



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 02:23 AM UTC

PDB ID : 8XVF / pdb_00008xvf
Title : Globular domain of Trichinella spiralis calreticulin
Authors : Zhu, X.P.; Jia, Z.H.; Yu, W.
Deposited on : 2024-01-15
Resolution : 2.76 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
Xtriage (Phenix)	:	2.0
EDS	:	3.0
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

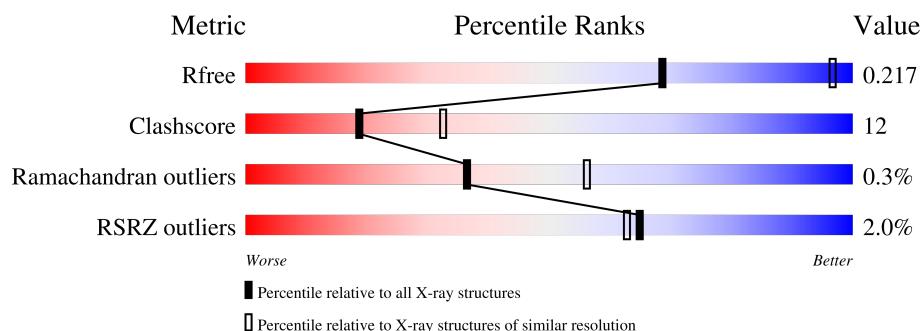
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.76 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	1009 (2.76-2.76)
Clashscore	190562	1044 (2.76-2.76)
Ramachandran outliers	187476	1024 (2.76-2.76)
RSRZ outliers	180081	1009 (2.76-2.76)

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	273	
1	B	273	
1	C	273	
1	D	273	
1	E	273	
1	F	273	

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Mol	Chain	Length	Quality of chain
1	G	273	62%28%8%
1	H	273	3% 63%27%8%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 16893 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 Calreticulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	251	Total 2061	C 1317	N 337	O 399	S 8	0	0	0
1	B	251	Total 2061	C 1317	N 337	O 399	S 8	0	0	0
1	C	251	Total 2061	C 1317	N 337	O 399	S 8	0	0	0
1	D	251	Total 2061	C 1317	N 337	O 399	S 8	0	0	0
1	E	251	Total 2061	C 1317	N 337	O 399	S 8	0	0	0
1	F	251	Total 2061	C 1317	N 337	O 399	S 8	0	0	0
1	G	251	Total 2061	C 1317	N 337	O 399	S 8	0	0	0
1	H	251	Total 2061	C 1317	N 337	O 399	S 8	0	0	0

There are 192 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	initiating methionine	UNP A0A0V1BA72
A	1	GLY	-	expression tag	UNP A0A0V1BA72
A	2	SER	-	expression tag	UNP A0A0V1BA72
A	3	SER	-	expression tag	UNP A0A0V1BA72
A	4	HIS	-	expression tag	UNP A0A0V1BA72
A	5	HIS	-	expression tag	UNP A0A0V1BA72
A	6	HIS	-	expression tag	UNP A0A0V1BA72
A	7	HIS	-	expression tag	UNP A0A0V1BA72
A	8	HIS	-	expression tag	UNP A0A0V1BA72
A	9	HIS	-	expression tag	UNP A0A0V1BA72
A	10	SER	-	expression tag	UNP A0A0V1BA72
A	11	SER	-	expression tag	UNP A0A0V1BA72
A	12	GLY	-	expression tag	UNP A0A0V1BA72

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Chain	Residue	Modelled	Actual	Comment	Reference
A	13	GLU	-	expression tag	UNP A0A0V1BA72
A	14	ASN	-	expression tag	UNP A0A0V1BA72
A	15	LEU	-	expression tag	UNP A0A0V1BA72
A	16	TYR	-	expression tag	UNP A0A0V1BA72
A	17	PHE	-	expression tag	UNP A0A0V1BA72
A	18	GLN	-	expression tag	UNP A0A0V1BA72
A	19	GLY	-	expression tag	UNP A0A0V1BA72
A	20	GLY	-	expression tag	UNP A0A0V1BA72
A	210	GLY	-	linker	UNP A0A0V1BA72
A	211	SER	-	linker	UNP A0A0V1BA72
A	212	GLY	-	linker	UNP A0A0V1BA72
B	0	MET	-	initiating methionine	UNP A0A0V1BA72
B	1	GLY	-	expression tag	UNP A0A0V1BA72
B	2	SER	-	expression tag	UNP A0A0V1BA72
B	3	SER	-	expression tag	UNP A0A0V1BA72
B	4	HIS	-	expression tag	UNP A0A0V1BA72
B	5	HIS	-	expression tag	UNP A0A0V1BA72
B	6	HIS	-	expression tag	UNP A0A0V1BA72
B	7	HIS	-	expression tag	UNP A0A0V1BA72
B	8	HIS	-	expression tag	UNP A0A0V1BA72
B	9	HIS	-	expression tag	UNP A0A0V1BA72
B	10	SER	-	expression tag	UNP A0A0V1BA72
B	11	SER	-	expression tag	UNP A0A0V1BA72
B	12	GLY	-	expression tag	UNP A0A0V1BA72
B	13	GLU	-	expression tag	UNP A0A0V1BA72
B	14	ASN	-	expression tag	UNP A0A0V1BA72
B	15	LEU	-	expression tag	UNP A0A0V1BA72
B	16	TYR	-	expression tag	UNP A0A0V1BA72
B	17	PHE	-	expression tag	UNP A0A0V1BA72
B	18	GLN	-	expression tag	UNP A0A0V1BA72
B	19	GLY	-	expression tag	UNP A0A0V1BA72
B	20	GLY	-	expression tag	UNP A0A0V1BA72
B	210	GLY	-	linker	UNP A0A0V1BA72
B	211	SER	-	linker	UNP A0A0V1BA72
B	212	GLY	-	linker	UNP A0A0V1BA72
C	0	MET	-	initiating methionine	UNP A0A0V1BA72
C	1	GLY	-	expression tag	UNP A0A0V1BA72
C	2	SER	-	expression tag	UNP A0A0V1BA72
C	3	SER	-	expression tag	UNP A0A0V1BA72
C	4	HIS	-	expression tag	UNP A0A0V1BA72
C	5	HIS	-	expression tag	UNP A0A0V1BA72
C	6	HIS	-	expression tag	UNP A0A0V1BA72

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Chain	Residue	Modelled	Actual	Comment	Reference
C	7	HIS	-	expression tag	UNP A0A0V1BA72
C	8	HIS	-	expression tag	UNP A0A0V1BA72
C	9	HIS	-	expression tag	UNP A0A0V1BA72
C	10	SER	-	expression tag	UNP A0A0V1BA72
C	11	SER	-	expression tag	UNP A0A0V1BA72
C	12	GLY	-	expression tag	UNP A0A0V1BA72
C	13	GLU	-	expression tag	UNP A0A0V1BA72
C	14	ASN	-	expression tag	UNP A0A0V1BA72
C	15	LEU	-	expression tag	UNP A0A0V1BA72
C	16	TYR	-	expression tag	UNP A0A0V1BA72
C	17	PHE	-	expression tag	UNP A0A0V1BA72
C	18	GLN	-	expression tag	UNP A0A0V1BA72
C	19	GLY	-	expression tag	UNP A0A0V1BA72
C	20	GLY	-	expression tag	UNP A0A0V1BA72
C	210	GLY	-	linker	UNP A0A0V1BA72
C	211	SER	-	linker	UNP A0A0V1BA72
C	212	GLY	-	linker	UNP A0A0V1BA72
D	0	MET	-	initiating methionine	UNP A0A0V1BA72
D	1	GLY	-	expression tag	UNP A0A0V1BA72
D	2	SER	-	expression tag	UNP A0A0V1BA72
D	3	SER	-	expression tag	UNP A0A0V1BA72
D	4	HIS	-	expression tag	UNP A0A0V1BA72
D	5	HIS	-	expression tag	UNP A0A0V1BA72
D	6	HIS	-	expression tag	UNP A0A0V1BA72
D	7	HIS	-	expression tag	UNP A0A0V1BA72
D	8	HIS	-	expression tag	UNP A0A0V1BA72
D	9	HIS	-	expression tag	UNP A0A0V1BA72
D	10	SER	-	expression tag	UNP A0A0V1BA72
D	11	SER	-	expression tag	UNP A0A0V1BA72
D	12	GLY	-	expression tag	UNP A0A0V1BA72
D	13	GLU	-	expression tag	UNP A0A0V1BA72
D	14	ASN	-	expression tag	UNP A0A0V1BA72
D	15	LEU	-	expression tag	UNP A0A0V1BA72
D	16	TYR	-	expression tag	UNP A0A0V1BA72
D	17	PHE	-	expression tag	UNP A0A0V1BA72
D	18	GLN	-	expression tag	UNP A0A0V1BA72
D	19	GLY	-	expression tag	UNP A0A0V1BA72
D	20	GLY	-	expression tag	UNP A0A0V1BA72
D	210	GLY	-	linker	UNP A0A0V1BA72
D	211	SER	-	linker	UNP A0A0V1BA72
D	212	GLY	-	linker	UNP A0A0V1BA72
E	0	MET	-	initiating methionine	UNP A0A0V1BA72

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Chain	Residue	Modelled	Actual	Comment	Reference
E	1	GLY	-	expression tag	UNP A0A0V1BA72
E	2	SER	-	expression tag	UNP A0A0V1BA72
E	3	SER	-	expression tag	UNP A0A0V1BA72
E	4	HIS	-	expression tag	UNP A0A0V1BA72
E	5	HIS	-	expression tag	UNP A0A0V1BA72
E	6	HIS	-	expression tag	UNP A0A0V1BA72
E	7	HIS	-	expression tag	UNP A0A0V1BA72
E	8	HIS	-	expression tag	UNP A0A0V1BA72
E	9	HIS	-	expression tag	UNP A0A0V1BA72
E	10	SER	-	expression tag	UNP A0A0V1BA72
E	11	SER	-	expression tag	UNP A0A0V1BA72
E	12	GLY	-	expression tag	UNP A0A0V1BA72
E	13	GLU	-	expression tag	UNP A0A0V1BA72
E	14	ASN	-	expression tag	UNP A0A0V1BA72
E	15	LEU	-	expression tag	UNP A0A0V1BA72
E	16	TYR	-	expression tag	UNP A0A0V1BA72
E	17	PHE	-	expression tag	UNP A0A0V1BA72
E	18	GLN	-	expression tag	UNP A0A0V1BA72
E	19	GLY	-	expression tag	UNP A0A0V1BA72
E	20	GLY	-	expression tag	UNP A0A0V1BA72
E	210	GLY	-	linker	UNP A0A0V1BA72
E	211	SER	-	linker	UNP A0A0V1BA72
E	212	GLY	-	linker	UNP A0A0V1BA72
F	0	MET	-	initiating methionine	UNP A0A0V1BA72
F	1	GLY	-	expression tag	UNP A0A0V1BA72
F	2	SER	-	expression tag	UNP A0A0V1BA72
F	3	SER	-	expression tag	UNP A0A0V1BA72
F	4	HIS	-	expression tag	UNP A0A0V1BA72
F	5	HIS	-	expression tag	UNP A0A0V1BA72
F	6	HIS	-	expression tag	UNP A0A0V1BA72
F	7	HIS	-	expression tag	UNP A0A0V1BA72
F	8	HIS	-	expression tag	UNP A0A0V1BA72
F	9	HIS	-	expression tag	UNP A0A0V1BA72
F	10	SER	-	expression tag	UNP A0A0V1BA72
F	11	SER	-	expression tag	UNP A0A0V1BA72
F	12	GLY	-	expression tag	UNP A0A0V1BA72
F	13	GLU	-	expression tag	UNP A0A0V1BA72
F	14	ASN	-	expression tag	UNP A0A0V1BA72
F	15	LEU	-	expression tag	UNP A0A0V1BA72
F	16	TYR	-	expression tag	UNP A0A0V1BA72
F	17	PHE	-	expression tag	UNP A0A0V1BA72
F	18	GLN	-	expression tag	UNP A0A0V1BA72

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Chain	Residue	Modelled	Actual	Comment	Reference
F	19	GLY	-	expression tag	UNP A0A0V1BA72
F	20	GLY	-	expression tag	UNP A0A0V1BA72
F	210	GLY	-	linker	UNP A0A0V1BA72
F	211	SER	-	linker	UNP A0A0V1BA72
F	212	GLY	-	linker	UNP A0A0V1BA72
G	0	MET	-	initiating methionine	UNP A0A0V1BA72
G	1	GLY	-	expression tag	UNP A0A0V1BA72
G	2	SER	-	expression tag	UNP A0A0V1BA72
G	3	SER	-	expression tag	UNP A0A0V1BA72
G	4	HIS	-	expression tag	UNP A0A0V1BA72
G	5	HIS	-	expression tag	UNP A0A0V1BA72
G	6	HIS	-	expression tag	UNP A0A0V1BA72
G	7	HIS	-	expression tag	UNP A0A0V1BA72
G	8	HIS	-	expression tag	UNP A0A0V1BA72
G	9	HIS	-	expression tag	UNP A0A0V1BA72
G	10	SER	-	expression tag	UNP A0A0V1BA72
G	11	SER	-	expression tag	UNP A0A0V1BA72
G	12	GLY	-	expression tag	UNP A0A0V1BA72
G	13	GLU	-	expression tag	UNP A0A0V1BA72
G	14	ASN	-	expression tag	UNP A0A0V1BA72
G	15	LEU	-	expression tag	UNP A0A0V1BA72
G	16	TYR	-	expression tag	UNP A0A0V1BA72
G	17	PHE	-	expression tag	UNP A0A0V1BA72
G	18	GLN	-	expression tag	UNP A0A0V1BA72
G	19	GLY	-	expression tag	UNP A0A0V1BA72
G	20	GLY	-	expression tag	UNP A0A0V1BA72
G	210	GLY	-	linker	UNP A0A0V1BA72
G	211	SER	-	linker	UNP A0A0V1BA72
G	212	GLY	-	linker	UNP A0A0V1BA72
H	0	MET	-	initiating methionine	UNP A0A0V1BA72
H	1	GLY	-	expression tag	UNP A0A0V1BA72
H	2	SER	-	expression tag	UNP A0A0V1BA72
H	3	SER	-	expression tag	UNP A0A0V1BA72
H	4	HIS	-	expression tag	UNP A0A0V1BA72
H	5	HIS	-	expression tag	UNP A0A0V1BA72
H	6	HIS	-	expression tag	UNP A0A0V1BA72
H	7	HIS	-	expression tag	UNP A0A0V1BA72
H	8	HIS	-	expression tag	UNP A0A0V1BA72
H	9	HIS	-	expression tag	UNP A0A0V1BA72
H	10	SER	-	expression tag	UNP A0A0V1BA72
H	11	SER	-	expression tag	UNP A0A0V1BA72
H	12	GLY	-	expression tag	UNP A0A0V1BA72

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Chain	Residue	Modelled	Actual	Comment	Reference
H	13	GLU	-	expression tag	UNP A0A0V1BA72
H	14	ASN	-	expression tag	UNP A0A0V1BA72
H	15	LEU	-	expression tag	UNP A0A0V1BA72
H	16	TYR	-	expression tag	UNP A0A0V1BA72
H	17	PHE	-	expression tag	UNP A0A0V1BA72
H	18	GLN	-	expression tag	UNP A0A0V1BA72
H	19	GLY	-	expression tag	UNP A0A0V1BA72
H	20	GLY	-	expression tag	UNP A0A0V1BA72
H	210	GLY	-	linker	UNP A0A0V1BA72
H	211	SER	-	linker	UNP A0A0V1BA72
H	212	GLY	-	linker	UNP A0A0V1BA72

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Ca 1 1	0	0
2	B	1	Total Ca 1 1	0	0
2	C	1	Total Ca 1 1	0	0
2	D	1	Total Ca 1 1	0	0
2	E	1	Total Ca 1 1	0	0
2	F	1	Total Ca 1 1	0	0
2	G	1	Total Ca 1 1	0	0
2	H	1	Total Ca 1 1	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	68	Total O 68 68	0	0
3	B	57	Total O 57 57	0	0
3	C	55	Total O 55 55	0	0

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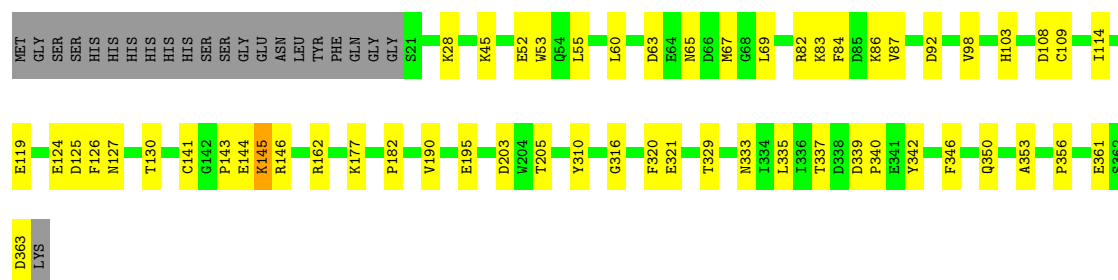
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	60	Total 60	O 60	0	0
3	E	41	Total 41	O 41	0	0
3	F	33	Total 33	O 33	0	0
3	G	44	Total 44	O 44	0	0
3	H	39	Total 39	O 39	0	0

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

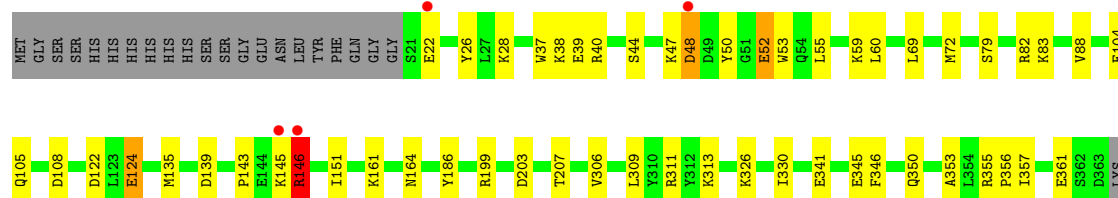
• Molecule 1: Calreticulin

Chain A: 



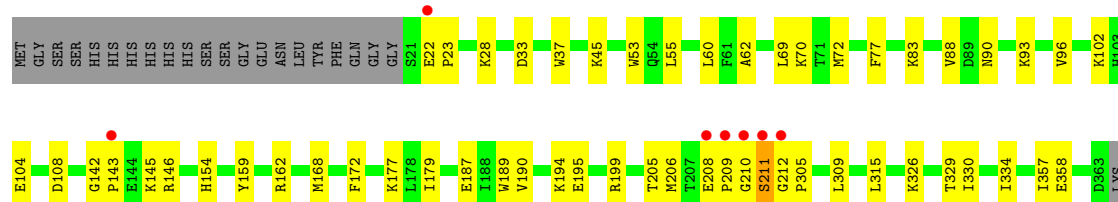
• Molecule 1: Calreticulin

Chain B: 



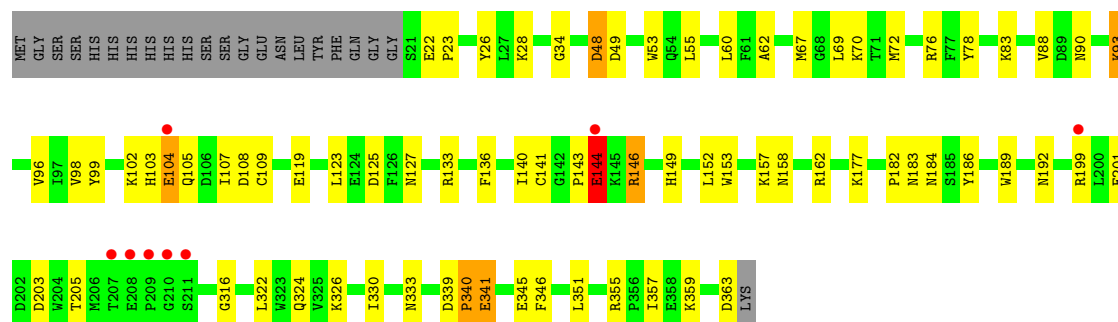
• Molecule 1: Calreticulin

Chain C: 

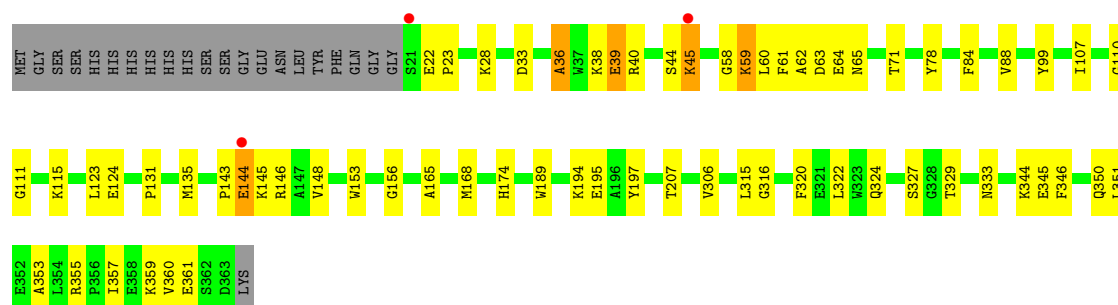


• Molecule 1: Calreticulin

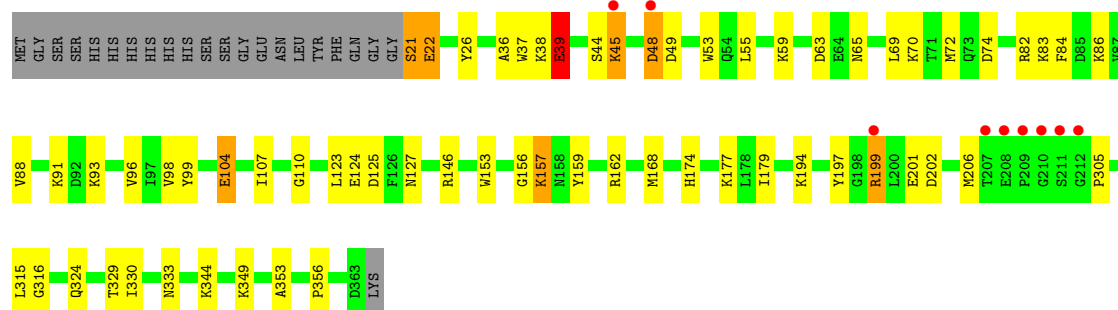
Chain D: 



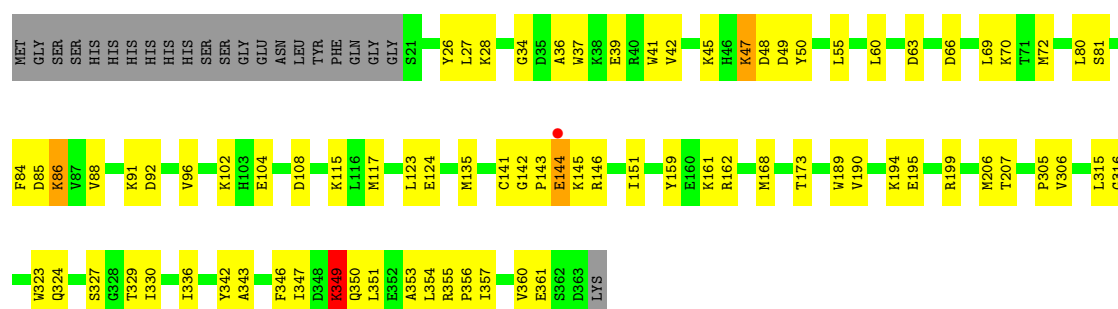
• Molecule 1: Calreticulin



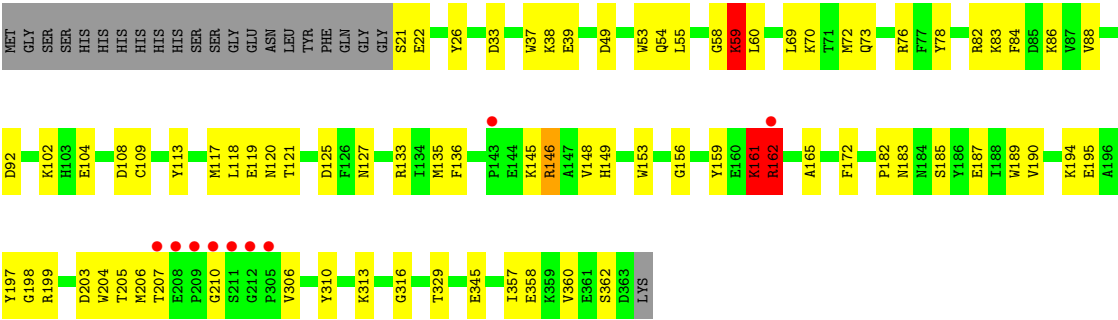
• Molecule 1: Calreticulin



• Molecule 1: Calreticulin



● Molecule 1: Calreticulin



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	83.52Å 154.07Å 100.78Å 90.00° 108.32° 90.00°	Depositor
Resolution (Å)	73.43 – 2.76 73.43 – 2.77	Depositor EDS
% Data completeness (in resolution range)	98.2 (73.43-2.76) 98.2 (73.43-2.77)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.33 (at 2.77Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.194 , 0.243 (Not available) , 0.217	Depositor DCC
R_{free} test set	2000 reflections (3.23%)	wwPDB-VP
Wilson B-factor (Å ²)	34.9	Xtriage
Anisotropy	0.425	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 41.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	16893	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.75% 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: CA

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	0.48	1/2115 (0.0%)	0.73	7/2860 (0.2%)
1	B	0.53	1/2115 (0.0%)	0.81	8/2860 (0.3%)
1	C	0.44	1/2115 (0.0%)	0.66	1/2860 (0.0%)
1	D	0.49	1/2115 (0.0%)	0.82	10/2860 (0.3%)
1	E	0.57	2/2115 (0.1%)	1.03	23/2860 (0.8%)
1	F	0.50	1/2115 (0.0%)	1.04	25/2860 (0.9%)
1	G	0.47	2/2115 (0.1%)	0.85	9/2860 (0.3%)
1	H	0.50	0/2115	1.02	19/2860 (0.7%)
All	All	0.50	9/16920 (0.1%)	0.88	102/22880 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	3
1	C	0	1
1	D	0	3
1	E	0	1
1	F	0	3
1	G	0	2
1	H	0	5
All	All	0	18

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	144	GLU	CB-CG	8.33	1.77	1.52
1	C	211	SER	CA-CB	7.23	1.64	1.53
1	G	86	LYS	CE-NZ	7.22	1.71	1.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	145	LYS	CE-NZ	5.96	1.67	1.49
1	G	86	LYS	CD-CE	5.64	1.69	1.52
1	F	39	GLU	CD-OE2	5.63	1.36	1.25
1	B	146	ARG	CB-CG	-5.32	1.36	1.52
1	E	359	LYS	CE-NZ	5.29	1.65	1.49
1	D	93	LYS	CE-NZ	5.12	1.64	1.49

All (102) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	161	LYS	CA-C-N	-15.46	94.42	122.09
1	H	161	LYS	C-N-CA	-15.46	94.42	122.09
1	F	38	LYS	CA-C-N	-12.21	100.32	122.54
1	F	38	LYS	C-N-CA	-12.21	100.32	122.54
1	G	144	GLU	N-CA-C	12.12	128.23	112.72
1	B	48	ASP	N-CA-C	-12.01	95.76	112.45
1	E	45	LYS	CA-CB-CG	11.86	137.81	114.10
1	E	58	GLY	CA-C-N	-11.44	103.73	120.89
1	E	58	GLY	C-N-CA	-11.44	103.73	120.89
1	F	45	LYS	CG-CD-CE	-10.78	86.50	111.30
1	H	162	ARG	CD-NE-CZ	10.74	139.43	124.40
1	E	59	LYS	CD-CE-NZ	-10.29	78.97	111.90
1	G	47	LYS	CB-CG-CD	-10.26	87.71	111.30
1	F	157	LYS	CB-CG-CD	10.21	134.79	111.30
1	H	59	LYS	CB-CG-CD	10.03	134.36	111.30
1	E	44	SER	CA-C-N	9.35	136.58	120.68
1	E	44	SER	C-N-CA	9.35	136.58	120.68
1	F	86	LYS	CB-CG-CD	-9.29	89.94	111.30
1	H	162	ARG	CA-CB-CG	-9.22	95.66	114.10
1	H	345	GLU	CB-CA-C	-9.13	96.54	110.88
1	D	340	PRO	CA-C-N	-8.42	107.36	121.92
1	D	340	PRO	C-N-CA	-8.42	107.36	121.92
1	G	146	ARG	CA-CB-CG	8.36	130.81	114.10
1	H	162	ARG	N-CA-C	-8.26	103.33	113.41
1	E	59	LYS	CB-CG-CD	-8.06	92.75	111.30
1	H	345	GLU	CA-CB-CG	8.03	130.17	114.10
1	F	21	SER	CA-C-N	-7.93	102.46	121.80
1	F	21	SER	C-N-CA	-7.93	102.46	121.80
1	E	144	GLU	CG-CD-OE2	-7.89	100.24	118.40
1	D	93	LYS	CB-CG-CD	-7.89	93.15	111.30
1	F	39	GLU	CA-CB-CG	7.87	129.84	114.10
1	G	144	GLU	N-CA-CB	-7.83	96.81	110.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	341	GLU	CA-CB-CG	-7.73	98.64	114.10
1	F	45	LYS	CB-CA-C	7.68	128.66	110.99
1	E	144	GLU	CG-CD-OE1	7.67	136.04	118.40
1	F	157	LYS	CG-CD-CE	-7.52	94.00	111.30
1	E	345	GLU	CA-CB-CG	7.45	129.01	114.10
1	D	157	LYS	CB-CG-CD	7.44	128.42	111.30
1	F	157	LYS	CD-CE-NZ	7.41	135.60	111.90
1	F	86	LYS	CA-CB-CG	7.38	128.86	114.10
1	D	48	ASP	CB-CA-C	-7.26	94.82	109.99
1	H	162	ARG	CB-CG-CD	7.24	127.95	111.30
1	B	48	ASP	CB-CG-OD2	-7.16	101.93	118.40
1	F	39	GLU	CB-CG-CD	7.16	124.77	112.60
1	G	48	ASP	CB-CG-OD2	-7.04	102.21	118.40
1	B	52	GLU	CB-CA-C	-6.99	96.59	109.46
1	E	39	GLU	CA-CB-CG	6.90	127.90	114.10
1	E	45	LYS	CB-CA-C	-6.88	96.35	110.38
1	E	45	LYS	CG-CD-CE	-6.74	95.81	111.30
1	E	39	GLU	CB-CG-CD	-6.69	101.23	112.60
1	E	39	GLU	N-CA-C	-6.68	105.29	113.50
1	H	145	LYS	CA-C-N	6.65	132.44	122.67
1	H	145	LYS	C-N-CA	6.65	132.44	122.67
1	F	157	LYS	N-CA-CB	-6.53	99.37	111.52
1	D	103	HIS	CA-C-N	-6.48	111.42	122.56
1	D	103	HIS	C-N-CA	-6.48	111.42	122.56
1	G	349	LYS	CA-CB-CG	6.35	126.79	114.10
1	A	86	LYS	CG-CD-CE	-6.20	97.04	111.30
1	F	22	GLU	CA-CB-CG	-6.18	101.75	114.10
1	F	45	LYS	CD-CE-NZ	6.13	131.51	111.90
1	G	144	GLU	CB-CA-C	6.11	120.39	110.37
1	F	104	GLU	CA-CB-CG	6.09	126.28	114.10
1	G	143	PRO	CA-C-N	6.07	133.39	122.46
1	G	143	PRO	C-N-CA	6.07	133.39	122.46
1	A	52	GLU	CA-CB-CG	6.02	126.13	114.10
1	F	86	LYS	CG-CD-CE	5.99	125.06	111.30
1	H	86	LYS	CA-CB-CG	5.95	125.99	114.10
1	E	145	LYS	CD-CE-NZ	-5.83	93.25	111.90
1	E	59	LYS	CA-CB-CG	5.78	125.65	114.10
1	B	22	GLU	N-CA-C	-5.68	100.78	109.64
1	D	104	GLU	CA-CB-CG	5.66	125.42	114.10
1	A	145	LYS	CA-CB-CG	5.60	125.30	114.10
1	H	162	ARG	CA-C-N	5.60	130.90	123.00
1	H	162	ARG	C-N-CA	5.60	130.90	123.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	45	LYS	CA-C-N	5.58	129.53	120.60
1	F	45	LYS	C-N-CA	5.58	129.53	120.60
1	H	86	LYS	CB-CA-C	5.56	119.26	109.80
1	B	146	ARG	CA-C-N	-5.53	113.97	122.87
1	B	146	ARG	C-N-CA	-5.53	113.97	122.87
1	E	344	LYS	CA-C-N	-5.48	112.51	120.29
1	E	344	LYS	C-N-CA	-5.48	112.51	120.29
1	E	144	GLU	CA-C-N	-5.43	114.96	122.30
1	E	144	GLU	C-N-CA	-5.43	114.96	122.30
1	H	86	LYS	N-CA-CB	-5.36	101.82	110.71
1	A	144	GLU	CB-CG-CD	5.33	121.66	112.60
1	B	48	ASP	CB-CG-OD1	5.31	130.62	118.40
1	B	124	GLU	CG-CD-OE2	-5.30	106.22	118.40
1	F	349	LYS	CB-CG-CD	-5.28	99.16	111.30
1	F	104	GLU	CB-CG-CD	-5.28	103.63	112.60
1	F	86	LYS	CD-CE-NZ	-5.23	95.17	111.90
1	E	59	LYS	N-CA-CB	5.22	117.52	110.10
1	C	211	SER	CB-CA-C	5.17	118.67	109.72
1	A	145	LYS	CA-C-N	-5.17	114.81	123.04
1	A	145	LYS	C-N-CA	-5.17	114.81	123.04
1	F	22	GLU	CB-CG-CD	5.15	121.36	112.60
1	H	162	ARG	NE-CZ-NH1	5.15	126.65	121.50
1	A	144	GLU	N-CA-C	5.14	118.02	111.69
1	F	48	ASP	CB-CA-C	-5.13	102.77	110.92
1	H	162	ARG	N-CA-CB	5.11	117.94	110.47
1	H	345	GLU	CG-CD-OE1	5.07	130.06	118.40
1	E	345	GLU	N-CA-CB	5.04	117.63	110.16
1	D	93	LYS	CD-CE-NZ	-5.02	95.83	111.90

There are no chirality outliers.

All (18) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	146	ARG	Sidechain
1	B	47	LYS	Peptide
1	B	48	ASP	Peptide
1	C	199	ARG	Sidechain
1	D	144	GLU	Peptide
1	D	146	ARG	Sidechain
1	D	199	ARG	Sidechain
1	E	195	GLU	Sidechain
1	F	199	ARG	Sidechain

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Mol	Chain	Res	Type	Group
1	F	39	GLU	Peptide
1	F	48	ASP	Peptide
1	G	144	GLU	Sidechain
1	G	199	ARG	Sidechain
1	H	146	ARG	Sidechain
1	H	161	LYS	Peptide
1	H	162	ARG	Sidechain
1	H	39	GLU	Peptide
1	H	58	GLY	Peptide

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2061	0	1968	34	0
1	B	2061	0	1968	40	0
1	C	2061	0	1968	47	0
1	D	2061	0	1968	48	0
1	E	2061	0	1968	50	0
1	F	2061	0	1968	43	0
1	G	2061	0	1968	66	0
1	H	2061	0	1968	66	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
3	A	68	0	0	6	0
3	B	57	0	0	9	1
3	C	55	0	0	11	0
3	D	60	0	0	5	1
3	E	41	0	0	9	0
3	F	33	0	0	7	0
3	G	44	0	0	8	0
3	H	39	0	0	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	16893	0	15744	385	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 12.

All (385) 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:144:GLU:CB	1:E:144:GLU:CG	1.77	1.57
1:G:86:LYS:NZ	1:G:86:LYS:CE	1.71	1.50
1:H:59:LYS:HG2	1:H:60:LEU:N	1.63	1.09
1:H:59:LYS:HG2	1:H:60:LEU:H	1.13	1.08
1:H:119:GLU:HG3	1:H:120:ASN:HD22	1.23	1.04
1:F:199:ARG:HH11	1:F:201:GLU:HB2	1.24	1.02
1:H:162:ARG:HB2	1:H:203:ASP:O	1.67	0.94
1:H:59:LYS:HB3	1:H:104:GLU:OE1	1.70	0.90
1:G:168:MET:O	3:G:501:HOH:O	1.92	0.88
1:F:45:LYS:NZ	3:F:501:HOH:O	1.84	0.86
1:B:122:ASP:OD1	1:B:124:GLU:OE1	1.92	0.85
1:F:74:ASP:O	3:F:502:HOH:O	1.94	0.85
1:E:329:THR:O	3:E:501:HOH:O	1.95	0.83
1:C:154:HIS:HB2	1:C:206:MET:HE3	1.58	0.82
1:F:162:ARG:NH2	1:F:202:ASP:O	2.11	0.82
1:H:59:LYS:CG	1:H:60:LEU:N	2.42	0.81
1:C:211:SER:O	3:C:501:HOH:O	1.98	0.81
1:E:123:LEU:O	3:E:502:HOH:O	1.99	0.81
1:C:142:GLY:O	1:C:146:ARG:NH1	2.12	0.81
1:H:161:LYS:HG2	1:H:162:ARG:H	1.46	0.78
1:H:121:THR:OG1	3:H:501:HOH:O	1.83	0.78
1:C:22:GLU:OE2	1:C:23:PRO:HD2	1.84	0.77
1:G:85:ASP:O	1:G:86:LYS:HG2	1.85	0.77
1:G:60:LEU:HD21	1:G:361:GLU:HG3	1.67	0.77
1:E:45:LYS:HB2	1:E:124:GLU:O	1.86	0.76
1:G:346:PHE:HA	1:G:349:LYS:HD3	1.67	0.76
1:A:55:LEU:HB3	1:A:67:MET:HE3	1.69	0.75
1:H:161:LYS:HG2	1:H:162:ARG:N	2.02	0.75
1:F:305:PRO:O	3:F:503:HOH:O	2.05	0.74
1:H:183:ASN:O	1:H:199:ARG:NH2	2.20	0.74
1:H:117:MET:HE2	1:H:121:THR:HG21	1.68	0.74
1:B:357:ILE:O	3:B:502:HOH:O	2.04	0.74
1:B:161:LYS:NZ	3:B:509:HOH:O	2.20	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:78:TYR:O	3:H:504:HOH:O	2.05	0.74
1:H:362:SER:O	3:H:505:HOH:O	2.06	0.73
1:B:143:PRO:O	1:B:146:ARG:NH2	2.22	0.72
1:C:162:ARG:NH1	3:C:505:HOH:O	2.22	0.72
1:A:340:PRO:O	3:A:501:HOH:O	2.07	0.72
1:E:39:GLU:HB2	3:E:504:HOH:O	1.90	0.72
1:D:192:ASN:OD1	3:D:501:HOH:O	2.08	0.72
1:D:22:GLU:O	3:D:502:HOH:O	2.08	0.72
1:G:47:LYS:HB2	1:G:50:TYR:CE2	2.25	0.71
1:E:36:ALA:O	1:E:40:ARG:NH1	2.24	0.71
1:D:143:PRO:O	1:D:146:ARG:NH1	2.22	0.70
1:H:119:GLU:HG3	1:H:120:ASN:ND2	2.01	0.70
1:A:130:THR:O	3:A:502:HOH:O	2.10	0.70
1:H:313:LYS:N	3:H:506:HOH:O	2.12	0.70
1:A:363:ASP:O	3:A:503:HOH:O	2.10	0.69
1:B:38:LYS:O	3:B:503:HOH:O	2.10	0.69
1:B:52:GLU:HB3	1:B:72:MET:HB2	1.72	0.69
1:B:164:ASN:OD1	3:B:504:HOH:O	2.10	0.69
1:C:22:GLU:OE2	1:C:23:PRO:CD	2.39	0.69
1:D:177:LYS:HB3	1:D:189:TRP:HB2	1.76	0.68
1:G:195:GLU:OE2	3:G:502:HOH:O	2.12	0.68
1:B:145:LYS:HG3	1:B:145:LYS:O	1.94	0.67
1:D:102:LYS:HE2	1:D:104:GLU:OE1	1.94	0.67
1:B:139:ASP:OD1	3:B:505:HOH:O	2.13	0.67
1:E:60:LEU:HD21	1:E:361:GLU:HG3	1.77	0.67
1:C:212:GLY:HA3	3:C:501:HOH:O	1.93	0.67
1:G:63:ASP:HB3	1:G:66:ASP:HB2	1.78	0.66
1:F:83:LYS:HD3	1:F:123:LEU:HD11	1.78	0.66
1:A:82:ARG:NH1	1:A:83:LYS:O	2.28	0.66
1:C:22:GLU:OE2	1:C:22:GLU:HA	1.95	0.66
1:H:108:ASP:OD1	1:H:109:CYS:N	2.30	0.65
1:C:33:ASP:OD1	3:C:503:HOH:O	2.14	0.65
1:A:92:ASP:OD2	3:A:504:HOH:O	2.14	0.64
1:B:40:ARG:O	3:B:507:HOH:O	2.14	0.64
1:A:146:ARG:NH1	3:A:509:HOH:O	2.30	0.64
1:F:156:GLY:O	1:F:157:LYS:HG2	1.98	0.64
1:F:199:ARG:HH11	1:F:201:GLU:CB	2.07	0.63
1:G:142:GLY:HA2	1:H:197:TYR:CE2	2.33	0.63
1:D:201:GLU:OE2	1:D:201:GLU:N	2.22	0.63
1:G:207:THR:HG21	1:G:306:VAL:HG21	1.81	0.63
1:C:168:MET:O	3:C:504:HOH:O	2.16	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:45:LYS:HE3	1:F:124:GLU:HB3	1.81	0.62
1:H:205:THR:O	1:H:205:THR:HG22	2.00	0.61
1:H:125:ASP:O	1:H:127:ASN:ND2	2.31	0.61
1:G:36:ALA:O	1:G:39:GLU:HG2	2.02	0.60
1:B:309:LEU:O	3:B:508:HOH:O	2.15	0.60
1:G:104:GLU:OE2	3:G:503:HOH:O	2.16	0.60
1:G:145:LYS:HE2	1:H:199:ARG:HD2	1.84	0.60
1:E:59:LYS:O	1:E:61:PHE:HD1	1.84	0.60
1:G:353:ALA:O	1:G:356:PRO:HD2	2.03	0.59
1:B:207:THR:HG21	1:B:306:VAL:HG21	1.85	0.59
1:D:70:LYS:HD3	1:D:72:MET:SD	2.43	0.59
1:F:49:ASP:OD1	1:F:49:ASP:N	2.36	0.59
1:D:133:ARG:NH1	3:D:505:HOH:O	2.23	0.59
1:E:131:PRO:HA	3:E:520:HOH:O	2.02	0.59
1:D:90:ASN:HA	1:D:93:LYS:HD2	1.85	0.58
1:F:59:LYS:HE3	1:F:104:GLU:OE1	2.03	0.58
1:D:107:ILE:HD11	1:D:324:GLN:HG2	1.85	0.58
1:B:122:ASP:CG	1:B:124:GLU:OE1	2.46	0.58
1:B:145:LYS:C	1:B:146:ARG:HG3	2.27	0.58
1:G:85:ASP:C	1:G:86:LYS:HG2	2.28	0.58
1:H:26:TYR:CE2	1:H:88:VAL:HA	2.39	0.58
1:F:88:VAL:HB	1:F:315:LEU:HB3	1.84	0.58
1:F:199:ARG:NH1	1:F:201:GLU:HB2	2.08	0.58
1:C:162:ARG:CZ	1:C:205:THR:HG22	2.35	0.57
1:G:27:LEU:HD13	1:G:84:PHE:HA	1.87	0.57
1:A:84:PHE:CD2	1:A:316:GLY:HA2	2.38	0.57
1:E:22:GLU:HG3	1:E:23:PRO:HD2	1.87	0.57
1:C:143:PRO:HA	1:C:146:ARG:HH12	1.70	0.57
1:E:148:VAL:HB	1:E:165:ALA:HB3	1.86	0.56
1:F:21:SER:OG	1:F:22:GLU:N	2.34	0.56
1:B:341:GLU:O	1:B:345:GLU:HG3	2.05	0.56
1:E:71:THR:N	3:E:501:HOH:O	2.23	0.56
1:D:140:ILE:HG23	1:D:146:ARG:HH21	1.69	0.56
1:F:70:LYS:HD3	1:F:72:MET:SD	2.46	0.56
1:F:96:VAL:HG22	1:F:179:ILE:HG23	1.88	0.56
1:E:351:LEU:HD23	1:E:355:ARG:NE	2.21	0.56
1:H:190:VAL:HG23	1:H:195:GLU:HG3	1.87	0.56
1:E:39:GLU:N	3:E:504:HOH:O	2.20	0.56
1:F:125:ASP:O	1:F:127:ASN:ND2	2.35	0.55
1:H:159:TYR:CD2	1:H:206:MET:HG2	2.41	0.55
1:C:145:LYS:O	1:C:146:ARG:HD3	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:49:ASP:O	3:G:504:HOH:O	2.18	0.55
1:B:60:LEU:HD11	1:B:361:GLU:HG3	1.89	0.55
1:A:141:CYS:HB3	1:A:145:LYS:HD3	1.89	0.54
1:D:351:LEU:O	1:D:355:ARG:HG2	2.07	0.54
1:E:144:GLU:CB	1:E:144:GLU:OE2	2.55	0.54
1:E:353:ALA:O	1:E:357:ILE:HD12	2.07	0.54
1:G:28:LYS:NZ	1:G:342:TYR:OH	2.38	0.54
1:B:59:LYS:HE2	1:B:104:GLU:HB2	1.89	0.54
1:A:45:LYS:HE3	1:A:124:GLU:CD	2.32	0.54
1:B:135:MET:HB3	1:B:151:ILE:HB	1.89	0.54
1:F:39:GLU:O	1:F:39:GLU:HG2	2.07	0.54
1:C:96:VAL:HG13	1:C:179:ILE:HG12	1.88	0.54
1:G:47:LYS:HB2	1:G:50:TYR:CZ	2.42	0.54
1:D:62:ALA:HB2	1:D:357:ILE:HD12	1.90	0.54
1:G:72:MET:HA	1:G:72:MET:HE3	1.89	0.53
1:E:351:LEU:HB3	1:E:355:ARG:HG2	1.91	0.53
1:F:37:TRP:CD1	1:F:55:LEU:HD11	2.43	0.53
1:H:70:LYS:HA	1:H:329:THR:O	2.09	0.53
1:H:182:PRO:HA	1:H:310:TYR:CD2	2.43	0.53
1:A:98:VAL:HG23	1:A:177:LYS:HB2	1.90	0.53
1:B:207:THR:CG2	1:B:306:VAL:HG21	2.38	0.53
1:G:207:THR:CG2	1:G:306:VAL:HG21	2.39	0.53
1:D:78:TYR:HB2	1:D:322:LEU:HD12	1.91	0.53
1:H:21:SER:OG	1:H:22:GLU:N	2.42	0.53
1:E:45:LYS:HG2	1:E:124:GLU:HG2	1.91	0.53
1:F:36:ALA:O	3:F:504:HOH:O	2.19	0.53
1:F:82:ARG:NH1	1:F:83:LYS:O	2.38	0.53
1:F:146:ARG:NE	3:F:505:HOH:O	2.24	0.53
1:H:357:ILE:O	1:H:360:VAL:HG22	2.09	0.53
1:G:70:LYS:NZ	3:G:503:HOH:O	2.26	0.52
1:D:60:LEU:HD22	1:D:357:ILE:HG22	1.92	0.52
1:E:144:GLU:CG	1:E:144:GLU:CA	2.82	0.52
1:G:357:ILE:O	1:G:360:VAL:HG22	2.10	0.52
1:B:69:LEU:O	1:B:330:ILE:HA	2.10	0.52
1:E:78:TYR:O	3:E:503:HOH:O	2.19	0.51
1:A:28:LYS:NZ	1:A:342:TYR:OH	2.43	0.51
1:H:161:LYS:HG3	1:H:204:TRP:CZ2	2.45	0.51
1:D:55:LEU:HB3	1:D:67:MET:HE3	1.92	0.51
1:F:70:LYS:HA	1:F:329:THR:O	2.11	0.51
1:F:99:TYR:HA	1:F:333:ASN:O	2.11	0.51
1:C:60:LEU:O	1:C:102:LYS:NZ	2.43	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:70:LYS:HA	1:C:329:THR:O	2.12	0.50
1:E:22:GLU:HG3	1:E:23:PRO:CD	2.42	0.50
1:A:126:PHE:CE2	1:A:321:GLU:HB2	2.47	0.50
1:A:162:ARG:NH2	1:A:205:THR:HG22	2.27	0.50
1:E:84:PHE:CD2	1:E:316:GLY:HA2	2.47	0.50
1:F:168:MET:HG3	1:F:174:HIS:CE1	2.46	0.50
1:E:144:GLU:CB	1:E:144:GLU:CD	2.77	0.50
1:G:47:LYS:HD3	1:G:49:ASP:OD1	2.11	0.50
1:H:189:TRP:CE2	1:H:194:LYS:HG3	2.47	0.50
1:D:133:ARG:HD3	1:D:153:TRP:CG	2.47	0.50
1:G:346:PHE:O	1:G:349:LYS:HB3	2.11	0.50
1:A:103:HIS:ND1	1:A:329:THR:OG1	2.37	0.50
1:D:48:ASP:N	1:D:48:ASP:OD1	2.45	0.50
1:D:93:LYS:O	1:D:182:PRO:HG3	2.11	0.50
1:E:144:GLU:OE2	1:E:144:GLU:HB3	2.11	0.49
1:G:34:GLY:N	3:G:509:HOH:O	2.44	0.49
1:H:53:TRP:CE3	1:H:69:LEU:HG	2.47	0.49
1:H:207:THR:HG21	1:H:306:VAL:HG21	1.94	0.49
1:A:60:LEU:HD11	1:A:361:GLU:HG2	1.94	0.49
1:C:70:LYS:HD3	1:C:72:MET:SD	2.52	0.49
1:E:115:LYS:HE2	1:E:135:MET:HE2	1.94	0.49
1:C:37:TRP:CE2	1:C:55:LEU:HD13	2.47	0.49
1:D:359:LYS:O	1:D:363:ASP:OD1	2.31	0.49
1:G:88:VAL:HB	1:G:315:LEU:HB3	1.94	0.49
1:H:207:THR:CG2	1:H:306:VAL:HG21	2.43	0.49
1:G:69:LEU:O	1:G:330:ILE:HA	2.13	0.49
1:H:54:GLN:HB3	1:H:72:MET:SD	2.53	0.49
1:D:108:ASP:OD1	1:D:109:CYS:N	2.42	0.49
1:E:189:TRP:CE2	1:E:194:LYS:HG3	2.47	0.49
1:D:98:VAL:HB	1:D:177:LYS:HG3	1.95	0.49
1:F:194:LYS:HD3	1:F:197:TYR:HB3	1.95	0.49
1:A:337:THR:HG23	1:A:339:ASP:H	1.78	0.49
1:A:190:VAL:HG23	1:A:195:GLU:HG3	1.94	0.48
1:E:111:GLY:C	1:E:322:LEU:HD23	2.37	0.48
1:D:83:LYS:HE3	1:D:123:LEU:HD13	1.95	0.48
1:H:37:TRP:CD1	1:H:55:LEU:HD11	2.47	0.48
1:E:99:TYR:HA	1:E:333:ASN:O	2.13	0.48
1:C:190:VAL:HG23	1:C:195:GLU:HG3	1.95	0.48
1:A:125:ASP:O	1:A:127:ASN:ND2	2.45	0.48
1:G:84:PHE:CD2	1:G:316:GLY:HA2	2.48	0.48
1:B:353:ALA:O	1:B:356:PRO:HD2	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:28:LYS:HD2	1:D:346:PHE:CE1	2.49	0.48
1:D:162:ARG:NH1	1:D:205:THR:HG23	2.29	0.48
1:A:143:PRO:HA	1:A:146:ARG:HH21	1.77	0.48
1:E:88:VAL:HB	1:E:315:LEU:HB3	1.94	0.48
1:F:53:TRP:CE3	1:F:69:LEU:HG	2.49	0.48
1:G:189:TRP:CZ2	1:G:194:LYS:HG3	2.49	0.48
1:A:203:ASP:HB3	3:A:513:HOH:O	2.13	0.47
1:H:189:TRP:CZ2	1:H:194:LYS:HG3	2.50	0.47
1:H:82:ARG:NH1	1:H:83:LYS:O	2.34	0.47
1:H:161:LYS:CG	1:H:162:ARG:N	2.73	0.47
1:B:28:LYS:HD2	1:B:346:PHE:CE1	2.49	0.47
1:D:119:GLU:HB3	1:D:316:GLY:HA3	1.95	0.47
1:E:357:ILE:O	1:E:360:VAL:HG12	2.15	0.47
1:F:353:ALA:O	1:F:356:PRO:HD2	2.15	0.47
1:G:81:SER:OG	1:G:123:LEU:HD22	2.15	0.47
1:C:143:PRO:CA	1:C:146:ARG:HH12	2.28	0.47
1:E:320:PHE:HB3	1:E:322:LEU:CD1	2.44	0.47
1:F:159:TYR:CG	1:F:206:MET:HG2	2.50	0.47
1:G:70:LYS:HA	1:G:329:THR:O	2.15	0.47
1:H:49:ASP:OD1	1:H:49:ASP:N	2.44	0.47
1:C:208:GLU:CD	1:C:209:PRO:HD2	2.39	0.47
1:A:126:PHE:CZ	1:A:321:GLU:HB2	2.50	0.47
1:A:353:ALA:O	1:A:356:PRO:HD2	2.14	0.47
1:D:34:GLY:O	3:D:504:HOH:O	2.21	0.47
1:G:28:LYS:HD2	1:G:346:PHE:CE1	2.50	0.47
1:C:62:ALA:HA	1:C:357:ILE:HD12	1.96	0.46
1:C:172:PHE:HB3	1:C:358:GLU:OE1	2.14	0.46
1:D:125:ASP:O	1:D:127:ASN:ND2	2.42	0.46
1:D:136:PHE:HA	1:D:149:HIS:O	2.15	0.46
1:G:190:VAL:HG23	1:G:195:GLU:HG3	1.97	0.46
1:A:182:PRO:HA	1:A:310:TYR:CD2	2.50	0.46
1:C:108:ASP:OD1	1:C:326:LYS:HB2	2.16	0.46
1:H:37:TRP:CZ2	1:H:38:LYS:HG2	2.51	0.46
1:E:324:GLN:NE2	1:E:327:SER:HA	2.31	0.46
1:B:53:TRP:CE3	1:B:69:LEU:HG	2.50	0.46
1:D:23:PRO:HB3	1:D:339:ASP:CB	2.45	0.46
1:A:53:TRP:CE3	1:A:69:LEU:HG	2.51	0.46
1:C:177:LYS:HB3	1:C:189:TRP:HB2	1.98	0.46
1:C:28:LYS:HA	1:C:334:ILE:O	2.16	0.46
1:C:70:LYS:NZ	3:C:508:HOH:O	2.25	0.46
1:D:99:TYR:HA	1:D:333:ASN:O	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:349:LYS:HB3	1:G:350:GLN:H	1.54	0.46
1:C:60:LEU:HD11	1:C:357:ILE:CG2	2.46	0.46
1:G:49:ASP:OD1	1:G:49:ASP:N	2.48	0.46
1:F:110:GLY:HA3	1:F:324:GLN:HG2	1.98	0.46
1:H:33:ASP:OD1	1:H:33:ASP:N	2.42	0.46
1:C:69:LEU:O	1:C:330:ILE:HA	2.15	0.45
1:E:107:ILE:HD11	1:E:110:GLY:HA3	1.96	0.45
1:E:153:TRP:CZ2	1:E:156:GLY:HA2	2.50	0.45
1:H:84:PHE:CD2	1:H:316:GLY:HA2	2.51	0.45
1:C:206:MET:HB2	3:C:538:HOH:O	2.15	0.45
1:D:177:LYS:NZ	3:D:503:HOH:O	2.12	0.45
1:G:47:LYS:NZ	1:H:92:ASP:HB3	2.31	0.45
1:D:69:LEU:O	1:D:330:ILE:HA	2.15	0.45
1:G:45:LYS:HD3	1:G:124:GLU:OE2	2.16	0.45
1:G:162:ARG:HB2	3:G:507:HOH:O	2.17	0.45
1:G:347:ILE:HD13	1:G:347:ILE:HA	1.82	0.45
1:H:159:TYR:CG	1:H:206:MET:HG2	2.52	0.45
1:H:161:LYS:HG3	1:H:204:TRP:CH2	2.52	0.45
1:C:206:MET:HE2	1:C:309:LEU:HD11	1.98	0.45
1:E:60:LEU:HD21	1:E:361:GLU:CG	2.46	0.45
1:E:36:ALA:C	3:E:504:HOH:O	2.59	0.45
1:G:42:VAL:C	1:G:80:LEU:HD12	2.42	0.45
1:C:45:LYS:HE2	1:C:45:LYS:HB3	1.82	0.45
1:B:82:ARG:NH1	1:B:83:LYS:O	2.43	0.45
1:D:341:GLU:O	1:D:345:GLU:HG3	2.16	0.45
1:A:114:ILE:HG22	1:A:320:PHE:CE1	2.52	0.45
1:C:83:LYS:NZ	3:C:509:HOH:O	2.25	0.45
1:E:333:ASN:ND2	1:E:350:GLN:HE21	2.15	0.44
1:F:44:SER:OG	1:F:45:LYS:N	2.50	0.44
1:B:311:ARG:NH2	1:B:313:LYS:HD2	2.32	0.44
1:C:208:GLU:HA	3:C:515:HOH:O	2.18	0.44
1:B:186:TYR:OH	1:B:203:ASP:OD2	2.29	0.44
1:E:33:ASP:CG	1:E:36:ALA:HB3	2.42	0.44
1:G:37:TRP:NE1	1:G:55:LEU:HD13	2.32	0.44
1:G:145:LYS:HD3	1:H:199:ARG:HG3	2.00	0.44
1:C:159:TYR:CG	1:C:206:MET:HG2	2.53	0.44
1:F:84:PHE:CD2	1:F:316:GLY:HA2	2.52	0.44
1:F:199:ARG:HE	1:F:202:ASP:CG	2.26	0.44
1:H:135:MET:HE2	1:H:149:HIS:CG	2.53	0.44
1:C:88:VAL:HB	1:C:315:LEU:HB3	2.00	0.44
1:G:142:GLY:HA2	1:H:197:TYR:CZ	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:194:LYS:HD3	1:E:197:TYR:HB3	1.99	0.44
1:G:42:VAL:O	1:G:80:LEU:HD12	2.18	0.44
1:B:52:GLU:CB	1:B:72:MET:HB2	2.46	0.44
1:C:208:GLU:N	3:C:515:HOH:O	2.51	0.44
1:A:162:ARG:HH21	1:A:205:THR:HG22	1.83	0.43
1:E:59:LYS:O	1:E:60:LEU:C	2.57	0.43
1:E:63:ASP:C	1:E:65:ASN:H	2.26	0.43
1:G:60:LEU:CD2	1:G:361:GLU:HG3	2.43	0.43
1:C:53:TRP:CE3	1:C:69:LEU:HG	2.54	0.43
1:D:144:GLU:O	1:D:144:GLU:CG	2.66	0.43
1:E:61:PHE:HE2	1:E:64:GLU:HB2	1.83	0.43
1:G:91:LYS:HE2	1:G:92:ASP:OD2	2.18	0.43
1:D:53:TRP:HB3	1:D:69:LEU:HD11	2.00	0.43
1:D:23:PRO:HB3	1:D:339:ASP:HB2	1.99	0.43
1:E:71:THR:HG22	1:E:78:TYR:CE2	2.53	0.43
1:H:37:TRP:CH2	1:H:38:LYS:HE2	2.54	0.43
1:H:136:PHE:HA	1:H:149:HIS:O	2.19	0.43
1:B:105:GLN:O	1:B:326:LYS:NZ	2.46	0.43
1:B:199:ARG:HE	1:B:199:ARG:HB3	1.63	0.43
1:G:115:LYS:O	1:G:117:MET:HE2	2.18	0.43
1:D:105:GLN:O	1:D:326:LYS:HE2	2.18	0.43
1:G:47:LYS:NZ	1:H:92:ASP:CB	2.82	0.43
1:H:59:LYS:HB3	1:H:104:GLU:CD	2.38	0.43
1:H:362:SER:C	3:H:505:HOH:O	2.56	0.43
1:G:159:TYR:CG	1:G:206:MET:HG2	2.54	0.42
1:G:323:TRP:CD1	1:G:323:TRP:C	2.97	0.42
1:H:26:TYR:CD2	1:H:88:VAL:HG22	2.54	0.42
1:H:148:VAL:HB	1:H:165:ALA:HB3	2.01	0.42
1:E:168:MET:HG3	1:E:174:HIS:ND1	2.33	0.42
1:F:63:ASP:C	1:F:65:ASN:H	2.26	0.42
1:H:113:TYR:CD1	1:H:135:MET:HE3	2.55	0.42
1:D:186:TYR:OH	1:D:203:ASP:OD2	2.17	0.42
1:F:344:LYS:HE3	3:F:511:HOH:O	2.20	0.42
1:G:161:LYS:NZ	3:G:507:HOH:O	2.38	0.42
1:A:87:VAL:HG21	1:A:119:GLU:OE1	2.19	0.42
1:B:355:ARG:O	3:B:510:HOH:O	2.21	0.42
1:H:172:PHE:HB3	1:H:358:GLU:OE1	2.20	0.42
1:A:346:PHE:O	1:A:350:GLN:HG2	2.19	0.42
1:B:26:TYR:CG	1:B:88:VAL:HG22	2.54	0.42
1:D:357:ILE:H	1:D:357:ILE:HG13	1.70	0.42
1:E:143:PRO:HA	1:E:146:ARG:HE	1.85	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:26:TYR:CE1	1:F:88:VAL:HA	2.54	0.42
1:H:153:TRP:CZ2	1:H:156:GLY:HA2	2.55	0.42
1:C:77:PHE:CZ	1:G:305:PRO:HD2	2.55	0.42
1:G:102:LYS:HE3	1:G:173:THR:OG1	2.19	0.42
1:H:102:LYS:HD3	1:H:104:GLU:HB3	2.01	0.42
1:H:194:LYS:HD3	1:H:197:TYR:HB3	2.02	0.42
1:A:45:LYS:HG2	1:B:39:GLU:HG3	2.02	0.42
1:B:50:TYR:CE1	1:B:79:SER:HB2	2.55	0.42
1:D:53:TRP:CE3	1:D:69:LEU:HG	2.55	0.42
1:A:333:ASN:ND2	1:A:350:GLN:HE21	2.18	0.42
1:F:91:LYS:O	1:F:93:LYS:HG3	2.20	0.42
1:A:63:ASP:C	1:A:65:ASN:H	2.28	0.41
1:B:346:PHE:O	1:B:350:GLN:HG2	2.20	0.41
1:F:153:TRP:CZ2	1:F:156:GLY:HA2	2.55	0.41
1:G:70:LYS:HD3	1:G:72:MET:SD	2.59	0.41
1:A:108:ASP:OD1	1:A:109:CYS:N	2.49	0.41
1:H:185:SER:HB2	1:H:198:GLY:O	2.20	0.41
1:B:44:SER:HB3	1:B:79:SER:HB3	2.02	0.41
1:C:187:GLU:OE2	1:C:194:LYS:HE2	2.21	0.41
1:F:344:LYS:CE	3:F:511:HOH:O	2.68	0.41
1:B:108:ASP:OD1	3:B:512:HOH:O	2.22	0.41
1:G:145:LYS:HG3	1:G:145:LYS:O	2.19	0.41
1:C:22:GLU:OE2	1:C:23:PRO:HD3	2.20	0.41
1:F:69:LEU:O	1:F:330:ILE:HA	2.20	0.41
1:G:96:VAL:HG11	1:G:343:ALA:CB	2.51	0.41
1:B:37:TRP:CE2	1:B:55:LEU:HD13	2.55	0.41
1:C:90:ASN:HA	1:C:93:LYS:HG3	2.03	0.41
1:D:152:LEU:O	1:D:158:ASN:HA	2.19	0.41
1:F:107:ILE:HD11	1:F:110:GLY:HA3	2.03	0.41
1:G:26:TYR:CD2	1:G:336:ILE:HG22	2.56	0.41
1:H:133:ARG:HA	1:H:133:ARG:HD3	1.42	0.41
1:H:146:ARG:HG2	3:H:515:HOH:O	2.20	0.41
1:C:205:THR:O	1:C:205:THR:OG1	2.35	0.41
1:E:38:LYS:N	3:E:504:HOH:O	2.52	0.41
1:A:335:LEU:HD22	1:A:346:PHE:CD2	2.56	0.41
1:C:53:TRP:HB3	1:C:69:LEU:HD11	2.01	0.41
1:D:26:TYR:CG	1:D:88:VAL:HG22	2.55	0.41
1:E:62:ALA:HB2	1:E:357:ILE:HD13	2.03	0.41
1:G:86:LYS:NZ	1:G:86:LYS:CD	2.73	0.41
1:D:49:ASP:OD1	1:D:49:ASP:N	2.54	0.41
1:D:76:ARG:NH2	1:F:49:ASP:OD2	2.47	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:109:CYS:HA	1:D:141:CYS:HA	2.01	0.41
1:E:207:THR:CG2	1:E:306:VAL:HG21	2.51	0.41
1:G:108:ASP:OD2	1:G:108:ASP:N	2.49	0.41
1:H:73:GLN:HB2	1:H:76:ARG:HG3	2.02	0.41
1:G:324:GLN:OE1	1:G:327:SER:HA	2.22	0.40
1:D:183:ASN:O	1:D:184:ASN:HB2	2.22	0.40
1:F:98:VAL:HB	1:F:177:LYS:HG3	2.04	0.40
1:G:41:TRP:HB3	1:G:80:LEU:HD11	2.03	0.40
1:G:141:CYS:HB3	1:G:145:LYS:HE3	2.03	0.40
1:B:60:LEU:HD23	1:B:60:LEU:HA	1.89	0.40
1:C:212:GLY:HA2	1:C:305:PRO:HD3	1.80	0.40
1:G:354:LEU:HD12	1:G:354:LEU:HA	1.86	0.40
1:B:145:LYS:O	1:B:146:ARG:HG3	2.22	0.40
1:B:161:LYS:HZ3	1:B:203:ASP:HB3	1.86	0.40
1:C:104:GLU:OE1	3:C:506:HOH:O	2.22	0.40
1:G:135:MET:HB3	1:G:151:ILE:HB	2.02	0.40
1:G:351:LEU:O	1:G:355:ARG:HG2	2.21	0.40
1:H:82:ARG:HD3	1:H:83:LYS:O	2.21	0.40
1:C:159:TYR:CD1	1:C:206:MET:HG2	2.57	0.40
1:D:96:VAL:HG21	1:D:340:PRO:HA	2.04	0.40
1:E:28:LYS:HD2	1:E:346:PHE:CE2	2.55	0.40
1:E:63:ASP:C	1:E:65:ASN:N	2.78	0.40
1:H:118:LEU:HD23	1:H:118:LEU:HA	1.86	0.40
1:H:187:GLU:OE2	1:H:194:LYS:HE2	2.20	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:548:HOH:O	3:D:559:HOH:O[1_655]	1.97	0.23

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

The Analysed column shows the number of residues for which the backbone conformation was

analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	249/273 (91%)	235 (94%)	14 (6%)	0	100	100
1	B	249/273 (91%)	237 (95%)	12 (5%)	0	100	100
1	C	249/273 (91%)	234 (94%)	14 (6%)	1 (0%)	30	50
1	D	249/273 (91%)	235 (94%)	13 (5%)	1 (0%)	30	50
1	E	249/273 (91%)	231 (93%)	17 (7%)	1 (0%)	30	50
1	F	249/273 (91%)	232 (93%)	17 (7%)	0	100	100
1	G	249/273 (91%)	234 (94%)	14 (6%)	1 (0%)	30	50
1	H	249/273 (91%)	230 (92%)	17 (7%)	2 (1%)	16	30
All	All	1992/2184 (91%)	1868 (94%)	118 (6%)	6 (0%)	36	56

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	144	GLU
1	H	59	LYS
1	G	349	LYS
1	C	210	GLY
1	H	210	GLY
1	E	36	ALA

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 8 ligands modelled in this entry, 8 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	251/273 (91%)	-0.51	0 100 100	14, 25, 41, 67	0
1	B	251/273 (91%)	-0.34	4 (1%) 70 69	17, 27, 47, 70	0
1	C	251/273 (91%)	-0.22	7 (2%) 55 53	19, 28, 54, 97	0
1	D	251/273 (91%)	-0.24	8 (3%) 50 48	15, 27, 54, 105	0
1	E	251/273 (91%)	-0.07	3 (1%) 76 76	19, 35, 65, 79	0
1	F	251/273 (91%)	-0.09	9 (3%) 46 44	15, 29, 59, 96	0
1	G	251/273 (91%)	-0.15	1 (0%) 88 89	22, 35, 54, 69	0
1	H	251/273 (91%)	-0.01	9 (3%) 46 44	21, 35, 60, 110	0
All	All	2008/2184 (91%)	-0.20	41 (2%) 65 63	14, 30, 56, 110	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	209	PRO	4.2
1	F	211	SER	3.9
1	F	209	PRO	3.5
1	F	212	GLY	3.4
1	H	210	GLY	3.4
1	F	199	ARG	3.4
1	C	209	PRO	3.0
1	H	211	SER	3.0
1	D	211	SER	3.0
1	E	45	LYS	3.0
1	B	146	ARG	2.9
1	H	208	GLU	2.8
1	D	199	ARG	2.8
1	E	21	SER	2.7
1	B	145	LYS	2.7
1	H	212	GLY	2.7

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Mol	Chain	Res	Type	RSRZ
1	G	144	GLU	2.7
1	C	212	GLY	2.6
1	D	210	GLY	2.6
1	B	22	GLU	2.6
1	C	211	SER	2.5
1	C	210	GLY	2.5
1	F	207	THR	2.4
1	H	207	THR	2.4
1	D	208	GLU	2.3
1	H	209	PRO	2.3
1	C	22	GLU	2.3
1	B	48	ASP	2.3
1	D	104	GLU	2.2
1	H	305	PRO	2.2
1	D	144	GLU	2.2
1	H	162	ARG	2.2
1	C	143	PRO	2.2
1	F	48	ASP	2.2
1	F	210	GLY	2.2
1	C	208	GLU	2.1
1	E	144	GLU	2.1
1	F	208	GLU	2.1
1	H	143	PRO	2.1
1	F	45	LYS	2.0
1	D	207	THR	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 ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

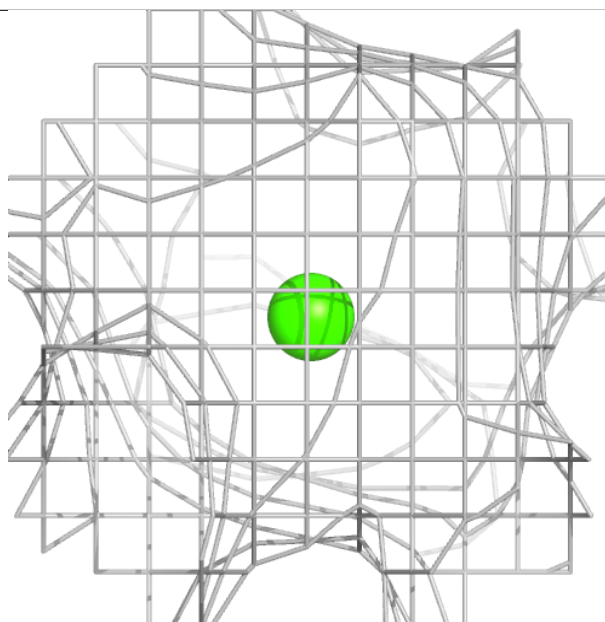
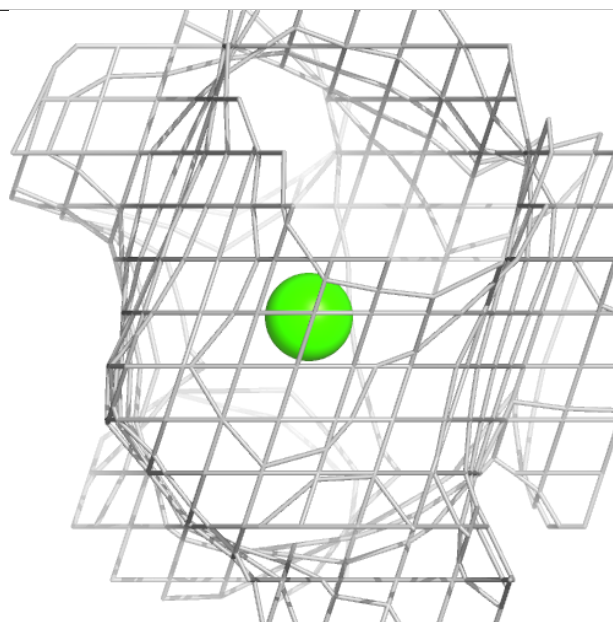
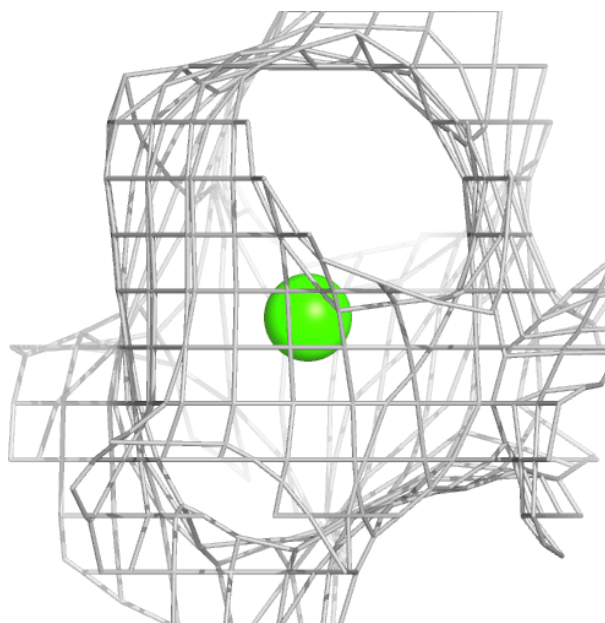
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	CA	A	401	1/1	0.96	0.04	31,31,31,31	0
2	CA	E	401	1/1	0.97	0.04	68,68,68,68	0
2	CA	F	401	1/1	0.97	0.04	47,47,47,47	0
2	CA	G	401	1/1	0.97	0.04	60,60,60,60	0
2	CA	D	401	1/1	0.98	0.03	38,38,38,38	0
2	CA	H	401	1/1	0.98	0.03	42,42,42,42	0
2	CA	C	401	1/1	0.99	0.02	37,37,37,37	0
2	CA	B	401	1/1	0.99	0.02	38,38,38,38	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.

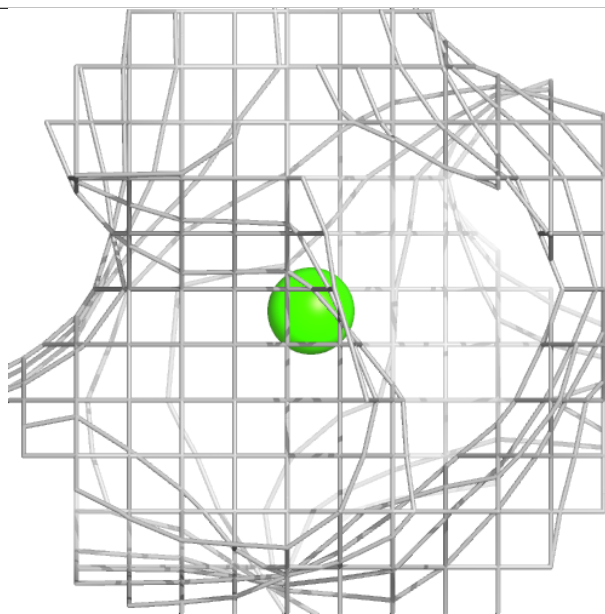
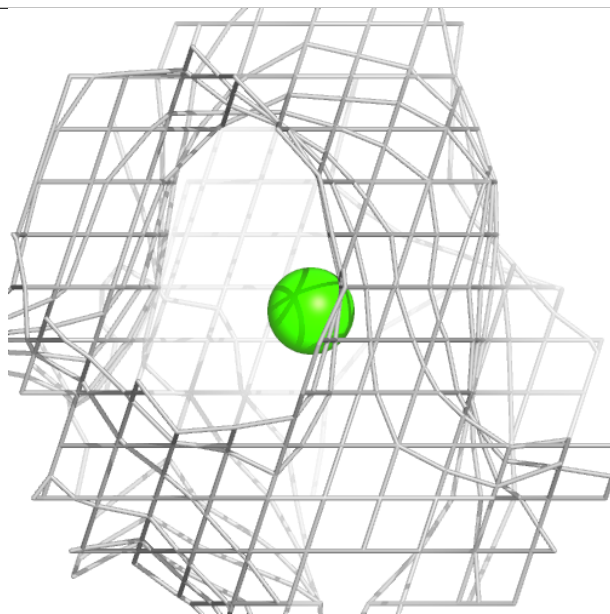
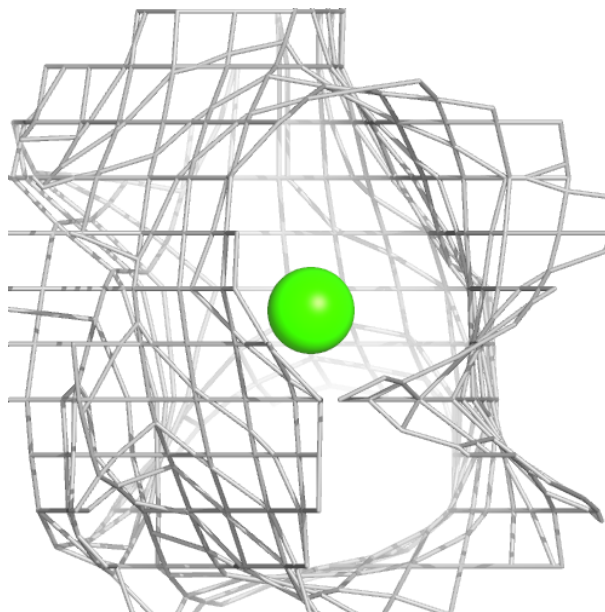
Electron density around CA A 401:

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



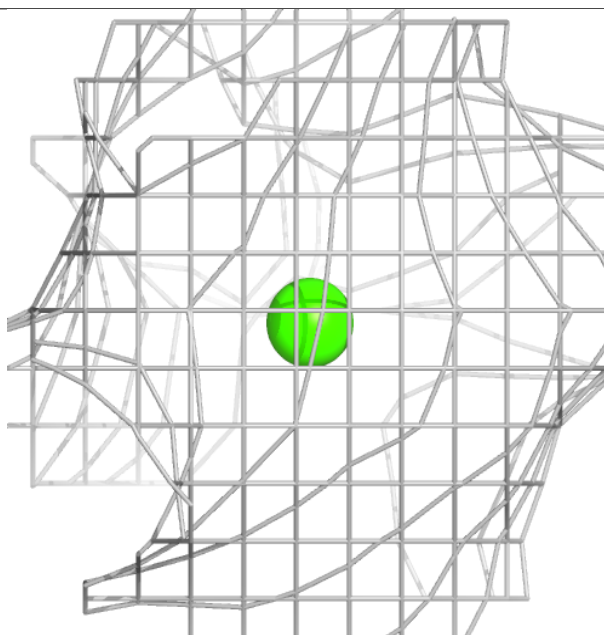
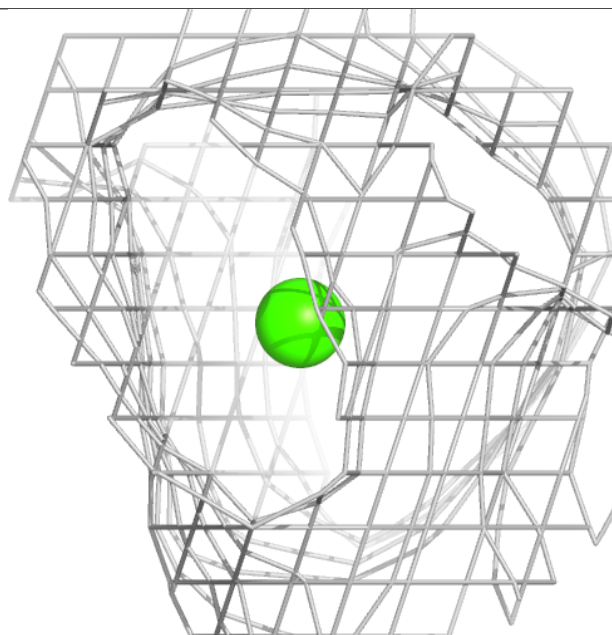
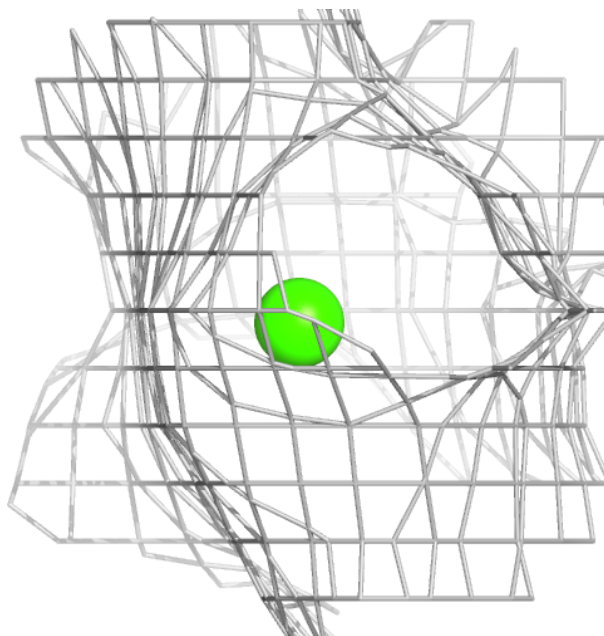
Electron density around CA E 401:

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



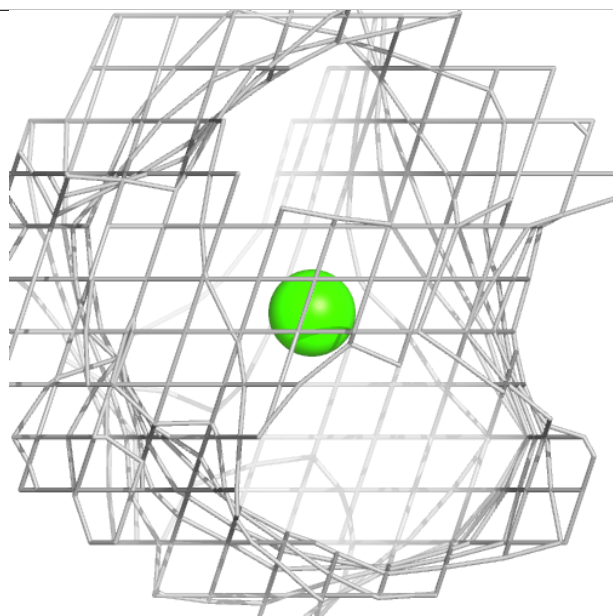
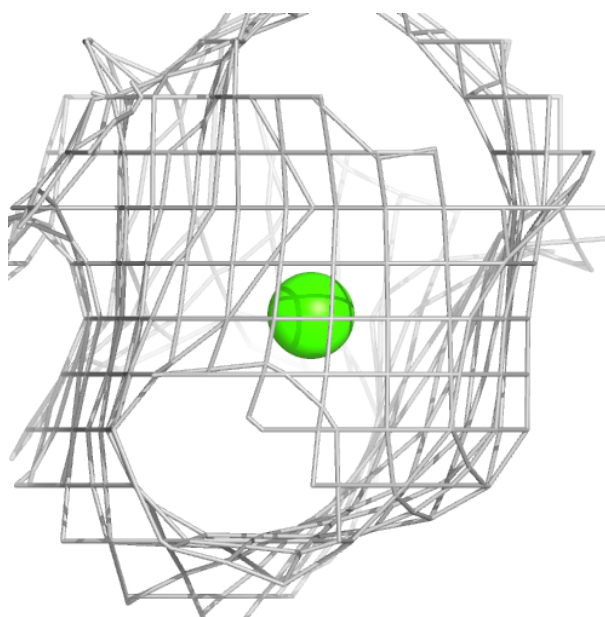
Electron density around CA F 401:

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



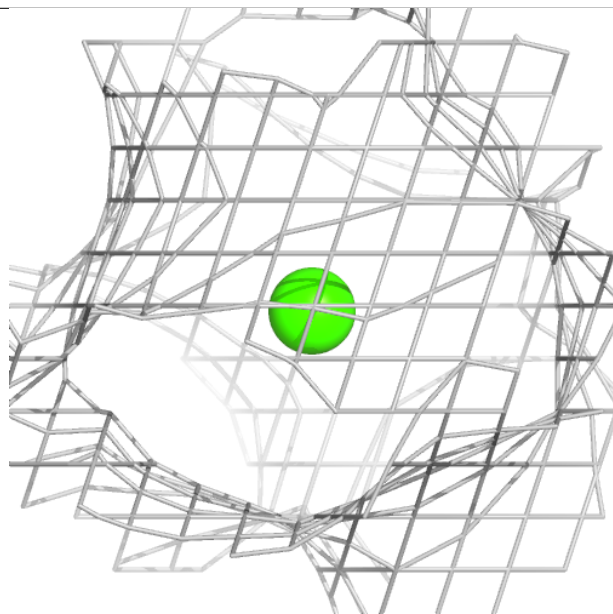
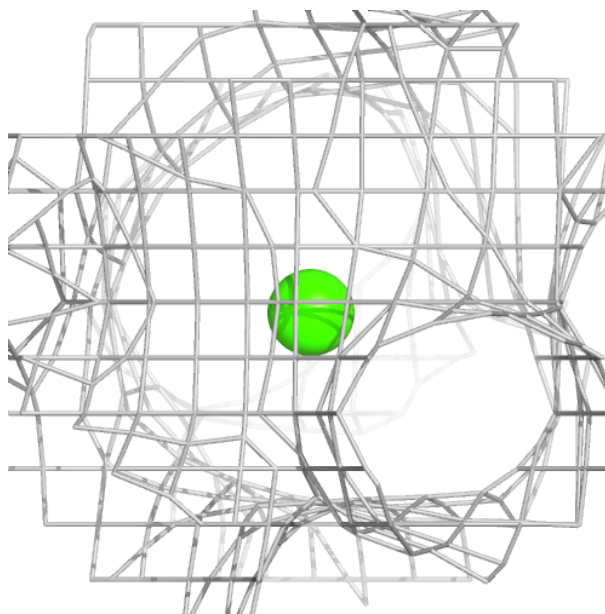
Electron density around CA G 401:

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



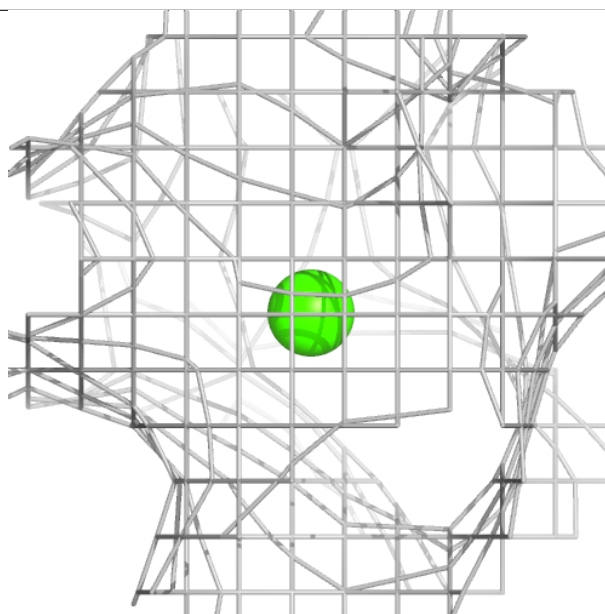
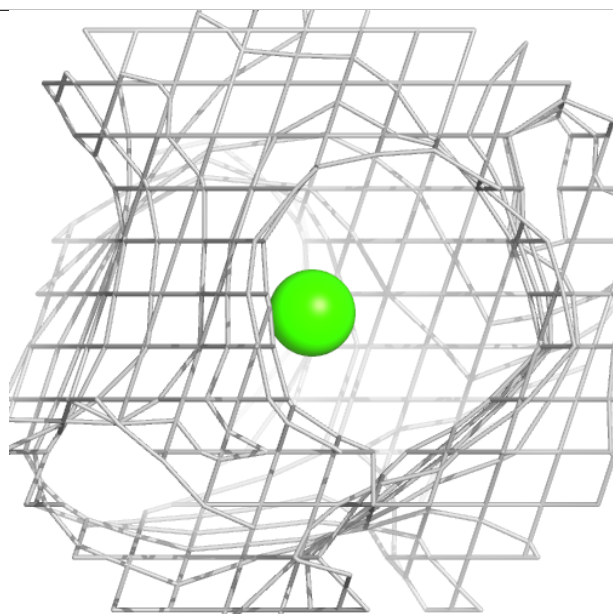
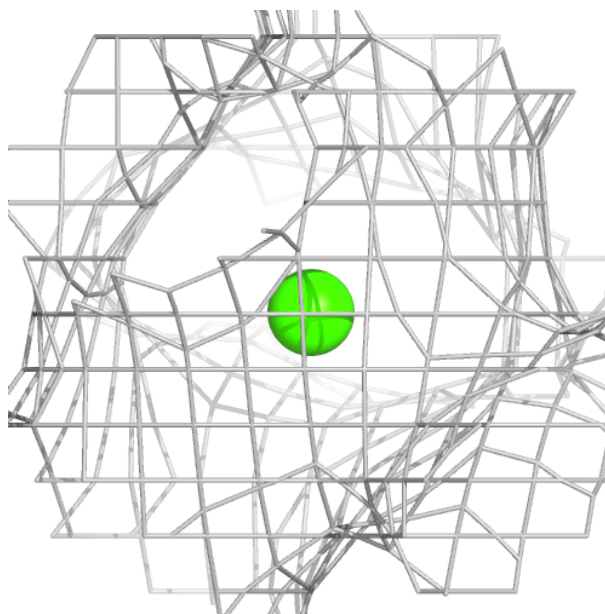
Electron density around CA D 401:

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



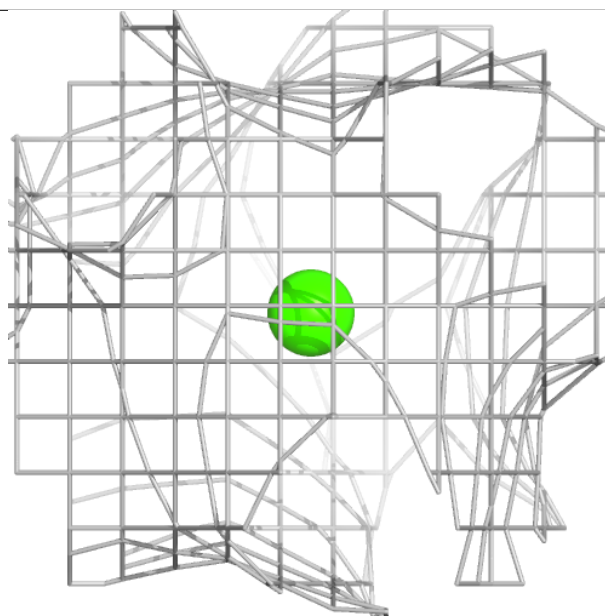
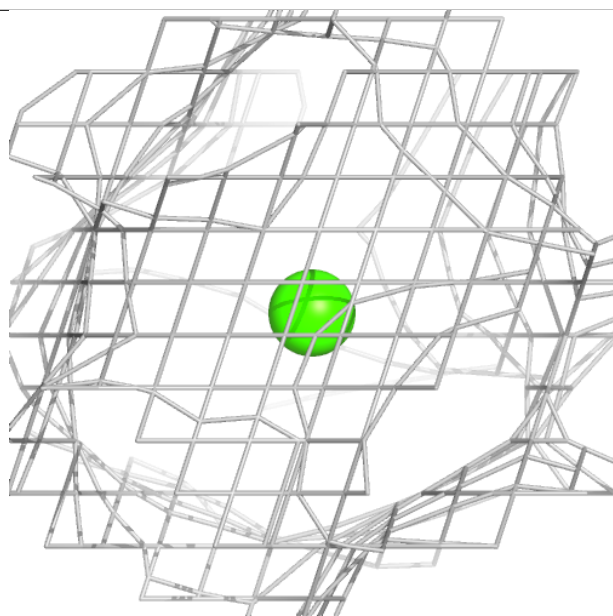
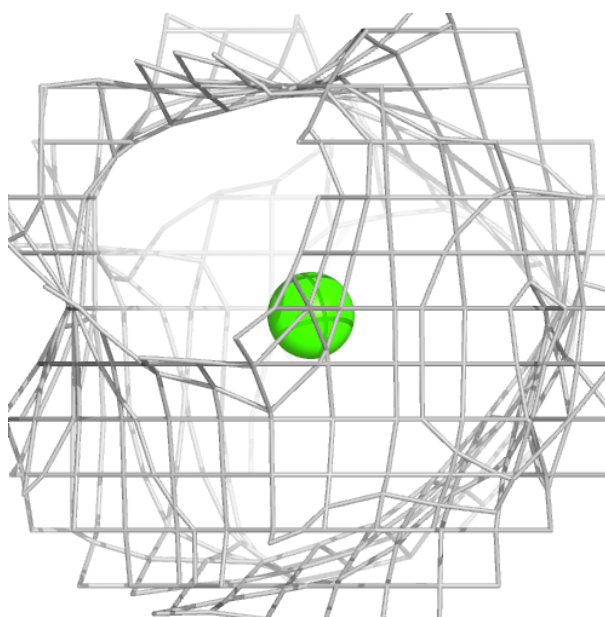
Electron density around CA H 401:

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



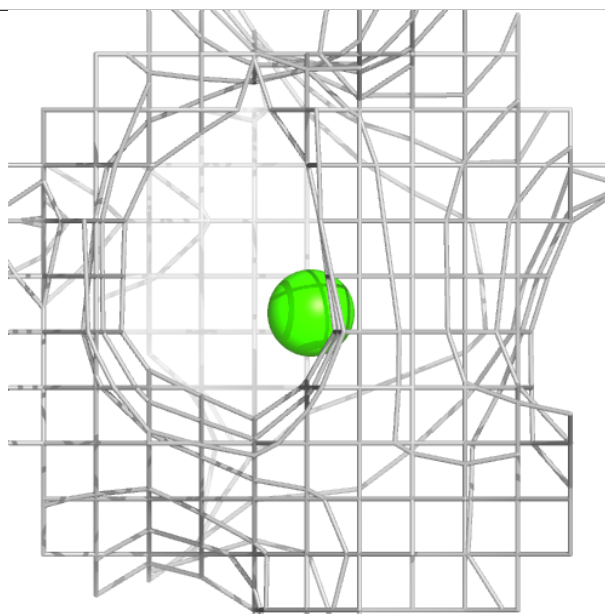
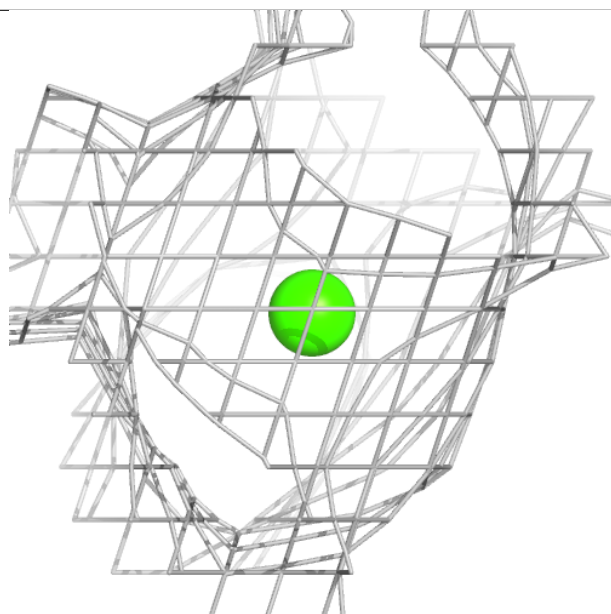
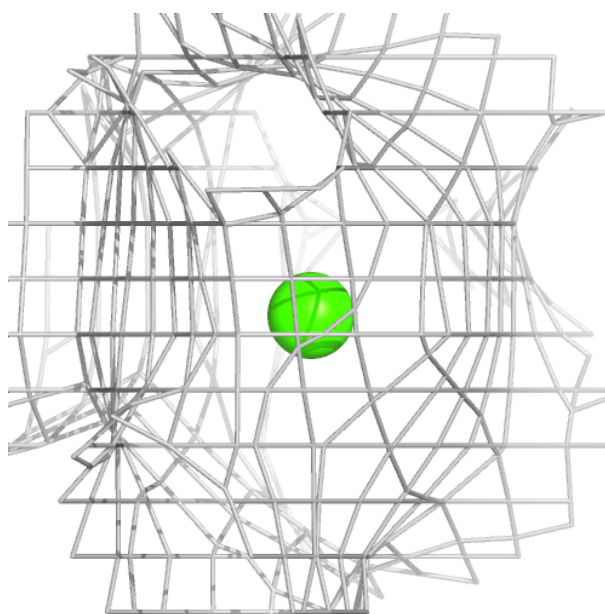
Electron density around CA C 401:

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



Electron density around CA B 401:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.