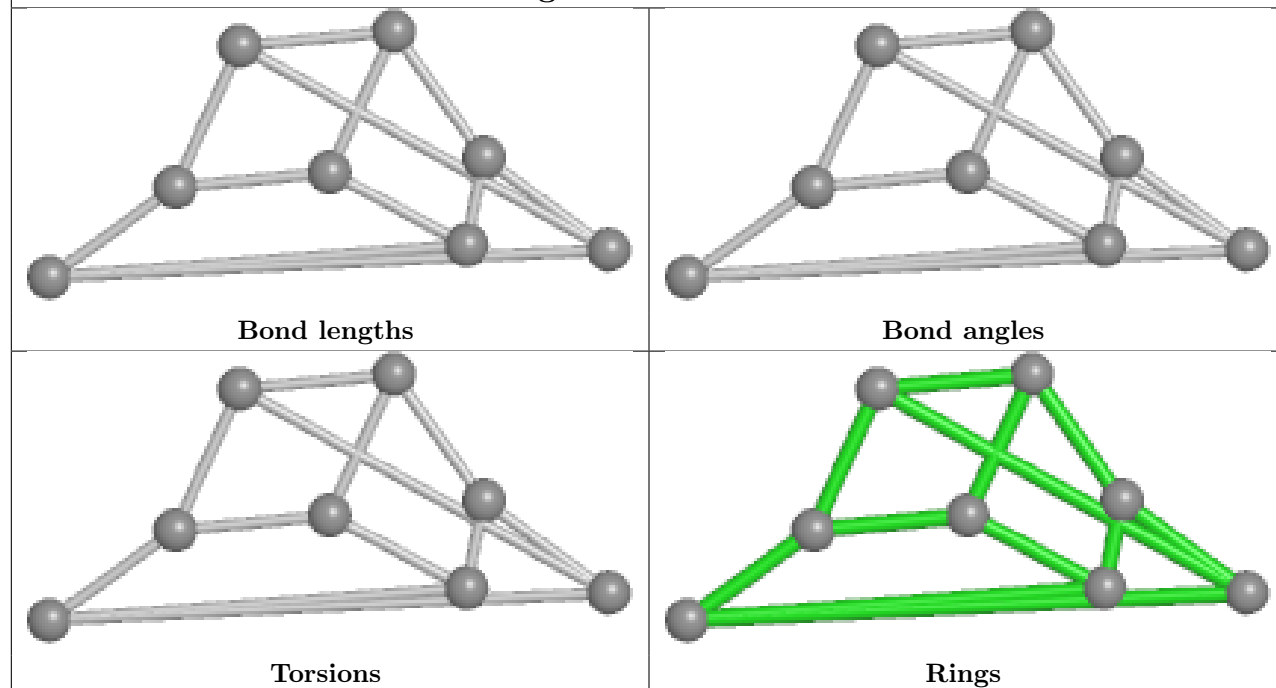
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	601	SF4	2	0
2	C	601	SF4	5	0
3	E	602	GOL	2	0
2	A	601	SF4	1	0
2	E	601	SF4	1	0
2	F	601	SF4	3	0
3	B	602	GOL	2	0

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

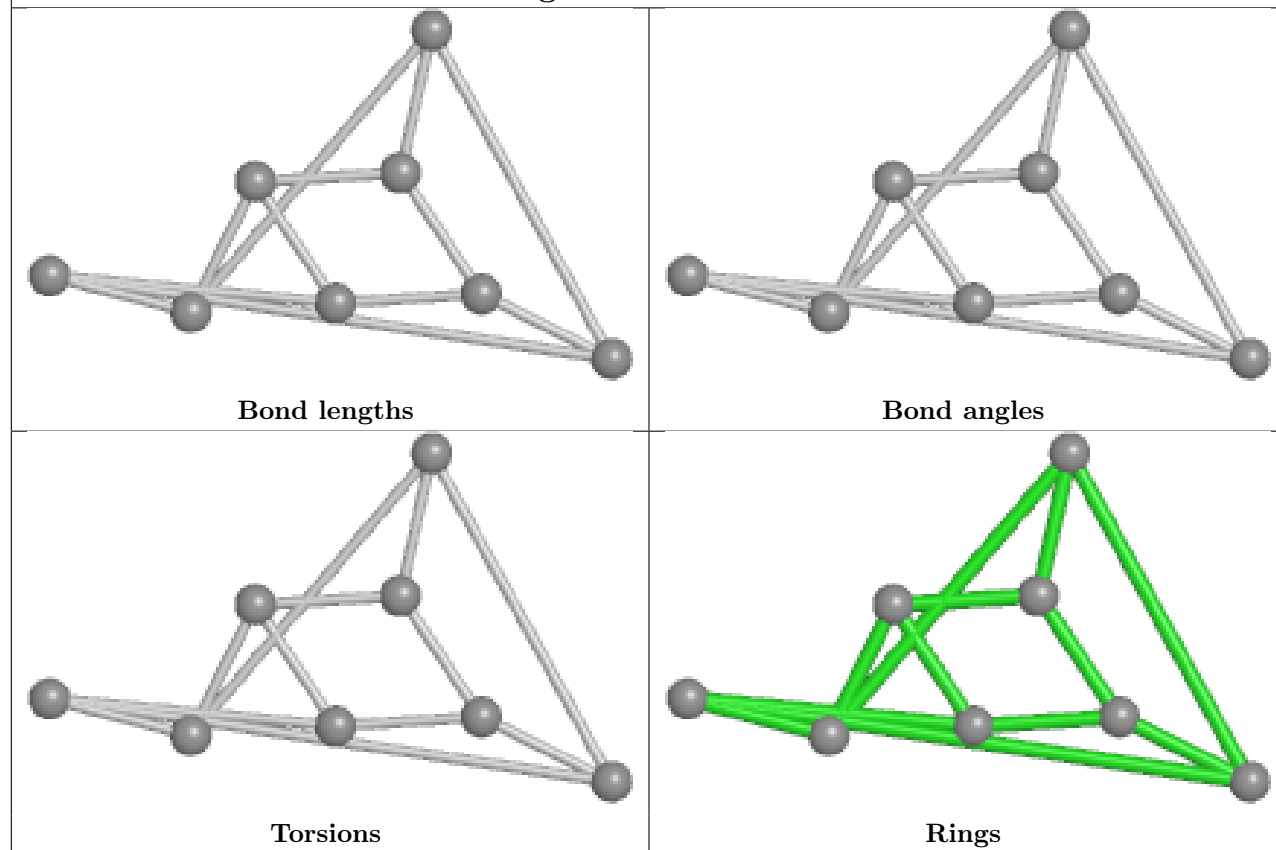
highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified 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.

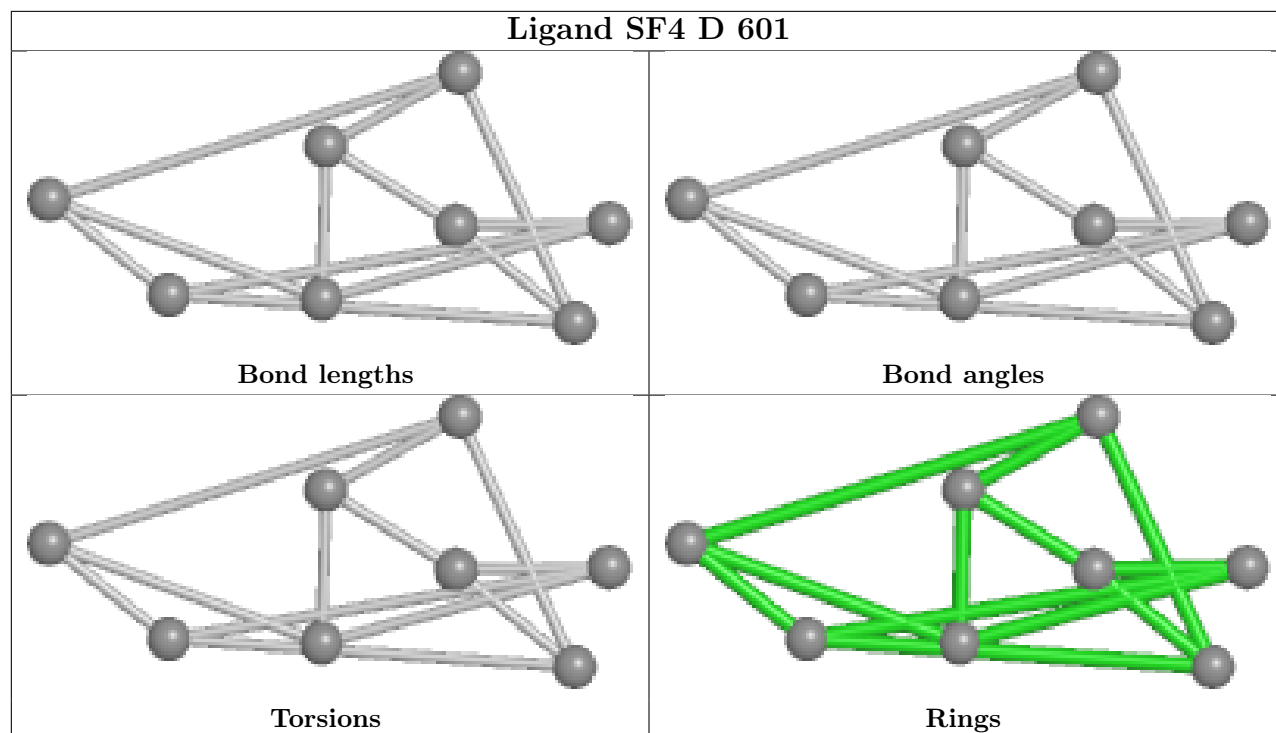
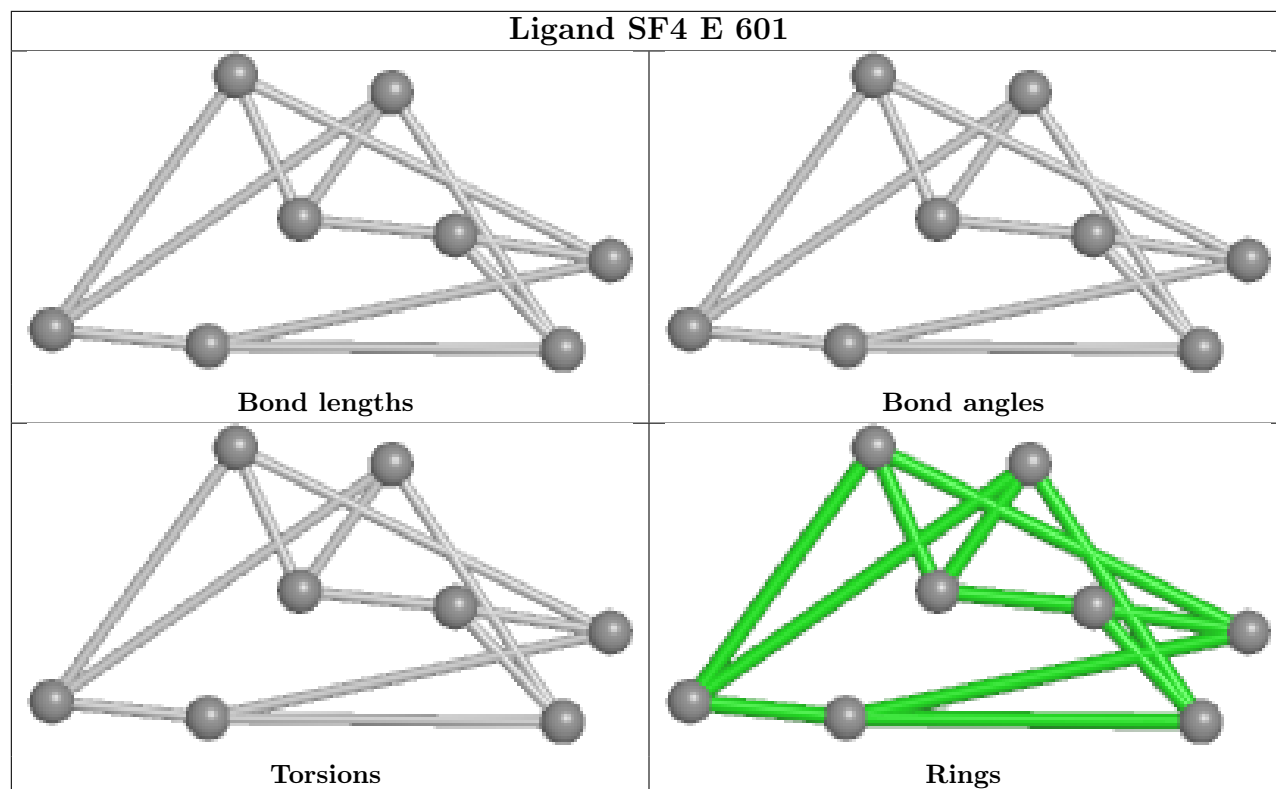


Ligand SF4 H 601

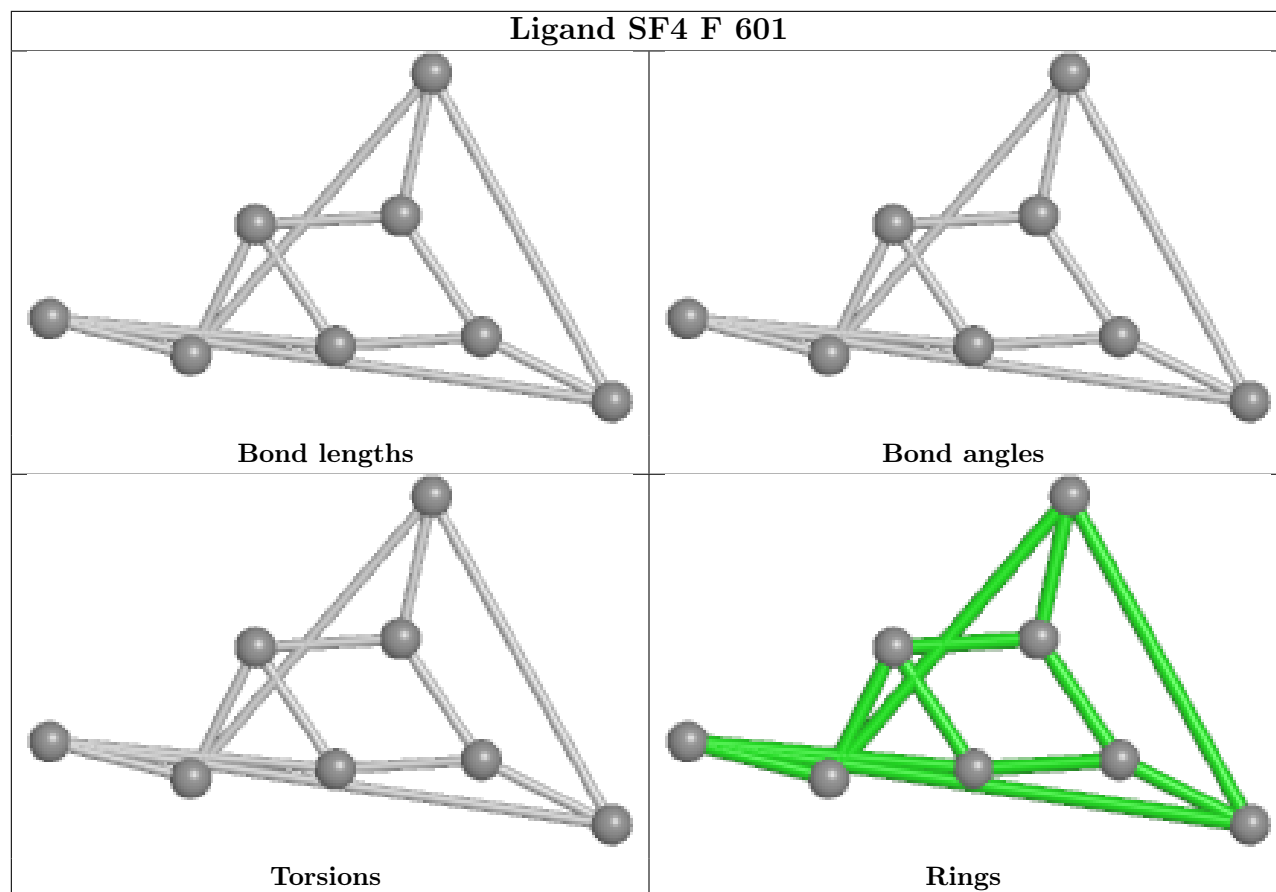


Ligand SF4 A 601

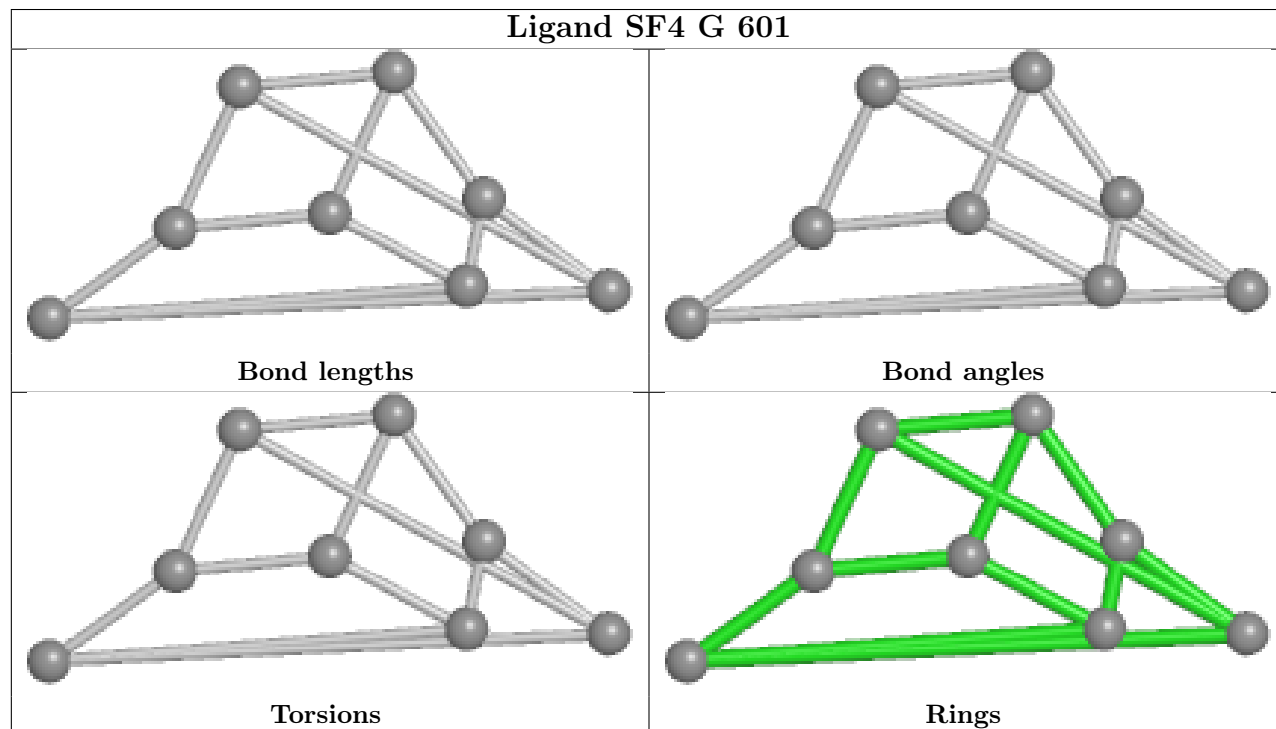




Ligand SF4 F 601



Ligand SF4 G 601



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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	539/604 (89%)	-0.50	1 (0%) 91 85	23, 58, 89, 106	0
1	B	534/604 (88%)	-0.53	0 100 100	24, 50, 79, 89	0
1	C	532/604 (88%)	-0.41	1 (0%) 91 85	37, 80, 91, 100	0
1	D	530/604 (87%)	-0.40	1 (0%) 91 85	29, 72, 89, 102	0
1	E	531/604 (87%)	-0.47	0 100 100	34, 65, 92, 103	0
1	F	531/604 (87%)	-0.54	0 100 100	34, 56, 71, 77	1 (0%)
1	G	532/604 (88%)	-0.45	0 100 100	28, 66, 91, 102	0
1	H	531/604 (87%)	-0.49	1 (0%) 91 85	22, 65, 88, 111	0
All	All	4260/4832 (88%)	-0.47	4 (0%) 92 89	22, 64, 89, 111	1 (0%)

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	250	ALA	2.7
1	A	540	ALA	2.3
1	C	454	LYS	2.1
1	H	509	TYR	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

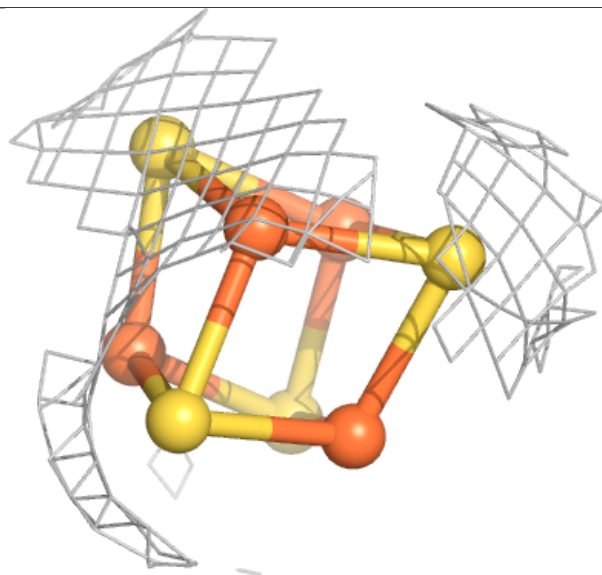
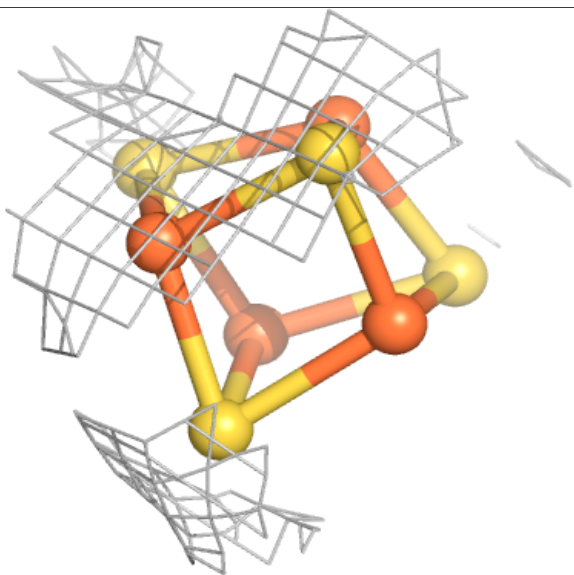
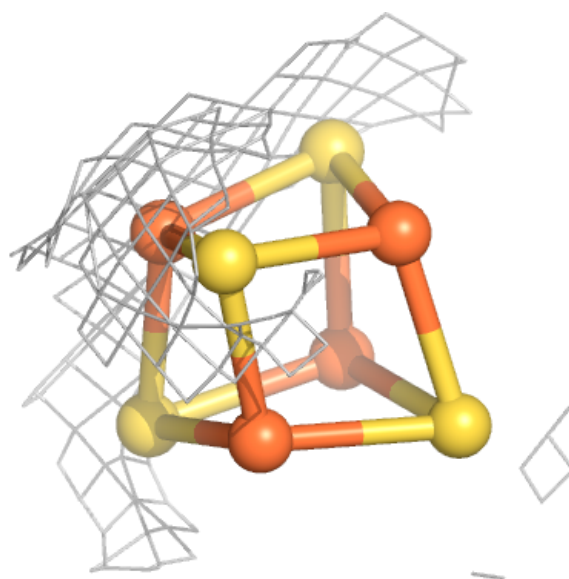
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	GOL	E	602	6/6	0.97	0.05	58,59,65,65	0
3	GOL	B	602	6/6	0.98	0.09	35,37,41,41	0
2	SF4	H	601	8/8	0.99	0.03	65,70,94,94	0
2	SF4	C	601	8/8	0.99	0.02	80,84,105,116	0
2	SF4	G	601	8/8	0.99	0.04	48,80,92,96	0
2	SF4	F	601	8/8	1.00	0.03	51,77,94,94	0
2	SF4	B	601	8/8	1.00	0.03	29,40,52,74	0
2	SF4	A	601	8/8	1.00	0.03	53,65,77,81	0
2	SF4	D	601	8/8	1.00	0.02	52,59,78,85	0
2	SF4	E	601	8/8	1.00	0.02	55,69,89,96	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.

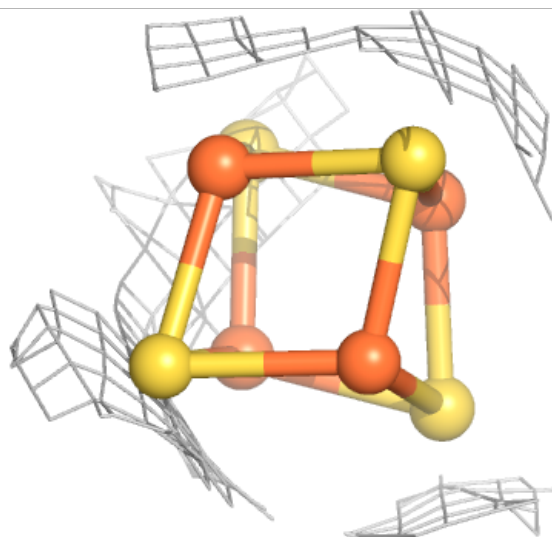
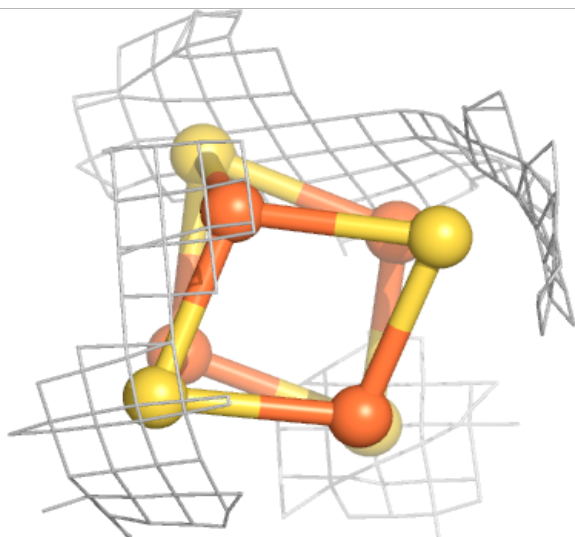
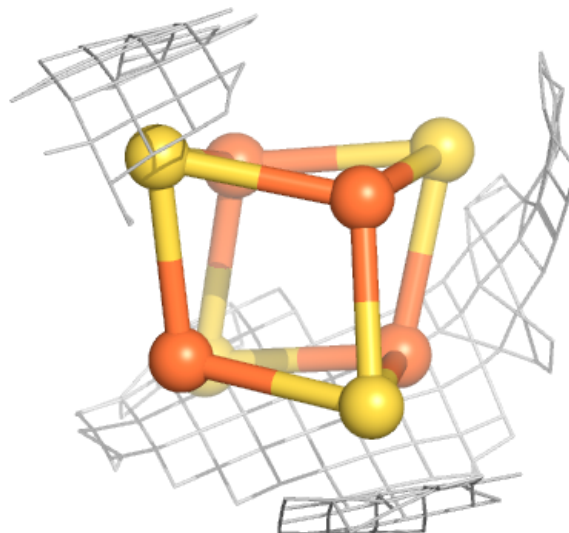
Electron density around SF4 H 601:

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



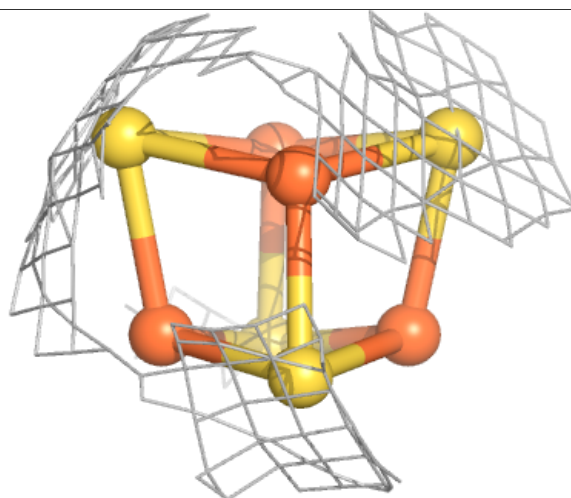
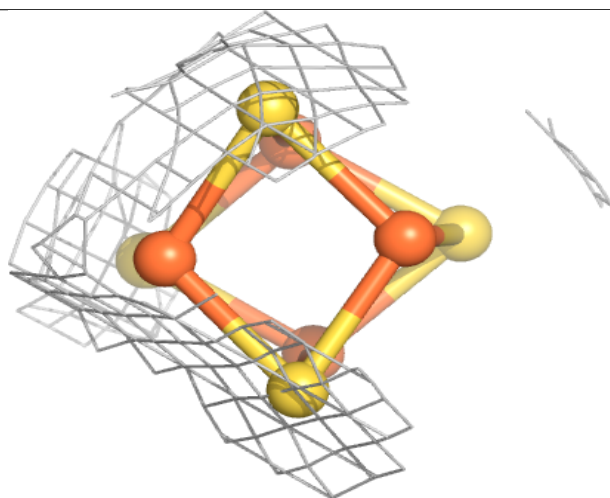
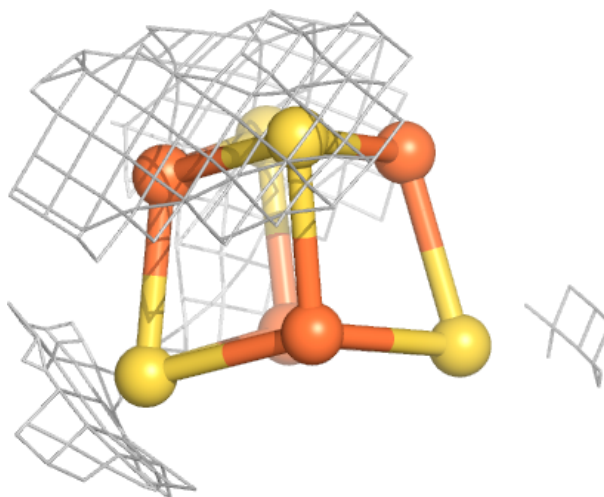
Electron density around SF4 C 601:

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and green (positive)



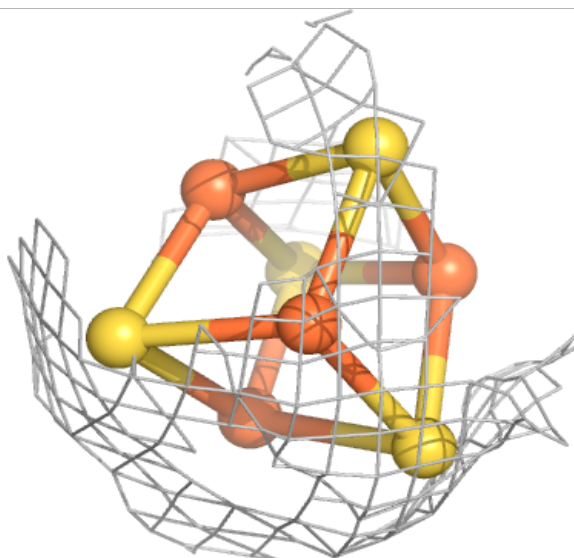
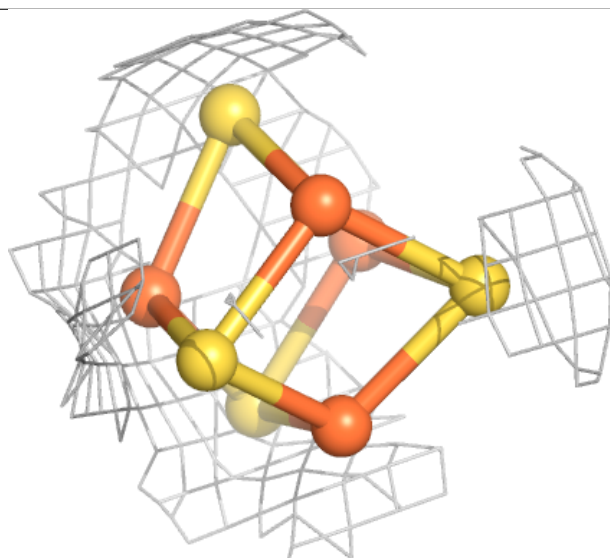
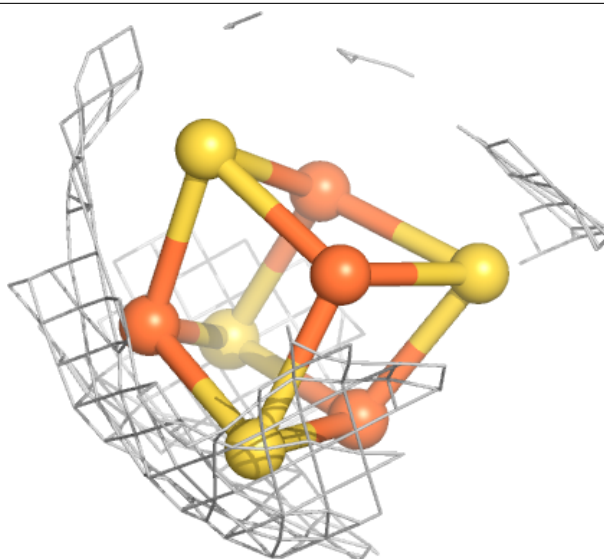
Electron density around SF4 G 601:

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



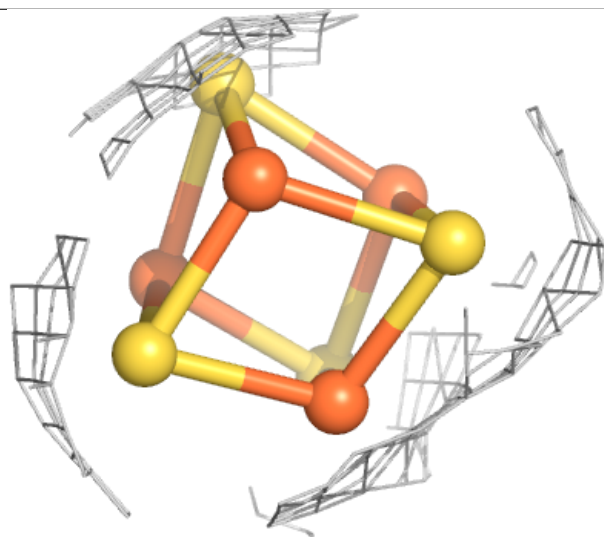
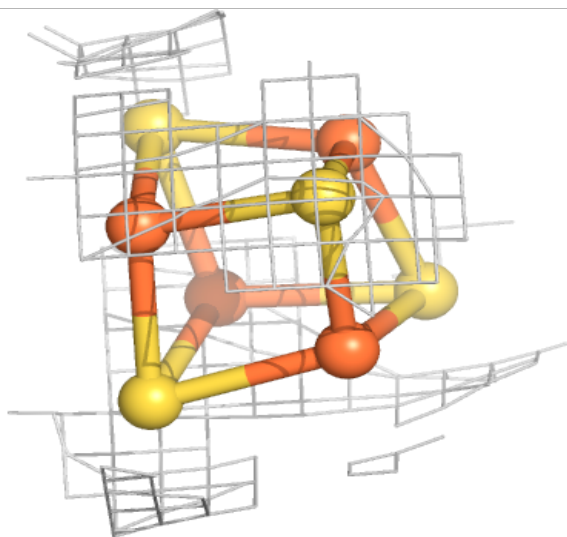
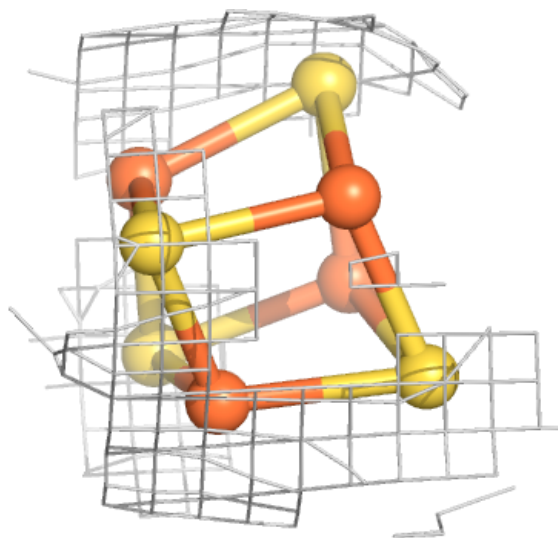
Electron density around SF4 F 601:

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



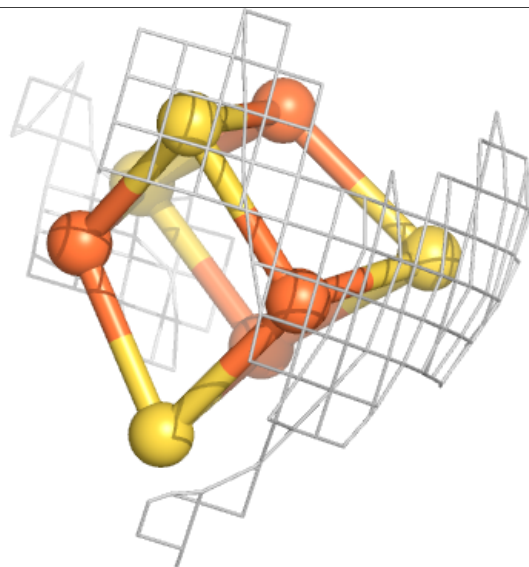
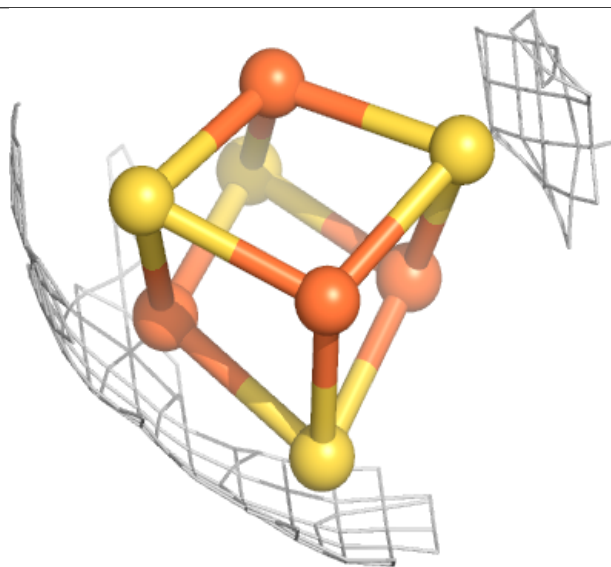
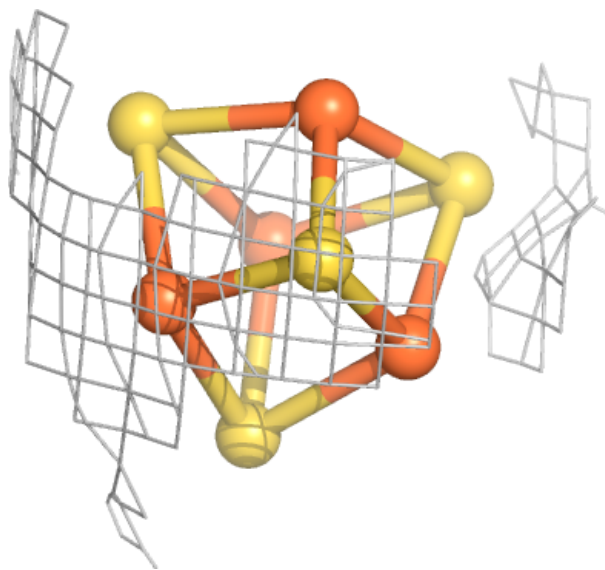
Electron density around SF4 B 601:

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and green (positive)



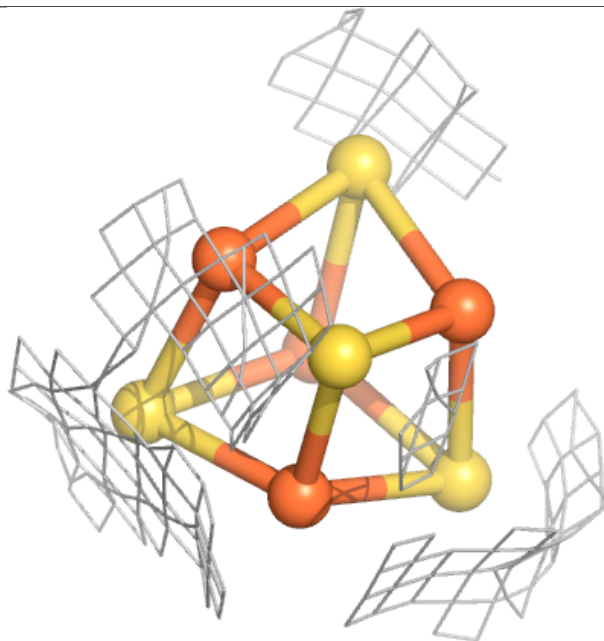
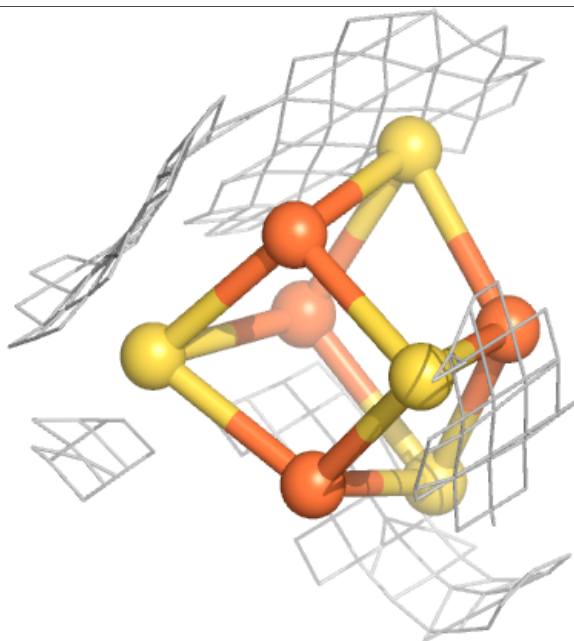
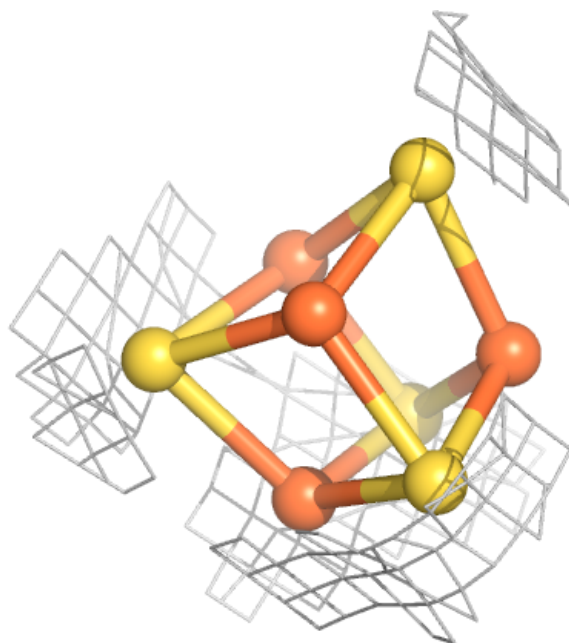
Electron density around SF4 A 601:

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



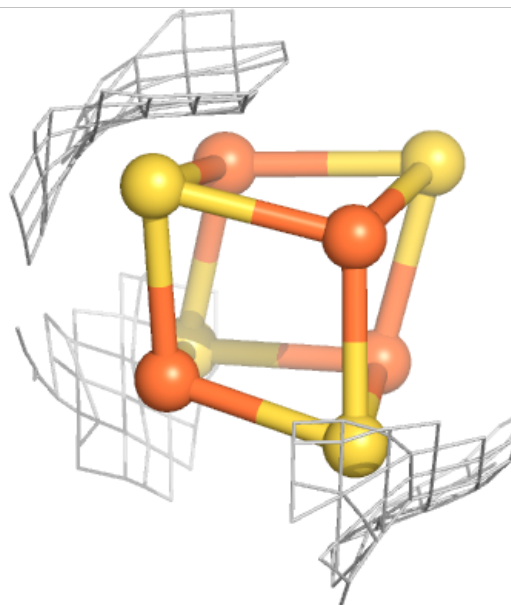
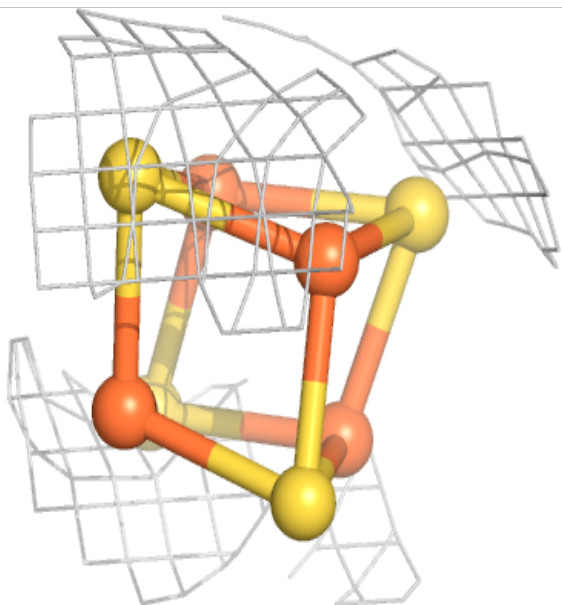
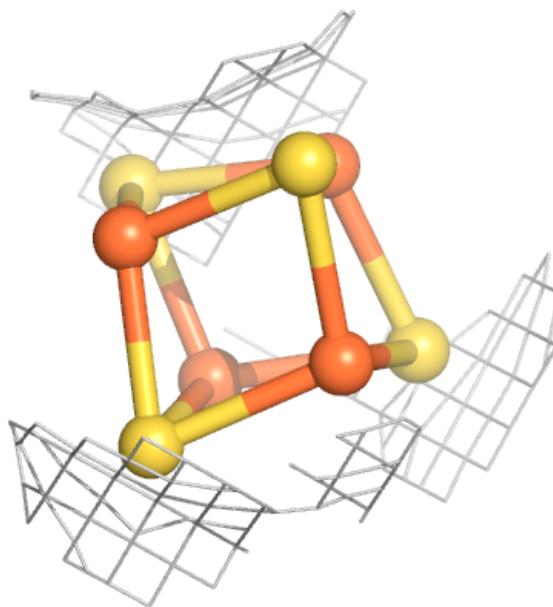
Electron density around SF4 D 601:

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



Electron density around SF4 E 601:

$2mF_o-DF_c$ (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.