



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:47 PM UTC

PDB ID : 6WA6 / pdb_00006wa6
Title : Structure of the Chlamydia pneumoniae CdsV protein
Authors : Jensen, J.L.; Spiller, B.W.
Deposited on : 2020-03-24
Resolution : 2.80 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
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

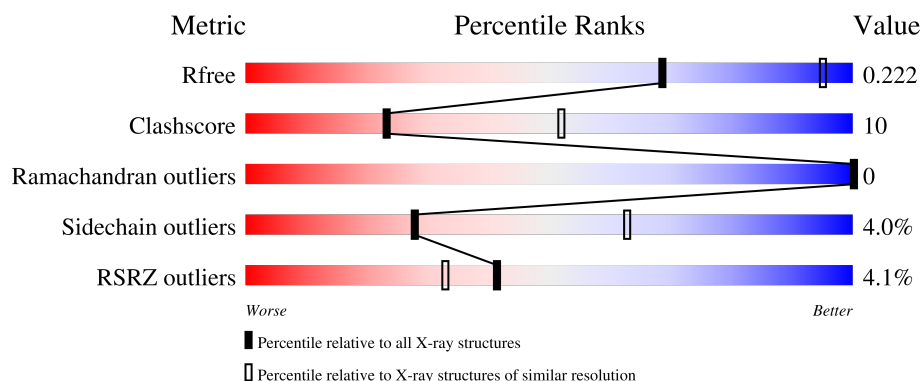
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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.



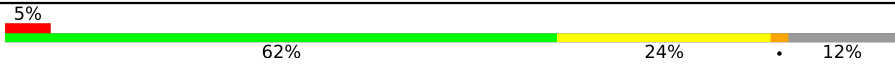
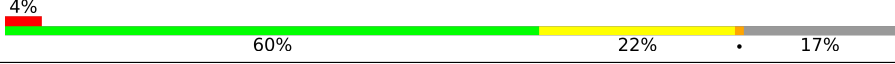
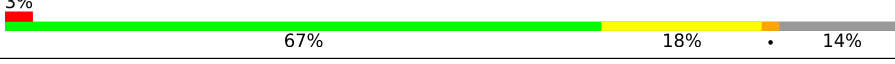
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	387	2% 72% 16% 10%
1	B	387	4% 67% 21% 10%
1	C	387	4% 67% 21% 12%
1	D	387	3% 72% 16% 11%
1	E	387	3% 66% 21% 11%

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Mol	Chain	Length	Quality of chain
1	F	387	
1	G	387	
1	H	387	
1	I	387	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 24992 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 Low calcium response locus protein D.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	348	Total	C	N	O	S	0	5	0
			2845	1843	470	525	7			
1	B	347	Total	C	N	O	S	0	0	0
			2800	1811	461	521	7			
1	C	342	Total	C	N	O	S	0	0	0
			2767	1790	457	513	7			
1	D	343	Total	C	N	O	S	0	2	0
			2793	1808	461	517	7			
1	E	343	Total	C	N	O	S	0	0	0
			2779	1798	457	517	7			
1	F	340	Total	C	N	O	S	0	1	0
			2755	1783	453	512	7			
1	G	343	Total	C	N	O	S	0	1	0
			2773	1794	455	517	7			
1	H	320	Total	C	N	O	S	0	0	0
			2609	1692	425	485	7			
1	I	334	Total	C	N	O	S	0	3	0
			2735	1770	452	506	7			

There are 189 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	324	MET	-	initiating methionine	UNP Q9Z8L5
A	325	GLY	-	expression tag	UNP Q9Z8L5
A	326	SER	-	expression tag	UNP Q9Z8L5
A	327	SER	-	expression tag	UNP Q9Z8L5
A	328	HIS	-	expression tag	UNP Q9Z8L5
A	329	HIS	-	expression tag	UNP Q9Z8L5
A	330	HIS	-	expression tag	UNP Q9Z8L5
A	331	HIS	-	expression tag	UNP Q9Z8L5
A	332	HIS	-	expression tag	UNP Q9Z8L5
A	333	HIS	-	expression tag	UNP Q9Z8L5
A	334	SER	-	expression tag	UNP Q9Z8L5

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Chain	Residue	Modelled	Actual	Comment	Reference
A	335	SER	-	expression tag	UNP Q9Z8L5
A	336	GLY	-	expression tag	UNP Q9Z8L5
A	337	LEU	-	expression tag	UNP Q9Z8L5
A	338	VAL	-	expression tag	UNP Q9Z8L5
A	339	PRO	-	expression tag	UNP Q9Z8L5
A	340	ARG	-	expression tag	UNP Q9Z8L5
A	341	GLY	-	expression tag	UNP Q9Z8L5
A	342	SER	-	expression tag	UNP Q9Z8L5
A	343	HIS	-	expression tag	UNP Q9Z8L5
A	344	MET	-	expression tag	UNP Q9Z8L5
B	324	MET	-	initiating methionine	UNP Q9Z8L5
B	325	GLY	-	expression tag	UNP Q9Z8L5
B	326	SER	-	expression tag	UNP Q9Z8L5
B	327	SER	-	expression tag	UNP Q9Z8L5
B	328	HIS	-	expression tag	UNP Q9Z8L5
B	329	HIS	-	expression tag	UNP Q9Z8L5
B	330	HIS	-	expression tag	UNP Q9Z8L5
B	331	HIS	-	expression tag	UNP Q9Z8L5
B	332	HIS	-	expression tag	UNP Q9Z8L5
B	333	HIS	-	expression tag	UNP Q9Z8L5
B	334	SER	-	expression tag	UNP Q9Z8L5
B	335	SER	-	expression tag	UNP Q9Z8L5
B	336	GLY	-	expression tag	UNP Q9Z8L5
B	337	LEU	-	expression tag	UNP Q9Z8L5
B	338	VAL	-	expression tag	UNP Q9Z8L5
B	339	PRO	-	expression tag	UNP Q9Z8L5
B	340	ARG	-	expression tag	UNP Q9Z8L5
B	341	GLY	-	expression tag	UNP Q9Z8L5
B	342	SER	-	expression tag	UNP Q9Z8L5
B	343	HIS	-	expression tag	UNP Q9Z8L5
B	344	MET	-	expression tag	UNP Q9Z8L5
C	324	MET	-	initiating methionine	UNP Q9Z8L5
C	325	GLY	-	expression tag	UNP Q9Z8L5
C	326	SER	-	expression tag	UNP Q9Z8L5
C	327	SER	-	expression tag	UNP Q9Z8L5
C	328	HIS	-	expression tag	UNP Q9Z8L5
C	329	HIS	-	expression tag	UNP Q9Z8L5
C	330	HIS	-	expression tag	UNP Q9Z8L5
C	331	HIS	-	expression tag	UNP Q9Z8L5
C	332	HIS	-	expression tag	UNP Q9Z8L5
C	333	HIS	-	expression tag	UNP Q9Z8L5
C	334	SER	-	expression tag	UNP Q9Z8L5

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Chain	Residue	Modelled	Actual	Comment	Reference
C	335	SER	-	expression tag	UNP Q9Z8L5
C	336	GLY	-	expression tag	UNP Q9Z8L5
C	337	LEU	-	expression tag	UNP Q9Z8L5
C	338	VAL	-	expression tag	UNP Q9Z8L5
C	339	PRO	-	expression tag	UNP Q9Z8L5
C	340	ARG	-	expression tag	UNP Q9Z8L5
C	341	GLY	-	expression tag	UNP Q9Z8L5
C	342	SER	-	expression tag	UNP Q9Z8L5
C	343	HIS	-	expression tag	UNP Q9Z8L5
C	344	MET	-	expression tag	UNP Q9Z8L5
D	324	MET	-	initiating methionine	UNP Q9Z8L5
D	325	GLY	-	expression tag	UNP Q9Z8L5
D	326	SER	-	expression tag	UNP Q9Z8L5
D	327	SER	-	expression tag	UNP Q9Z8L5
D	328	HIS	-	expression tag	UNP Q9Z8L5
D	329	HIS	-	expression tag	UNP Q9Z8L5
D	330	HIS	-	expression tag	UNP Q9Z8L5
D	331	HIS	-	expression tag	UNP Q9Z8L5
D	332	HIS	-	expression tag	UNP Q9Z8L5
D	333	HIS	-	expression tag	UNP Q9Z8L5
D	334	SER	-	expression tag	UNP Q9Z8L5
D	335	SER	-	expression tag	UNP Q9Z8L5
D	336	GLY	-	expression tag	UNP Q9Z8L5
D	337	LEU	-	expression tag	UNP Q9Z8L5
D	338	VAL	-	expression tag	UNP Q9Z8L5
D	339	PRO	-	expression tag	UNP Q9Z8L5
D	340	ARG	-	expression tag	UNP Q9Z8L5
D	341	GLY	-	expression tag	UNP Q9Z8L5
D	342	SER	-	expression tag	UNP Q9Z8L5
D	343	HIS	-	expression tag	UNP Q9Z8L5
D	344	MET	-	expression tag	UNP Q9Z8L5
E	324	MET	-	initiating methionine	UNP Q9Z8L5
E	325	GLY	-	expression tag	UNP Q9Z8L5
E	326	SER	-	expression tag	UNP Q9Z8L5
E	327	SER	-	expression tag	UNP Q9Z8L5
E	328	HIS	-	expression tag	UNP Q9Z8L5
E	329	HIS	-	expression tag	UNP Q9Z8L5
E	330	HIS	-	expression tag	UNP Q9Z8L5
E	331	HIS	-	expression tag	UNP Q9Z8L5
E	332	HIS	-	expression tag	UNP Q9Z8L5
E	333	HIS	-	expression tag	UNP Q9Z8L5
E	334	SER	-	expression tag	UNP Q9Z8L5

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Chain	Residue	Modelled	Actual	Comment	Reference
E	335	SER	-	expression tag	UNP Q9Z8L5
E	336	GLY	-	expression tag	UNP Q9Z8L5
E	337	LEU	-	expression tag	UNP Q9Z8L5
E	338	VAL	-	expression tag	UNP Q9Z8L5
E	339	PRO	-	expression tag	UNP Q9Z8L5
E	340	ARG	-	expression tag	UNP Q9Z8L5
E	341	GLY	-	expression tag	UNP Q9Z8L5
E	342	SER	-	expression tag	UNP Q9Z8L5
E	343	HIS	-	expression tag	UNP Q9Z8L5
E	344	MET	-	expression tag	UNP Q9Z8L5
F	324	MET	-	initiating methionine	UNP Q9Z8L5
F	325	GLY	-	expression tag	UNP Q9Z8L5
F	326	SER	-	expression tag	UNP Q9Z8L5
F	327	SER	-	expression tag	UNP Q9Z8L5
F	328	HIS	-	expression tag	UNP Q9Z8L5
F	329	HIS	-	expression tag	UNP Q9Z8L5
F	330	HIS	-	expression tag	UNP Q9Z8L5
F	331	HIS	-	expression tag	UNP Q9Z8L5
F	332	HIS	-	expression tag	UNP Q9Z8L5
F	333	HIS	-	expression tag	UNP Q9Z8L5
F	334	SER	-	expression tag	UNP Q9Z8L5
F	335	SER	-	expression tag	UNP Q9Z8L5
F	336	GLY	-	expression tag	UNP Q9Z8L5
F	337	LEU	-	expression tag	UNP Q9Z8L5
F	338	VAL	-	expression tag	UNP Q9Z8L5
F	339	PRO	-	expression tag	UNP Q9Z8L5
F	340	ARG	-	expression tag	UNP Q9Z8L5
F	341	GLY	-	expression tag	UNP Q9Z8L5
F	342	SER	-	expression tag	UNP Q9Z8L5
F	343	HIS	-	expression tag	UNP Q9Z8L5
F	344	MET	-	expression tag	UNP Q9Z8L5
G	324	MET	-	initiating methionine	UNP Q9Z8L5
G	325	GLY	-	expression tag	UNP Q9Z8L5
G	326	SER	-	expression tag	UNP Q9Z8L5
G	327	SER	-	expression tag	UNP Q9Z8L5
G	328	HIS	-	expression tag	UNP Q9Z8L5
G	329	HIS	-	expression tag	UNP Q9Z8L5
G	330	HIS	-	expression tag	UNP Q9Z8L5
G	331	HIS	-	expression tag	UNP Q9Z8L5
G	332	HIS	-	expression tag	UNP Q9Z8L5
G	333	HIS	-	expression tag	UNP Q9Z8L5
G	334	SER	-	expression tag	UNP Q9Z8L5

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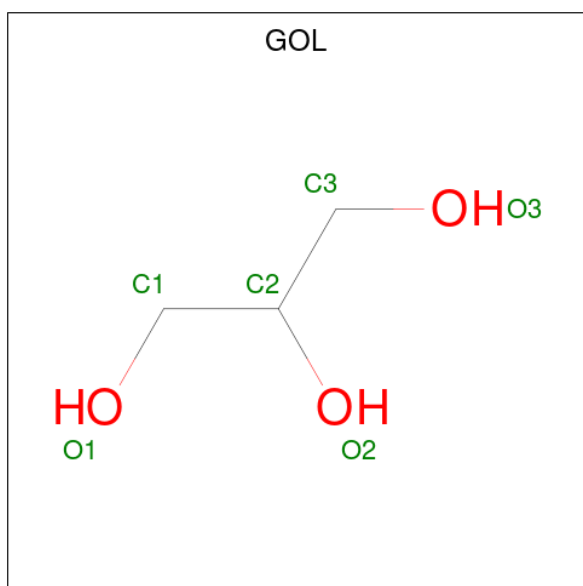
Chain	Residue	Modelled	Actual	Comment	Reference
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G	336	GLY	-	expression tag	UNP Q9Z8L5
G	337	LEU	-	expression tag	UNP Q9Z8L5
G	338	VAL	-	expression tag	UNP Q9Z8L5
G	339	PRO	-	expression tag	UNP Q9Z8L5
G	340	ARG	-	expression tag	UNP Q9Z8L5
G	341	GLY	-	expression tag	UNP Q9Z8L5
G	342	SER	-	expression tag	UNP Q9Z8L5
G	343	HIS	-	expression tag	UNP Q9Z8L5
G	344	MET	-	expression tag	UNP Q9Z8L5
H	324	MET	-	initiating methionine	UNP Q9Z8L5
H	325	GLY	-	expression tag	UNP Q9Z8L5
H	326	SER	-	expression tag	UNP Q9Z8L5
H	327	SER	-	expression tag	UNP Q9Z8L5
H	328	HIS	-	expression tag	UNP Q9Z8L5
H	329	HIS	-	expression tag	UNP Q9Z8L5
H	330	HIS	-	expression tag	UNP Q9Z8L5
H	331	HIS	-	expression tag	UNP Q9Z8L5
H	332	HIS	-	expression tag	UNP Q9Z8L5
H	333	HIS	-	expression tag	UNP Q9Z8L5
H	334	SER	-	expression tag	UNP Q9Z8L5
H	335	SER	-	expression tag	UNP Q9Z8L5
H	336	GLY	-	expression tag	UNP Q9Z8L5
H	337	LEU	-	expression tag	UNP Q9Z8L5
H	338	VAL	-	expression tag	UNP Q9Z8L5
H	339	PRO	-	expression tag	UNP Q9Z8L5
H	340	ARG	-	expression tag	UNP Q9Z8L5
H	341	GLY	-	expression tag	UNP Q9Z8L5
H	342	SER	-	expression tag	UNP Q9Z8L5
H	343	HIS	-	expression tag	UNP Q9Z8L5
H	344	MET	-	expression tag	UNP Q9Z8L5
I	324	MET	-	initiating methionine	UNP Q9Z8L5
I	325	GLY	-	expression tag	UNP Q9Z8L5
I	326	SER	-	expression tag	UNP Q9Z8L5
I	327	SER	-	expression tag	UNP Q9Z8L5
I	328	HIS	-	expression tag	UNP Q9Z8L5
I	329	HIS	-	expression tag	UNP Q9Z8L5
I	330	HIS	-	expression tag	UNP Q9Z8L5
I	331	HIS	-	expression tag	UNP Q9Z8L5
I	332	HIS	-	expression tag	UNP Q9Z8L5
I	333	HIS	-	expression tag	UNP Q9Z8L5
I	334	SER	-	expression tag	UNP Q9Z8L5

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Chain	Residue	Modelled	Actual	Comment	Reference
I	335	SER	-	expression tag	UNP Q9Z8L5
I	336	GLY	-	expression tag	UNP Q9Z8L5
I	337	LEU	-	expression tag	UNP Q9Z8L5
I	338	VAL	-	expression tag	UNP Q9Z8L5
I	339	PRO	-	expression tag	UNP Q9Z8L5
I	340	ARG	-	expression tag	UNP Q9Z8L5
I	341	GLY	-	expression tag	UNP Q9Z8L5
I	342	SER	-	expression tag	UNP Q9Z8L5
I	343	HIS	-	expression tag	UNP Q9Z8L5
I	344	MET	-	expression tag	UNP Q9Z8L5

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	C	1	Total C O 6 3 3	0	0
2	D	1	Total C O 6 3 3	0	0
2	E	1	Total C O 6 3 3	0	0
2	E	1	Total C O 6 3 3	0	0
2	E	1	Total C O 6 3 3	0	0
2	G	1	Total C O 6 3 3	0	0
2	I	1	Total C O 6 3 3	0	0

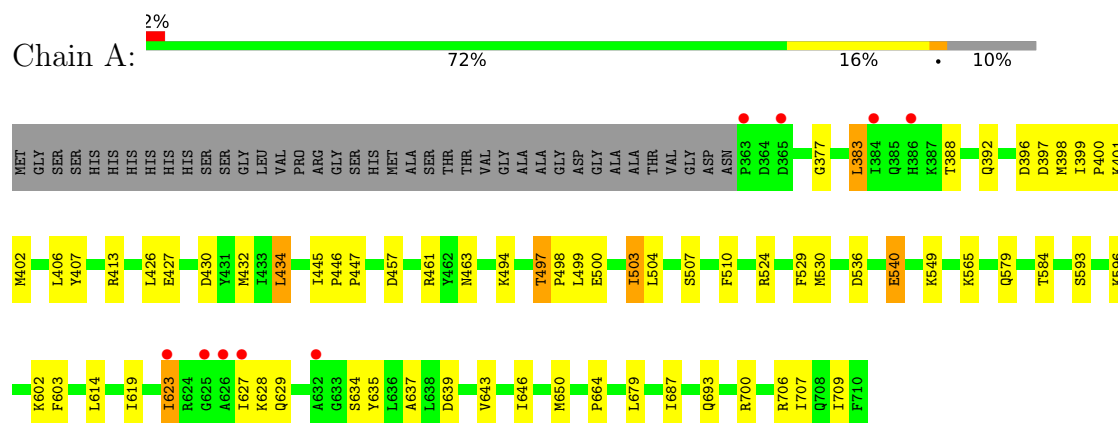
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	17	Total O 17 17	0	0
3	B	6	Total O 6 6	0	0
3	C	5	Total O 5 5	0	0
3	D	9	Total O 9 9	0	0
3	E	3	Total O 3 3	0	0
3	F	4	Total O 4 4	0	0
3	G	4	Total O 4 4	0	0
3	H	4	Total O 4 4	0	0
3	I	6	Total O 6 6	0	0

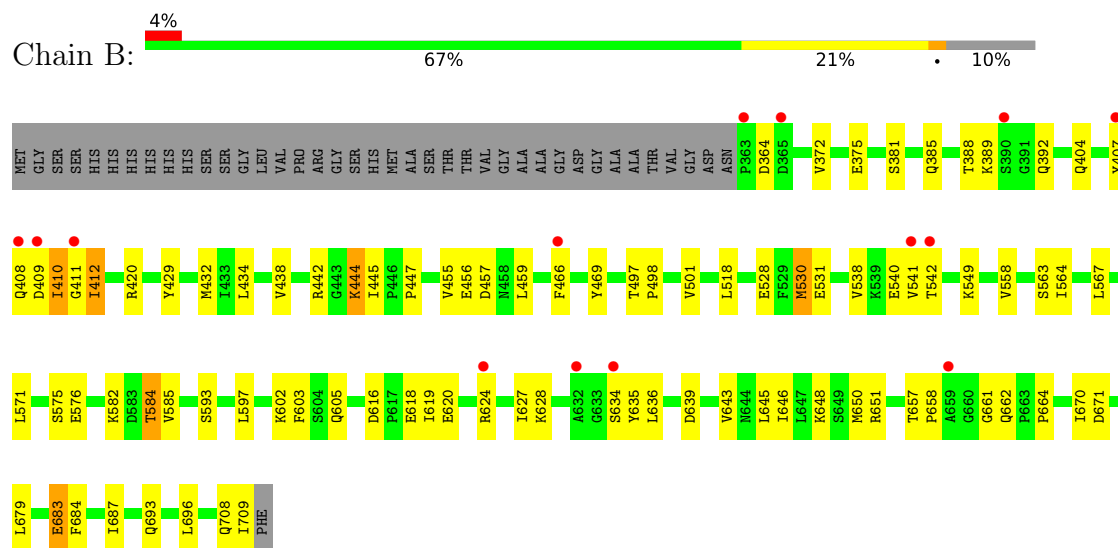
3 Residue-property plots

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.

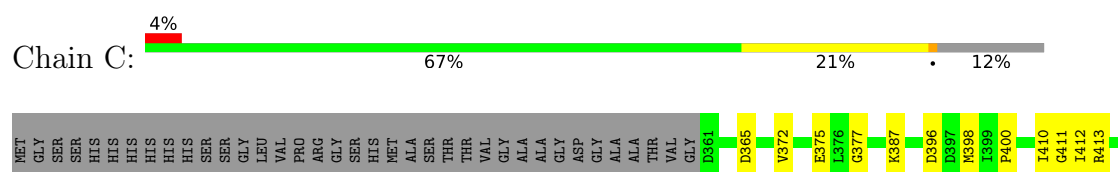
• Molecule 1: Low calcium response locus protein D

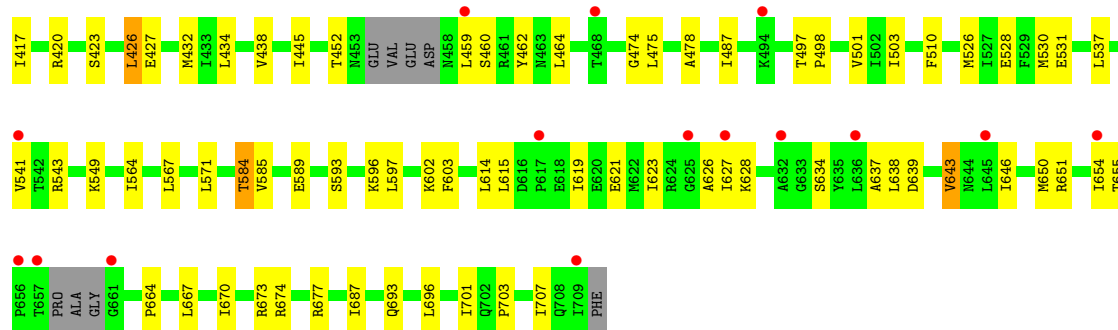


• Molecule 1: Low calcium response locus protein D

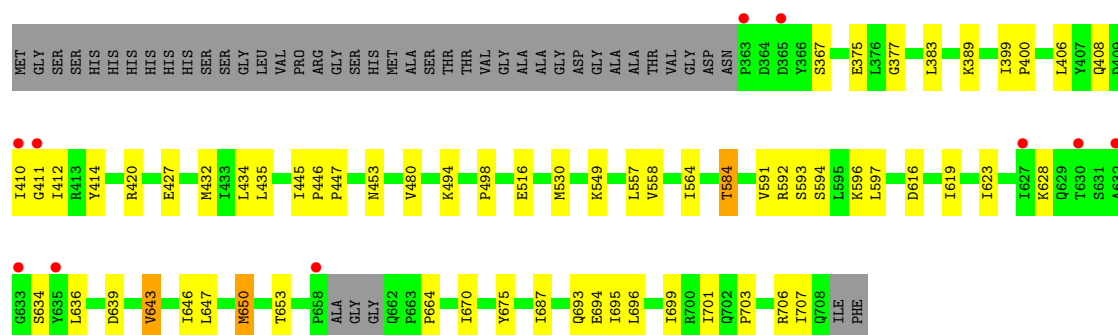


• Molecule 1: Low calcium response locus protein D

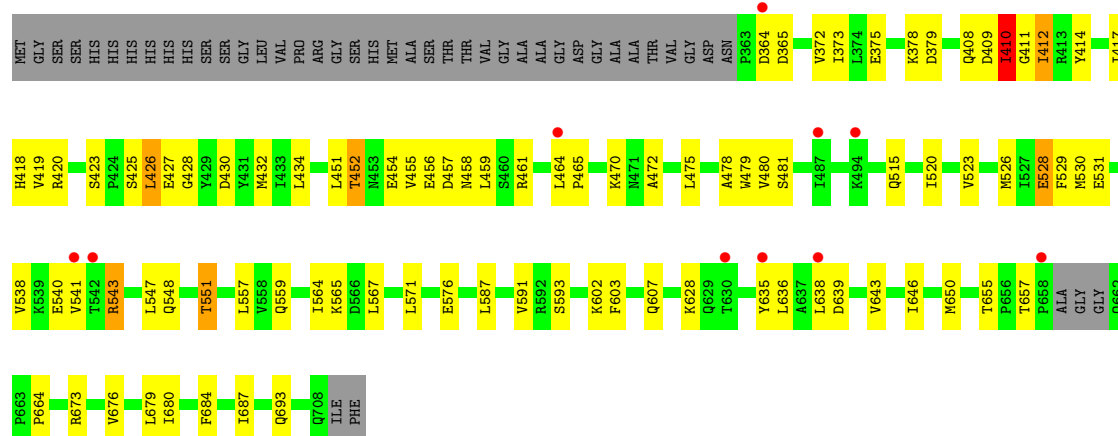




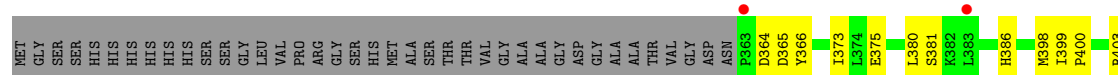
• Molecule 1: Low calcium response locus protein D

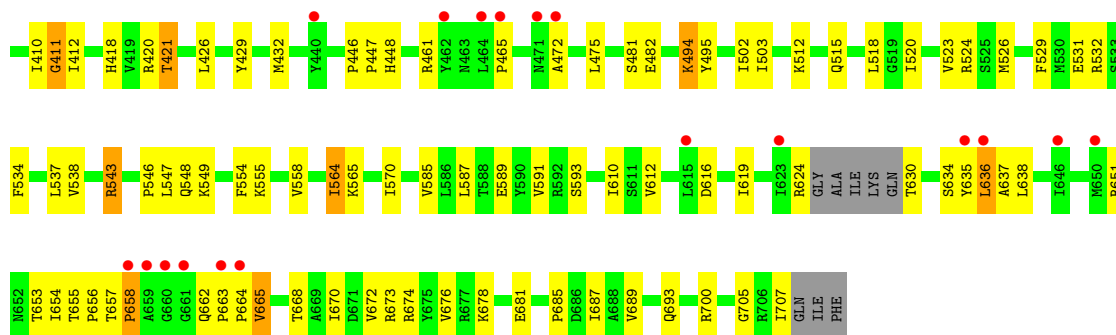


• Molecule 1: Low calcium response locus protein D



• Molecule 1: Low calcium response locus protein D





• Molecule 1: Low calcium response locus protein D

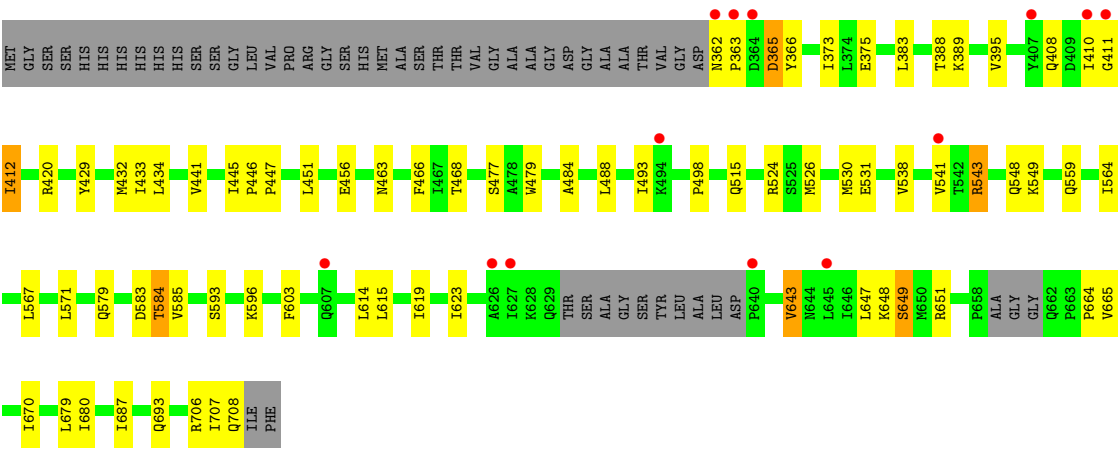


• Molecule 1: Low calcium response locus protein D



• Molecule 1: Low calcium response locus protein D





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	66.65Å 280.44Å 290.41Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	126.27 – 2.80 126.27 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.4 (126.27-2.80) 99.4 (126.27-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.31 (at 2.82Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.196 , 0.244 (Not available) , 0.222	Depositor DCC
R_{free} test set	6708 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	66.1	Xtriage
Anisotropy	0.729	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 62.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.011 for -h,l,k	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	24992	wwPDB-VP
Average B, all atoms (Å ²)	95.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.99% 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: 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	0.53	2/2923 (0.1%)	0.72	0/3965
1	B	0.45	0/2860	0.70	0/3881
1	C	0.39	0/2824	0.65	0/3829
1	D	0.39	0/2860	0.63	0/3880
1	E	0.48	3/2838 (0.1%)	0.68	1/3850 (0.0%)
1	F	0.47	2/2818 (0.1%)	0.66	4/3824 (0.1%)
1	G	0.43	0/2834	0.65	0/3845
1	H	0.41	0/2665	0.65	0/3614
1	I	0.44	1/2802 (0.0%)	0.66	1/3799 (0.0%)
All	All	0.45	8/25424 (0.0%)	0.67	6/34487 (0.0%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	410	ILE	C-O	-7.44	1.15	1.23
1	E	410	ILE	C-O	-7.05	1.14	1.23
1	I	412	ILE	C-O	-6.69	1.17	1.24
1	F	412	ILE	C-O	-5.68	1.17	1.24
1	E	412	ILE	C-O	-5.64	1.17	1.24
1	E	409	ASP	C-O	-5.46	1.17	1.24
1	A	504	LEU	C-O	-5.10	1.17	1.24
1	A	434	LEU	C-O	-5.02	1.18	1.24

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	409	ASP	N-CA-C	6.31	119.14	111.82
1	F	658	PRO	CB-CA-C	-6.20	103.82	111.64
1	F	412	ILE	N-CA-C	6.05	117.38	109.58
1	F	411	GLY	N-CA-C	5.83	121.02	113.27

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	705	GLY	CA-C-O	-5.82	118.26	122.45
1	I	365	ASP	CB-CA-C	-5.38	101.15	112.82

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	2845	0	2930	42	0
1	B	2800	0	2880	57	0
1	C	2767	0	2847	50	0
1	D	2793	0	2871	37	0
1	E	2779	0	2857	81	0
1	F	2755	0	2827	84	0
1	G	2773	0	2847	67	0
1	H	2609	0	2677	54	0
1	I	2735	0	2814	59	0
2	A	24	0	32	1	0
2	B	12	0	16	1	0
2	C	6	0	8	0	0
2	D	6	0	8	0	0
2	E	18	0	24	1	0
2	G	6	0	8	0	0
2	I	6	0	8	0	0
3	A	17	0	0	0	0
3	B	6	0	0	0	0
3	C	5	0	0	0	0
3	D	9	0	0	0	0
3	E	3	0	0	0	0
3	F	4	0	0	0	0
3	G	4	0	0	0	0
3	H	4	0	0	0	0
3	I	6	0	0	1	0
All	All	24992	0	25654	520	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 10.

All (520) 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:377:GLY:HA2	1:G:426:LEU:CD2	1.47	1.44
1:G:377:GLY:CA	1:G:426:LEU:HD22	1.52	1.38
1:E:464:LEU:CD1	1:E:465:PRO:HD3	1.54	1.33
1:G:412:ILE:HD11	1:G:558:VAL:CG2	1.69	1.20
1:E:464:LEU:HD12	1:E:465:PRO:CD	1.72	1.19
1:E:464:LEU:HD12	1:E:465:PRO:N	1.61	1.15
1:E:464:LEU:CG	1:E:465:PRO:HD3	1.80	1.12
1:G:412:ILE:HD12	1:G:514:SER:HB2	1.19	1.12
1:E:464:LEU:CD1	1:E:465:PRO:CD	2.26	1.11
1:F:546:PRO:HD2	1:F:549:LYS:HD2	1.20	1.11
1:G:412:ILE:CD1	1:G:558:VAL:HG21	1.82	1.09
1:F:612:VAL:HG12	1:F:665:VAL:HG11	1.15	1.08
1:F:653:THR:HG23	1:F:654:ILE:HD12	1.27	1.07
1:F:411:GLY:HA3	1:F:564:ILE:HG12	1.33	1.07
1:E:411:GLY:HA3	1:E:564:ILE:CG1	1.85	1.04
1:A:388:THR:OG1	1:A:392:GLN:HG2	1.61	1.00
1:C:628:LYS:HG3	1:C:634:SER:OG	1.62	0.99
1:F:612:VAL:CG1	1:F:665:VAL:HG11	1.93	0.99
1:E:464:LEU:HG	1:E:465:PRO:HD3	1.42	0.98
1:F:664:PRO:HG2	1:F:687:ILE:HA	1.44	0.97
1:C:541:VAL:HG11	1:C:571:LEU:HD22	1.49	0.94
1:B:624:ARG:HD3	1:B:671:ASP:OD2	1.67	0.94
1:E:411:GLY:HA3	1:E:564:ILE:HG13	1.47	0.94
1:F:653:THR:CG2	1:F:654:ILE:HD12	1.97	0.94
1:I:649:SER:HB3	1:I:707:ILE:HB	1.49	0.93
1:I:614:LEU:CD1	1:I:706:ARG:HH21	1.81	0.92
1:F:546:PRO:CD	1:F:549:LYS:HD2	2.00	0.92
1:E:411:GLY:HA3	1:E:564:ILE:HG12	1.52	0.90
1:D:549:LYS:HE2	1:D:584:THR:HG22	1.54	0.89
1:F:399:ILE:O	1:F:403:ARG:HG3	1.74	0.88
1:I:362:ASN:N	1:I:363:PRO:CD	2.36	0.88
1:H:647:LEU:HB3	1:H:651:ARG:HH12	1.38	0.88
1:G:426:LEU:CD1	1:G:428:GLY:H	1.87	0.88
1:G:412:ILE:HD11	1:G:558:VAL:HG21	0.89	0.87
1:I:614:LEU:HD13	1:I:706:ARG:HH21	1.37	0.87
1:F:670:ILE:HD12	1:F:693:GLN:HB2	1.59	0.84
1:E:464:LEU:HG	1:E:465:PRO:CD	2.08	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:411:GLY:HA3	1:F:564:ILE:CG1	2.08	0.84
1:B:531:GLU:HG3	1:B:538:VAL:HG11	1.60	0.83
1:F:612:VAL:HG12	1:F:665:VAL:CG1	2.04	0.83
1:G:412:ILE:CD1	1:G:514:SER:HB2	2.07	0.83
1:H:432:MET:HE3	1:H:434:LEU:HD11	1.61	0.82
1:I:541:VAL:HG11	1:I:571:LEU:HD22	1.60	0.82
1:E:464:LEU:CG	1:E:465:PRO:CD	2.57	0.82
1:E:464:LEU:HD11	1:E:465:PRO:HD3	1.60	0.82
1:I:614:LEU:HD13	1:I:706:ARG:NH2	1.95	0.81
1:F:657:THR:HG21	1:F:662:GLN:O	1.81	0.81
1:H:647:LEU:HB3	1:H:651:ARG:NH1	1.95	0.81
1:A:497:THR:HG22	1:A:500:GLU:H	1.46	0.80
1:I:614:LEU:HB2	1:I:706:ARG:HG2	1.64	0.80
1:I:410:ILE:HG21	1:I:412:ILE:HD12	1.64	0.80
1:B:645:LEU:HB3	1:B:709:ILE:HD13	1.64	0.78
1:B:388:THR:HG22	1:B:389:LYS:N	1.99	0.78
1:B:541:VAL:HG11	1:B:571:LEU:HD22	1.64	0.78
1:B:412:ILE:HD11	1:B:518:LEU:HD22	1.65	0.77
1:G:378:LYS:HG3	1:G:426:LEU:HD11	1.66	0.77
1:D:636:LEU:HD22	1:D:675:TYR:HB3	1.68	0.76
1:G:430:ASP:OD1	1:G:444:LYS:HG2	1.84	0.76
1:D:670:ILE:HD12	1:D:693:GLN:HB2	1.67	0.76
1:H:375:GLU:HG2	1:H:420:ARG:HB2	1.68	0.76
1:F:664:PRO:CG	1:F:687:ILE:HG12	2.15	0.75
1:I:411:GLY:HA3	1:I:564:ILE:HG13	1.66	0.75
1:B:549:LYS:HE2	1:B:584:THR:HG22	1.68	0.75
1:F:653:THR:HG23	1:F:654:ILE:CD1	2.13	0.75
1:G:411:GLY:HA3	1:G:564:ILE:HG13	1.68	0.75
1:F:612:VAL:HB	1:F:665:VAL:CG1	2.17	0.74
1:I:411:GLY:HA3	1:I:564:ILE:CG1	2.17	0.74
1:G:377:GLY:HA2	1:G:426:LEU:CG	2.17	0.74
1:I:362:ASN:N	1:I:363:PRO:HD2	2.02	0.74
1:H:523:VAL:HA	1:H:526:MET:HE2	1.70	0.74
1:F:657:THR:HB	1:F:662:GLN:HB2	1.70	0.74
1:E:372:VAL:HG21	1:E:417:ILE:HD13	1.70	0.73
1:I:549:LYS:HE3	1:I:584:THR:HG22	1.69	0.73
1:B:410:ILE:HG13	1:B:563:SER:HA	1.68	0.73
1:G:426:LEU:HD12	1:G:428:GLY:H	1.52	0.73
1:C:655:THR:O	1:C:655:THR:HG22	1.86	0.73
1:I:643:VAL:HG13	1:I:679:LEU:HD11	1.70	0.73
1:C:627:ILE:HG13	1:C:637:ALA:HB2	1.69	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:547:LEU:O	1:E:551:THR:HG23	1.90	0.72
1:I:543:ARG:HG2	1:I:543:ARG:O	1.90	0.71
1:C:664:PRO:HG2	1:C:687:ILE:HG12	1.72	0.71
1:E:452:THR:HG22	1:E:478:ALA:HB3	1.71	0.71
1:G:375:GLU:HG2	1:G:420:ARG:HB2	1.72	0.71
1:E:638:LEU:HB3	1:E:643:VAL:HG21	1.73	0.70
1:E:543:ARG:O	1:E:543:ARG:HG2	1.90	0.70
1:B:593:SER:HB2	1:B:693:GLN:HG2	1.73	0.70
1:F:664:PRO:HG3	1:F:687:ILE:HG12	1.71	0.70
1:A:398:MET:HE2	1:A:503:ILE:HD13	1.74	0.70
1:E:636:LEU:CD2	1:E:638:LEU:HG	2.22	0.69
1:G:412:ILE:HD12	1:G:514:SER:CB	2.12	0.69
1:B:432:MET:HE3	1:B:434:LEU:HD11	1.74	0.69
1:B:670:ILE:HD12	1:B:693:GLN:HB2	1.75	0.69
1:B:531:GLU:CG	1:B:538:VAL:HG11	2.22	0.69
1:F:494:LYS:HE3	1:F:495:TYR:H	1.56	0.69
1:F:612:VAL:CG1	1:F:665:VAL:CG1	2.65	0.68
1:F:612:VAL:CB	1:F:665:VAL:CG1	2.70	0.68
1:G:426:LEU:HD13	1:G:428:GLY:H	1.58	0.68
1:D:411:GLY:HA3	1:D:564:ILE:HG13	1.76	0.68
1:E:636:LEU:HD21	1:E:679:LEU:HD12	1.76	0.68
1:A:619:ILE:HD11	1:A:709:ILE:HG13	1.75	0.67
1:E:451:LEU:HD13	1:E:479:TRP:CE2	2.29	0.67
1:I:614:LEU:CD1	1:I:706:ARG:HE	2.08	0.67
1:D:664:PRO:HG2	1:D:687:ILE:HG12	1.77	0.66
1:F:612:VAL:HB	1:F:665:VAL:HG12	1.78	0.66
1:E:664:PRO:HG2	1:E:687:ILE:HG12	1.77	0.66
1:D:411:GLY:HA3	1:D:564:ILE:CG1	2.25	0.66
1:A:619:ILE:O	1:A:623:ILE:HG13	1.96	0.66
1:D:628:LYS:HA	1:D:634:SER:HB3	1.77	0.65
1:I:614:LEU:HD12	1:I:706:ARG:HE	1.61	0.65
1:C:432:MET:HE3	1:C:434:LEU:HD11	1.79	0.65
1:E:636:LEU:HD23	1:E:638:LEU:HG	1.79	0.65
1:H:378:LYS:H	1:H:426:LEU:HD22	1.61	0.65
1:C:543:ARG:O	1:C:543:ARG:HG2	1.95	0.64
1:G:410:ILE:HG23	1:G:412:ILE:H	1.61	0.64
1:B:381:SER:O	1:B:385:GLN:HG2	1.96	0.64
1:H:608:SER:HA	1:H:699:ILE:HD11	1.78	0.64
1:B:388:THR:HG22	1:B:389:LYS:H	1.61	0.64
1:A:549:LYS:HE2	1:A:584:THR:HG22	1.79	0.64
1:G:673:ARG:HG3	1:G:674:ARG:N	2.12	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:546:PRO:HD2	1:F:549:LYS:CD	2.13	0.63
1:D:695:ILE:HD12	1:D:701:ILE:HD11	1.78	0.63
1:C:549:LYS:HE2	1:C:584:THR:HG22	1.80	0.63
1:C:670:ILE:HD12	1:C:693:GLN:HB2	1.81	0.63
1:G:412:ILE:HD11	1:G:558:VAL:CB	2.29	0.63
1:H:560:GLU:OE2	1:H:677:ARG:HG2	1.98	0.62
1:B:404:GLN:O	1:B:408:GLN:HG3	1.99	0.62
1:F:472:ALA:HB3	1:F:475:LEU:HG	1.81	0.62
1:I:526:MET:HE3	1:I:567:LEU:HD11	1.81	0.62
1:A:627:ILE:HD12	1:A:635:TYR:HB2	1.82	0.62
1:H:549:LYS:HE2	1:H:584:THR:HG22	1.82	0.62
1:I:410:ILE:CG2	1:I:412:ILE:HD12	2.28	0.62
1:B:412:ILE:HD12	1:B:558:VAL:HG11	1.81	0.62
1:F:616:ASP:H	1:F:707:ILE:HG22	1.65	0.62
1:A:402:MET:HE3	1:A:507:SER:HA	1.80	0.61
1:B:388:THR:CG2	1:B:389:LYS:N	2.63	0.61
1:H:642:SER:HA	1:H:645:LEU:HD12	1.81	0.61
1:E:464:LEU:HD12	1:E:465:PRO:CA	2.29	0.61
1:G:377:GLY:HA3	1:G:426:LEU:HD13	1.82	0.61
1:D:375:GLU:HG2	1:D:420:ARG:HB3	1.83	0.61
1:D:406:LEU:O	1:D:410:ILE:HG22	2.01	0.61
1:E:455:VAL:HG12	1:E:456:GLU:N	2.16	0.61
1:I:614:LEU:HB2	1:I:706:ARG:CG	2.30	0.61
1:A:388:THR:OG1	1:A:392:GLN:CG	2.45	0.60
1:H:526:MET:HE1	1:H:554:PHE:CD2	2.37	0.60
1:E:375:GLU:HG2	1:E:420:ARG:HB2	1.83	0.60
1:A:623:ILE:HG22	1:A:623:ILE:O	2.02	0.59
1:B:388:THR:CG2	1:B:389:LYS:H	2.15	0.59
1:A:536:ASP:O	1:A:540[B]:GLU:HG2	2.03	0.59
1:I:664:PRO:HG2	1:I:687:ILE:HG12	1.85	0.59
1:F:612:VAL:CB	1:F:665:VAL:HG12	2.33	0.58
1:G:423:SER:HB3	1:G:426:LEU:HD23	1.83	0.58
1:G:432:MET:HE3	1:G:434:LEU:HD11	1.85	0.58
1:G:583:ASP:OD1	1:G:585:VAL:HG12	2.03	0.58
1:C:673:ARG:HG3	1:C:674:ARG:N	2.17	0.58
1:D:619:ILE:O	1:D:623:ILE:HG12	2.03	0.58
1:E:541:VAL:HG11	1:E:571:LEU:HD22	1.85	0.58
1:F:585:VAL:O	1:F:589:GLU:HG2	2.03	0.58
1:E:465:PRO:HG3	1:E:481:SER:H	1.68	0.58
1:I:433:ILE:HB	1:I:441:VAL:HG12	1.85	0.58
1:F:465:PRO:HB2	1:F:481:SER:H	1.68	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:373:ILE:HB	1:H:434:LEU:HB2	1.84	0.58
1:E:378:LYS:N	1:E:428:GLY:HA2	2.18	0.57
1:C:628:LYS:CG	1:C:634:SER:OG	2.46	0.57
1:H:402:MET:HE1	1:H:506:LEU:HB3	1.87	0.57
1:I:365:ASP:OD1	1:I:365:ASP:C	2.47	0.57
1:F:411:GLY:H	1:F:564:ILE:H	1.51	0.57
1:C:626:ALA:HA	1:C:637:ALA:HB3	1.86	0.57
1:G:377:GLY:CA	1:G:426:LEU:HD13	2.34	0.57
1:B:455:VAL:HG12	1:B:456:GLU:N	2.20	0.57
1:G:412:ILE:CG1	1:G:558:VAL:HG11	2.35	0.57
1:C:410:ILE:HG23	1:C:412:ILE:H	1.69	0.57
1:G:423:SER:N	1:G:426:LEU:HD23	2.20	0.57
1:B:412:ILE:CD1	1:B:558:VAL:HG11	2.35	0.56
1:E:425:SER:O	1:E:426:LEU:HD12	2.06	0.56
1:F:635:TYR:OH	1:F:678:LYS:HE2	2.06	0.56
1:E:464:LEU:CD1	1:E:465:PRO:N	2.51	0.56
1:G:411:GLY:HA3	1:G:564:ILE:CG1	2.36	0.56
1:I:614:LEU:CD1	1:I:706:ARG:NH2	2.57	0.56
1:D:701:ILE:O	1:D:703:PRO:HD3	2.05	0.56
1:H:560:GLU:OE1	1:H:592:ARG:NH1	2.39	0.56
1:C:593:SER:O	1:C:596:LYS:HG3	2.06	0.56
1:G:426:LEU:HD12	1:G:426:LEU:C	2.32	0.55
1:E:410:ILE:HD11	1:E:414:TYR:HE2	1.71	0.55
1:B:375:GLU:HG2	1:B:420:ARG:HB2	1.88	0.55
1:F:364:ASP:O	1:G:413:ARG:NH2	2.33	0.55
1:C:701:ILE:O	1:C:703:PRO:HD3	2.07	0.55
1:I:615:LEU:O	1:I:706:ARG:HD3	2.06	0.55
1:G:414:TYR:CD1	1:G:510:PHE:CD1	2.95	0.55
1:I:375:GLU:HG2	1:I:420:ARG:HB2	1.89	0.55
1:F:636:LEU:HD22	1:F:637:ALA:H	1.72	0.54
1:E:639:ASP:O	1:E:643:VAL:HG23	2.07	0.54
1:I:619:ILE:O	1:I:623:ILE:HG12	2.07	0.54
1:C:398:MET:HE2	1:C:503:ILE:HG13	1.89	0.54
1:I:614:LEU:CB	1:I:706:ARG:HG2	2.35	0.54
1:A:402:MET:CE	1:A:510:PHE:HD2	2.20	0.54
1:C:462:TYR:HB3	1:C:464:LEU:HG	1.90	0.54
1:C:377:GLY:HA3	1:C:427:GLU:O	2.08	0.54
1:E:373:ILE:HB	1:E:434:LEU:HB2	1.90	0.54
1:B:620:GLU:O	1:B:624:ARG:HG3	2.07	0.53
1:F:448:HIS:O	1:F:482:GLU:HG2	2.08	0.53
1:F:515:GLN:HE22	1:F:558:VAL:HG23	1.73	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:543:ARG:O	1:F:543:ARG:HG2	2.07	0.53
1:H:673:ARG:HG3	1:H:674:ARG:N	2.23	0.53
1:I:362:ASN:N	1:I:363:PRO:HD3	2.21	0.53
1:D:593:SER:HB2	1:D:693:GLN:HG2	1.91	0.53
1:I:593:SER:HB2	1:I:693:GLN:HG2	1.91	0.53
1:E:455:VAL:HG12	1:E:457:ASP:H	1.74	0.53
1:F:531:GLU:HG3	1:F:538:VAL:HG11	1.90	0.53
1:H:526:MET:HE1	1:H:554:PHE:HD2	1.74	0.53
1:A:413:ARG:HH21	1:I:365:ASP:HA	1.74	0.52
1:E:628:LYS:HD2	1:E:635:TYR:HB2	1.90	0.52
1:H:642:SER:HA	1:H:645:LEU:HB2	1.91	0.52
1:B:438:VAL:HG11	1:C:597:LEU:HG	1.90	0.52
1:F:662:GLN:N	1:F:663:PRO:HD3	2.24	0.52
1:H:444:LYS:HD2	1:H:469:TYR:OH	2.10	0.52
1:A:593:SER:HB2	1:A:693:GLN:HG2	1.92	0.52
1:A:627:ILE:HD11	1:A:637:ALA:HB2	1.91	0.52
1:I:647:LEU:HB3	1:I:651:ARG:NH1	2.24	0.52
1:F:616:ASP:HB3	1:F:619:ILE:HG12	1.92	0.52
1:G:664:PRO:HG2	1:G:687:ILE:HG12	1.90	0.52
1:A:402:MET:HE3	1:A:507:SER:CA	2.39	0.52
1:E:411:GLY:CA	1:E:564:ILE:HG12	2.33	0.52
1:F:658:PRO:HG2	1:F:662:GLN:HG3	1.90	0.52
1:B:459:LEU:HD21	1:B:466:PHE:HB3	1.92	0.51
1:B:597:LEU:HD12	1:B:696:LEU:HD11	1.92	0.51
1:D:592:ARG:NE	1:D:694:GLU:OE1	2.42	0.51
1:F:612:VAL:HA	1:F:665:VAL:HG12	1.92	0.51
1:H:378:LYS:N	1:H:426:LEU:HD22	2.24	0.51
1:G:379:ASP:HB3	1:G:429:TYR:CZ	2.45	0.51
1:B:576:GLU:HB2	2:B:801:GOL:H2	1.93	0.51
1:E:587:LEU:O	1:E:591:VAL:HG23	2.10	0.51
1:I:388:THR:HG22	3:I:905:HOH:O	2.10	0.51
1:E:628:LYS:HB2	1:E:635:TYR:H	1.75	0.51
1:H:531:GLU:HG3	1:H:538:VAL:HG11	1.92	0.51
1:H:646:ILE:HG22	1:H:650:MET:HE2	1.92	0.51
1:G:592:ARG:NE	1:G:673:ARG:HH21	2.09	0.51
1:E:372:VAL:CG2	1:E:417:ILE:HD13	2.40	0.50
1:G:372:VAL:HG21	1:G:417:ILE:HD13	1.94	0.50
1:C:537:LEU:O	1:C:541:VAL:HG23	2.12	0.50
1:A:619:ILE:HD12	1:A:646:ILE:HG12	1.94	0.50
1:I:411:GLY:HA3	1:I:564:ILE:HG12	1.90	0.50
1:B:648:LYS:HG3	1:B:651:ARG:HH21	1.76	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:658:PRO:HG2	1:B:661:GLY:HA3	1.93	0.50
1:F:465:PRO:CB	1:F:481:SER:H	2.24	0.50
1:I:583:ASP:OD1	1:I:585:VAL:HG12	2.12	0.50
1:E:646:ILE:HG22	1:E:650:MET:HE2	1.94	0.50
1:C:541:VAL:HG21	1:C:571:LEU:HB3	1.94	0.49
1:F:619:ILE:HG13	1:F:707:ILE:CG2	2.42	0.49
1:H:426:LEU:HD13	1:H:428:GLY:N	2.27	0.49
1:E:602:LYS:HD3	1:E:603:PHE:CZ	2.47	0.49
1:F:426:LEU:HD11	1:F:432:MET:HB2	1.95	0.49
1:I:432:MET:HE3	1:I:434:LEU:HD11	1.94	0.49
1:E:452:THR:HG23	1:E:454:GLU:H	1.78	0.49
1:E:464:LEU:HD12	1:E:464:LEU:C	2.34	0.49
1:D:593:SER:O	1:D:596:LYS:HG3	2.12	0.49
1:H:466:PHE:HA	1:H:479:TRP:O	2.13	0.49
1:E:372:VAL:HG23	1:E:417:ILE:HA	1.94	0.49
1:F:672:VAL:O	1:F:676:VAL:HG23	2.12	0.49
1:E:455:VAL:CG1	1:E:456:GLU:N	2.76	0.49
1:E:528:GLU:HG3	1:E:529:PHE:N	2.27	0.49
1:G:377:GLY:HA2	1:G:426:LEU:HD22	0.61	0.49
1:H:614:LEU:CD2	1:H:667:LEU:HB3	2.42	0.49
1:F:657:THR:CG2	1:F:662:GLN:O	2.58	0.49
1:H:410:ILE:HG23	1:H:412:ILE:H	1.77	0.49
1:A:396:ASP:O	1:A:400:PRO:HG2	2.13	0.48
1:G:380:LEU:HG	1:G:429:TYR:HA	1.94	0.48
1:H:375:GLU:HG2	1:H:420:ARG:CB	2.40	0.48
1:B:364:ASP:O	1:C:413:ARG:HD2	2.13	0.48
1:B:541:VAL:CG2	1:B:575:SER:HB2	2.43	0.48
1:E:636:LEU:O	1:E:638:LEU:HD12	2.13	0.48
1:H:445:ILE:HG21	1:H:498:PRO:HB3	1.96	0.48
1:G:412:ILE:HG12	1:G:558:VAL:HG11	1.94	0.48
1:B:541:VAL:HG11	1:B:571:LEU:CD2	2.40	0.48
1:C:674:ARG:O	1:C:677:ARG:HB3	2.13	0.48
1:E:523:VAL:HG11	1:E:551:THR:HG22	1.96	0.48
1:B:455:VAL:CG1	1:B:456:GLU:N	2.77	0.48
1:F:589:GLU:CD	1:F:674:ARG:HB3	2.39	0.48
1:F:657:THR:HB	1:F:662:GLN:CB	2.40	0.48
1:D:432:MET:HE3	1:D:434:LEU:HD11	1.96	0.47
1:E:636:LEU:HG	1:E:638:LEU:HG	1.96	0.47
1:F:524:ARG:HB2	1:F:547:LEU:HD21	1.95	0.47
1:G:430:ASP:CG	1:G:444:LYS:HG2	2.39	0.47
1:E:576:GLU:HB2	2:E:801:GOL:H32	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:429:TYR:CZ	1:F:447:PRO:HB3	2.49	0.47
1:G:394:PHE:HB2	1:G:499:LEU:HD13	1.96	0.47
1:H:402:MET:HE3	1:H:507:SER:HA	1.96	0.47
1:C:474:GLY:O	1:C:475:LEU:HD23	2.14	0.47
1:D:410:ILE:HD11	1:D:558:VAL:HB	1.95	0.47
1:G:540:GLU:OE1	1:G:540:GLU:HA	2.14	0.47
1:B:628:LYS:HD2	1:B:635:TYR:CE2	2.49	0.47
1:G:455:VAL:HG12	1:G:456:GLU:N	2.30	0.47
1:E:432:MET:HE3	1:E:434:LEU:HD11	1.96	0.47
1:F:523:VAL:HA	1:F:526:MET:CE	2.44	0.47
1:I:451:LEU:HD13	1:I:479:TRP:CE2	2.50	0.47
1:D:410:ILE:HG21	1:D:414:TYR:OH	2.15	0.47
1:D:411:GLY:HA3	1:D:564:ILE:HG12	1.96	0.47
1:G:593:SER:HB2	1:G:693:GLN:HG2	1.97	0.47
1:H:596:LYS:HE3	1:H:693:GLN:O	2.15	0.47
1:A:432:MET:HE3	1:A:434:LEU:HD11	1.95	0.47
1:A:445:ILE:HG21	1:A:498:PRO:HB3	1.96	0.47
1:A:579:GLN:NE2	1:I:548:GLN:HG3	2.30	0.47
1:A:664:PRO:HG2	1:A:687:ILE:HG12	1.96	0.47
1:B:388:THR:HG22	1:B:392:GLN:HB3	1.96	0.47
1:C:597:LEU:HD12	1:C:696:LEU:HD11	1.97	0.47
1:D:435:LEU:HD21	1:D:516:GLU:HG3	1.97	0.47
1:G:378:LYS:CG	1:G:426:LEU:HD11	2.41	0.47
1:A:377:GLY:HA3	1:A:427:GLU:O	2.15	0.47
1:C:452:THR:O	1:C:478:ALA:HB3	2.14	0.47
1:C:651:ARG:O	1:C:654:ILE:HG13	2.15	0.47
1:E:373:ILE:HD13	1:E:418:HIS:HB2	1.97	0.47
1:F:630:THR:HG22	1:F:634:SER:HA	1.96	0.47
1:H:524:ARG:HB2	1:H:547:LEU:HD21	1.97	0.47
1:I:488:LEU:HD22	1:I:493:ILE:HG21	1.96	0.47
1:F:366:TYR:HD2	1:G:407:TYR:OH	1.97	0.47
1:F:380:LEU:HD13	1:F:502:ILE:HD11	1.97	0.47
1:H:380:LEU:HD12	1:H:430:ASP:O	2.16	0.47
1:C:593:SER:HB2	1:C:693:GLN:HG2	1.97	0.46
1:D:646:ILE:HG22	1:D:650:MET:HE2	1.96	0.46
1:E:593:SER:HB2	1:E:693:GLN:HG2	1.97	0.46
1:F:523:VAL:HA	1:F:526:MET:HE3	1.98	0.46
1:F:373:ILE:HG12	1:F:418:HIS:HB2	1.96	0.46
1:G:399:ILE:HB	1:G:400:PRO:HD3	1.97	0.46
1:G:673:ARG:NH1	1:G:674:ARG:HA	2.31	0.46
1:H:672:VAL:HA	1:H:675:TYR:HD2	1.81	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:388:THR:HG23	1:H:391:GLY:HA3	1.98	0.46
1:G:512:LYS:NZ	1:G:686:ASP:OD1	2.44	0.46
1:A:643:VAL:HG22	1:A:679:LEU:HD11	1.98	0.46
1:D:653:THR:HG21	1:D:706:ARG:O	2.16	0.46
1:G:592:ARG:CD	1:G:673:ARG:HH21	2.29	0.46
1:A:596:LYS:HE3	1:A:693:GLN:O	2.16	0.46
1:E:458:ASN:O	1:E:461:ARG:HG3	2.16	0.46
1:H:431:TYR:CZ	1:H:443:GLY:HA3	2.51	0.46
1:A:407:TYR:CE2	1:A:565:LYS:HE2	2.51	0.46
1:G:423:SER:CA	1:G:426:LEU:HD23	2.46	0.46
1:H:589:GLU:HA	1:H:589:GLU:OE1	2.15	0.46
1:I:614:LEU:HD13	1:I:706:ARG:CZ	2.46	0.46
1:H:378:LYS:HG3	1:H:426:LEU:HD21	1.98	0.45
1:H:454:GLU:HG2	1:H:493:ILE:HG23	1.98	0.45
1:I:648:LYS:HG2	1:I:651:ARG:NH2	2.31	0.45
1:E:465:PRO:HB2	1:E:480:VAL:HA	1.98	0.45
1:F:411:GLY:CA	1:F:564:ILE:HG12	2.24	0.45
1:G:377:GLY:CA	1:G:426:LEU:CD2	2.42	0.45
1:A:623:ILE:HD11	1:A:646:ILE:HD13	1.98	0.45
1:E:531:GLU:HG3	1:E:538:VAL:HG11	1.98	0.45
1:F:515:GLN:NE2	1:F:558:VAL:HG23	2.31	0.45
1:I:670:ILE:CD1	1:I:693:GLN:HB2	2.47	0.45
1:D:410:ILE:HG12	1:D:412:ILE:HD12	1.99	0.45
1:H:411:GLY:HA3	1:H:564:ILE:HG13	1.98	0.45
1:A:383:LEU:HD12	1:A:383:LEU:HA	1.77	0.45
1:A:614:LEU:HD12	1:A:706:ARG:HE	1.81	0.45
1:G:523:VAL:HG21	1:G:551:THR:HG23	1.99	0.45
1:I:593:SER:O	1:I:596:LYS:HG3	2.16	0.45
1:D:628:LYS:HB3	1:D:628:LYS:HE3	1.78	0.45
1:E:452:THR:HB	1:E:478:ALA:O	2.17	0.45
1:F:635:TYR:OH	1:F:678:LYS:HG2	2.16	0.45
1:H:423:SER:O	1:H:426:LEU:HD23	2.16	0.45
1:B:684:PHE:HB3	1:B:687:ILE:HD12	1.99	0.45
1:F:518:LEU:HD23	1:F:555:LYS:HG2	1.98	0.45
1:A:457:ASP:O	1:A:461:ARG:HG3	2.16	0.45
1:E:520:ILE:HA	1:E:551:THR:HG21	1.99	0.45
1:F:589:GLU:OE1	1:F:674:ARG:HB3	2.16	0.45
1:F:681:GLU:HA	1:F:685:PRO:HA	1.99	0.45
1:G:373:ILE:HD13	1:G:418:HIS:HB2	1.97	0.45
1:F:365:ASP:HA	1:G:413:ARG:NH2	2.32	0.45
1:G:531:GLU:HG3	1:G:538:VAL:HG11	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:468:THR:HA	1:I:477:SER:O	2.17	0.45
1:C:655:THR:O	1:C:655:THR:CG2	2.58	0.44
1:I:456:GLU:HB3	1:I:466:PHE:CZ	2.52	0.44
1:C:445:ILE:HD13	1:C:498:PRO:HB3	1.98	0.44
1:H:380:LEU:HD23	1:H:383:LEU:HD12	1.99	0.44
1:A:627:ILE:HG22	1:A:629:GLN:HG3	1.98	0.44
1:B:497:THR:O	1:B:501:VAL:HG23	2.17	0.44
1:C:375:GLU:HG2	1:C:420:ARG:HB2	1.98	0.44
1:G:497:THR:OG1	1:G:500:GLU:HG3	2.17	0.44
1:H:530:MET:HG2	1:H:567:LEU:HD13	2.00	0.44
1:I:445:ILE:HG21	1:I:498:PRO:HB3	2.00	0.44
1:E:470:LYS:HG2	1:E:472:ALA:O	2.17	0.44
1:F:638:LEU:HD23	1:F:638:LEU:HA	1.86	0.44
1:E:567:LEU:HD23	1:E:567:LEU:HA	1.84	0.44
1:G:410:ILE:HG23	1:G:412:ILE:N	2.29	0.44
1:H:665:VAL:HG13	1:H:690:ILE:HD13	1.99	0.44
1:A:529:PHE:CE1	1:I:366:TYR:O	2.71	0.44
1:C:396:ASP:O	1:C:400:PRO:HG2	2.17	0.44
1:I:515:GLN:NE2	1:I:559:GLN:HB2	2.32	0.44
1:B:530:MET:HG2	1:B:567:LEU:HD13	2.00	0.44
1:H:402:MET:HE3	1:H:507:SER:CA	2.47	0.44
1:H:402:MET:HE2	1:H:406:LEU:HD11	2.00	0.44
1:H:494:LYS:HD2	1:H:496:TRP:CZ2	2.52	0.44
1:A:628:LYS:HE3	1:A:628:LYS:HB2	1.76	0.44
1:B:616:ASP:OD2	1:B:708:GLN:HG3	2.18	0.44
1:E:557:LEU:HD11	1:E:591:VAL:HG11	1.99	0.44
1:G:423:SER:CB	1:G:426:LEU:HD23	2.47	0.44
1:I:614:LEU:CD1	1:I:706:ARG:NE	2.78	0.44
1:C:615:LEU:HD23	1:C:707:ILE:HB	2.00	0.44
1:F:465:PRO:HB2	1:F:481:SER:N	2.32	0.44
1:I:429:TYR:CE2	1:I:447:PRO:HB3	2.53	0.44
1:A:399:ILE:HB	1:A:400:PRO:HD3	2.00	0.43
1:A:402:MET:HE2	1:A:510:PHE:HD2	1.83	0.43
1:C:423:SER:HB3	1:C:426:LEU:HB2	2.00	0.43
1:E:379:ASP:HB2	1:E:428:GLY:O	2.18	0.43
1:E:636:LEU:HD23	1:E:638:LEU:CG	2.48	0.43
1:G:616:ASP:HB2	1:G:707:ILE:O	2.19	0.43
1:H:368:LEU:HD12	1:H:368:LEU:HA	1.83	0.43
1:C:646:ILE:HG22	1:C:650:MET:HE2	2.00	0.43
1:D:616:ASP:HB2	1:D:707:ILE:O	2.18	0.43
1:F:446:PRO:HA	1:F:447:PRO:HD3	1.79	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:639:ASP:O	1:C:643:VAL:HG23	2.17	0.43
1:E:459:LEU:C	1:E:461:ARG:H	2.27	0.43
1:G:557:LEU:HD11	1:G:591:VAL:HG11	2.00	0.43
1:A:446:PRO:HA	1:A:447:PRO:HD3	1.92	0.43
1:E:520:ILE:HD11	1:E:548:GLN:HG2	2.01	0.43
1:G:619:ILE:HA	1:G:622:MET:HE2	2.01	0.43
1:B:388:THR:CG2	1:B:392:GLN:HB3	2.48	0.43
1:B:429:TYR:CD2	1:B:447:PRO:HG3	2.53	0.43
1:B:670:ILE:CD1	1:B:693:GLN:HB2	2.47	0.43
1:F:526:MET:HE1	1:F:554:PHE:HD2	1.84	0.43
1:B:664:PRO:HG2	1:B:687:ILE:HG12	2.01	0.43
1:F:624:ARG:C	1:F:637:ALA:HB3	2.44	0.43
1:I:484:ALA:O	1:I:488:LEU:HG	2.19	0.43
1:I:648:LYS:HA	1:I:651:ARG:NH2	2.34	0.43
1:B:582:LYS:HA	1:B:582:LYS:HD3	1.77	0.43
1:D:445:ILE:HG21	1:D:498:PRO:HB3	2.00	0.43
1:G:423:SER:H	1:G:426:LEU:HD23	1.84	0.43
1:B:636:LEU:HD21	1:B:679:LEU:HD12	2.01	0.42
1:E:364:ASP:CG	1:F:403:ARG:HH22	2.27	0.42
1:E:636:LEU:CG	1:E:638:LEU:HG	2.50	0.42
1:F:587:LEU:O	1:F:591:VAL:HG23	2.19	0.42
1:F:673:ARG:HG3	1:F:674:ARG:N	2.33	0.42
1:I:526:MET:CE	1:I:567:LEU:HD21	2.50	0.42
1:E:423:SER:OG	1:E:426:LEU:HD13	2.20	0.42
1:E:472:ALA:HB3	1:E:475:LEU:HG	2.01	0.42
1:E:523:VAL:HA	1:E:526:MET:HG2	2.01	0.42
1:F:375:GLU:HG2	1:F:420:ARG:HB2	2.00	0.42
1:F:570:ILE:HG23	1:F:591:VAL:HG11	2.02	0.42
1:F:612:VAL:CA	1:F:665:VAL:HG12	2.49	0.42
1:G:451:LEU:HD13	1:G:479:TRP:CE2	2.55	0.42
1:A:402:MET:HE2	1:A:406:LEU:HD11	2.02	0.42
1:D:399:ILE:HB	1:D:400:PRO:HD3	2.02	0.42
1:G:414:TYR:CE1	1:G:510:PHE:HB3	2.55	0.42
1:H:427:GLU:C	1:H:429:TYR:H	2.28	0.42
1:C:459:LEU:HD21	1:C:478:ALA:HB1	2.02	0.42
1:C:497:THR:O	1:C:501:VAL:HG23	2.19	0.42
1:E:452:THR:CG2	1:E:454:GLU:H	2.32	0.42
1:E:515:GLN:NE2	1:E:559:GLN:HB2	2.35	0.42
1:F:610:ILE:HG12	1:F:612:VAL:HG13	2.02	0.42
1:F:624:ARG:O	1:F:637:ALA:HB3	2.19	0.42
1:B:407:TYR:O	1:B:407:TYR:CG	2.73	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:411:GLY:HA3	1:C:564:ILE:HG13	2.02	0.42
1:D:377:GLY:HA3	1:D:427:GLU:O	2.20	0.42
1:F:398:MET:HB3	1:F:503:ILE:HG13	2.01	0.42
1:B:639:ASP:O	1:B:643:VAL:HG23	2.20	0.42
1:G:445:ILE:HD13	1:G:498:PRO:HB3	2.01	0.42
1:B:411:GLY:O	1:B:564:ILE:HG13	2.20	0.42
1:C:462:TYR:CZ	1:C:487:ILE:HG12	2.55	0.42
1:H:451:LEU:HD13	1:H:479:TRP:CE2	2.55	0.42
1:I:373:ILE:HB	1:I:434:LEU:HB2	2.02	0.42
1:A:397:ASP:O	1:A:401:LYS:HG2	2.20	0.41
1:C:614:LEU:HD23	1:C:667:LEU:HB3	2.01	0.41
1:F:381:SER:HB3	1:F:421:THR:HB	2.01	0.41
1:G:568:ARG:O	1:G:572:GLU:HB2	2.20	0.41
1:H:690:ILE:HG22	1:H:691:SER:N	2.35	0.41
2:A:802:GOL:C3	1:I:548:GLN:HE21	2.32	0.41
1:H:615:LEU:HG	1:H:666:LEU:HD11	2.02	0.41
1:B:445:ILE:HG21	1:B:498:PRO:HB3	2.02	0.41
1:D:557:LEU:HD11	1:D:591:VAL:HG11	2.03	0.41
1:E:419:VAL:HG13	1:E:419:VAL:O	2.20	0.41
1:E:455:VAL:CG1	1:E:456:GLU:H	2.33	0.41
1:B:542:THR:O	1:B:542:THR:HG22	2.20	0.41
1:C:526:MET:HE2	1:C:526:MET:HB3	1.79	0.41
1:C:589:GLU:HA	1:C:589:GLU:OE1	2.19	0.41
1:D:639:ASP:O	1:D:643:VAL:HG23	2.21	0.41
1:E:408:GLN:OE1	1:E:565:LYS:HE3	2.20	0.41
1:E:684:PHE:HB3	1:E:687:ILE:HD12	2.02	0.41
1:F:529:PHE:HD1	1:F:532:ARG:HH21	1.68	0.41
1:B:455:VAL:CG1	1:B:456:GLU:H	2.33	0.41
1:I:603:PHE:HB3	1:I:665:VAL:HG21	2.02	0.41
1:A:426:LEU:HD22	1:A:430:ASP:HB3	2.02	0.41
1:B:619:ILE:HD12	1:B:646:ILE:HG12	2.01	0.41
1:B:645:LEU:HD23	1:B:709:ILE:HG21	2.03	0.41
1:C:528:GLU:O	1:C:531:GLU:HB2	2.21	0.41
1:D:453:ASN:HB3	1:D:494:LYS:HE3	2.01	0.41
1:D:643:VAL:O	1:D:647:LEU:HD13	2.20	0.41
1:A:383:LEU:O	1:A:499:LEU:HD11	2.19	0.41
1:B:444:LYS:HB2	1:B:469:TYR:OH	2.21	0.41
1:D:592:ARG:HB3	1:D:694:GLU:OE1	2.21	0.41
1:C:567:LEU:HA	1:C:567:LEU:HD23	1.77	0.41
1:C:670:ILE:CD1	1:C:693:GLN:HB2	2.49	0.41
1:E:364:ASP:CG	1:E:364:ASP:O	2.63	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:380:LEU:HD23	1:F:380:LEU:HA	1.88	0.41
1:F:534:PHE:HB3	1:F:537:LEU:HB3	2.03	0.41
1:H:552:GLU:O	1:H:556:ARG:HG3	2.21	0.41
1:H:618:GLU:O	1:H:622:MET:HB2	2.20	0.41
1:C:438:VAL:HG11	1:D:597:LEU:HG	2.03	0.41
1:E:676:VAL:O	1:E:680:ILE:HG23	2.21	0.41
1:F:619:ILE:HG13	1:F:707:ILE:HG23	2.03	0.41
1:H:548:GLN:HG2	1:I:579:GLN:NE2	2.36	0.41
1:A:602:LYS:HD3	1:A:603:PHE:CZ	2.55	0.41
1:D:383:LEU:HA	1:D:383:LEU:HD12	1.78	0.41
1:E:412:ILE:HD12	1:E:412:ILE:H	1.86	0.41
1:D:446:PRO:HA	1:D:447:PRO:HD3	1.92	0.40
1:G:426:LEU:HD13	1:G:428:GLY:N	2.31	0.40
1:I:531:GLU:HG3	1:I:538:VAL:HG11	2.03	0.40
1:D:696:LEU:O	1:D:699:ILE:HG22	2.22	0.40
1:E:459:LEU:HD23	1:E:459:LEU:HA	1.79	0.40
1:B:411:GLY:HA3	1:B:564:ILE:HG13	2.04	0.40
1:C:602:LYS:HD3	1:C:603:PHE:CE1	2.57	0.40
1:C:619:ILE:O	1:C:623:ILE:HG12	2.21	0.40
1:F:399:ILE:HB	1:F:400:PRO:HD3	2.03	0.40
1:F:593:SER:HB2	1:F:693:GLN:HG2	2.04	0.40
1:H:593:SER:HB2	1:H:693:GLN:HG2	2.02	0.40
1:I:446:PRO:HA	1:I:447:PRO:HD3	1.91	0.40
1:B:602:LYS:HD3	1:B:603:PHE:CZ	2.56	0.40
1:B:651:ARG:NH1	1:B:683:GLU:OE2	2.54	0.40
1:A:650:MET:HG3	1:A:707:ILE:HD12	2.04	0.40
1:B:605:GLN:NE2	1:B:662:GLN:HB2	2.36	0.40
1:C:417:ILE:HD11	1:C:510:PHE:HE2	1.86	0.40
1:E:427:GLU:HG3	1:E:430:ASP:OD2	2.21	0.40
1:F:655:THR:HA	1:F:656:PRO:HD3	1.95	0.40
1:G:446:PRO:HA	1:G:447:PRO:HD3	1.95	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	351/387 (91%)	344 (98%)	7 (2%)	0	100	100
1	B	345/387 (89%)	331 (96%)	14 (4%)	0	100	100
1	C	336/387 (87%)	321 (96%)	15 (4%)	0	100	100
1	D	341/387 (88%)	329 (96%)	12 (4%)	0	100	100
1	E	339/387 (88%)	323 (95%)	16 (5%)	0	100	100
1	F	337/387 (87%)	313 (93%)	24 (7%)	0	100	100
1	G	340/387 (88%)	329 (97%)	11 (3%)	0	100	100
1	H	314/387 (81%)	300 (96%)	14 (4%)	0	100	100
1	I	331/387 (86%)	320 (97%)	11 (3%)	0	100	100
All	All	3034/3483 (87%)	2910 (96%)	124 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	323/345 (94%)	310 (96%)	13 (4%)	28	63
1	B	317/345 (92%)	299 (94%)	18 (6%)	18	49
1	C	314/345 (91%)	303 (96%)	11 (4%)	32	67
1	D	318/345 (92%)	309 (97%)	9 (3%)	38	73
1	E	316/345 (92%)	303 (96%)	13 (4%)	27	62
1	F	313/345 (91%)	297 (95%)	16 (5%)	21	54
1	G	314/345 (91%)	302 (96%)	12 (4%)	29	64
1	H	299/345 (87%)	290 (97%)	9 (3%)	36	72
1	I	313/345 (91%)	301 (96%)	12 (4%)	29	64
All	All	2827/3105 (91%)	2714 (96%)	113 (4%)	28	63

All (113) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	383	LEU
1	A	463	ASN
1	A	494	LYS
1	A	497	THR
1	A	503	ILE
1	A	524	ARG
1	A	530	MET
1	A	540[A]	GLU
1	A	540[B]	GLU
1	A	623	ILE
1	A	634	SER
1	A	639	ASP
1	A	700	ARG
1	B	372	VAL
1	B	409	ASP
1	B	410	ILE
1	B	412	ILE
1	B	442	ARG
1	B	444	LYS
1	B	457	ASP
1	B	528	GLU
1	B	530	MET
1	B	540	GLU
1	B	584	THR
1	B	585	VAL
1	B	618	GLU
1	B	627	ILE
1	B	634	SER
1	B	650	MET
1	B	657	THR
1	B	683	GLU
1	C	365	ASP
1	C	372	VAL
1	C	387	LYS
1	C	426	LEU
1	C	460	SER
1	C	530	MET
1	C	584	THR
1	C	585	VAL
1	C	621	GLU
1	C	638	LEU
1	C	643	VAL

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Mol	Chain	Res	Type
1	D	367	SER
1	D	389	LYS
1	D	408	GLN
1	D	480	VAL
1	D	530	MET
1	D	584	THR
1	D	594	SER
1	D	643	VAL
1	D	650	MET
1	E	365	ASP
1	E	410	ILE
1	E	426	LEU
1	E	452	THR
1	E	528	GLU
1	E	530	MET
1	E	540	GLU
1	E	543	ARG
1	E	551	THR
1	E	607	GLN
1	E	655	THR
1	E	657	THR
1	E	673	ARG
1	F	386	HIS
1	F	421	THR
1	F	461	ARG
1	F	494	LYS
1	F	512	LYS
1	F	520	ILE
1	F	543	ARG
1	F	548	GLN
1	F	564	ILE
1	F	565	LYS
1	F	636	LEU
1	F	651	ARG
1	F	665	VAL
1	F	668	THR
1	F	689	VAL
1	F	700	ARG
1	G	365	ASP
1	G	372	VAL
1	G	383	LEU
1	G	408	GLN

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Mol	Chain	Res	Type
1	G	410	ILE
1	G	412	ILE
1	G	524	ARG
1	G	542	THR
1	G	584	THR
1	G	643	VAL
1	G	662	GLN
1	G	673	ARG
1	H	372	VAL
1	H	388	THR
1	H	392	GLN
1	H	421	THR
1	H	422	ASP
1	H	455	VAL
1	H	575	SER
1	H	584	THR
1	H	585	VAL
1	I	383	LEU
1	I	389	LYS
1	I	395	VAL
1	I	408	GLN
1	I	524	ARG
1	I	530	MET
1	I	543	ARG
1	I	584	THR
1	I	643	VAL
1	I	649	SER
1	I	680	ILE
1	I	708	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	385	GLN
1	B	404	GLN
1	B	463	ASN
1	C	448	HIS
1	C	515	GLN
1	C	579	GLN
1	C	605	GLN
1	D	448	HIS
1	E	385	GLN

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Mol	Chain	Res	Type
1	E	511	HIS
1	E	652	ASN
1	F	548	GLN
1	F	607	GLN
1	F	652	ASN
1	G	708	GLN
1	H	471	ASN
1	H	644	ASN
1	I	448	HIS
1	I	548	GLN
1	I	605	GLN
1	I	708	GLN

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 [i](#)

13 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	GOL	G	801	-	5,5,5	0.08	0	5,5,5	0.32	0
2	GOL	B	802	-	5,5,5	0.09	0	5,5,5	0.31	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	GOL	A	802	-	5,5,5	0.09	0	5,5,5	0.31	0
2	GOL	D	801	-	5,5,5	0.09	0	5,5,5	0.31	0
2	GOL	E	802	-	5,5,5	0.08	0	5,5,5	0.31	0
2	GOL	B	801	-	5,5,5	0.08	0	5,5,5	0.31	0
2	GOL	I	801	-	5,5,5	0.09	0	5,5,5	0.31	0
2	GOL	C	801	-	5,5,5	0.09	0	5,5,5	0.32	0
2	GOL	A	804	-	5,5,5	0.08	0	5,5,5	0.31	0
2	GOL	A	803	-	5,5,5	0.08	0	5,5,5	0.31	0
2	GOL	A	801	-	5,5,5	0.09	0	5,5,5	0.32	0
2	GOL	E	801	-	5,5,5	0.08	0	5,5,5	0.31	0
2	GOL	E	803	-	5,5,5	0.09	0	5,5,5	0.31	0

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	GOL	G	801	-	-	2/4/4/4	-
2	GOL	B	802	-	-	2/4/4/4	-
2	GOL	A	802	-	-	2/4/4/4	-
2	GOL	D	801	-	-	2/4/4/4	-
2	GOL	E	802	-	-	4/4/4/4	-
2	GOL	B	801	-	-	2/4/4/4	-
2	GOL	I	801	-	-	2/4/4/4	-
2	GOL	C	801	-	-	0/4/4/4	-
2	GOL	A	804	-	-	3/4/4/4	-
2	GOL	A	803	-	-	3/4/4/4	-
2	GOL	A	801	-	-	2/4/4/4	-
2	GOL	E	801	-	-	1/4/4/4	-
2	GOL	E	803	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (27) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	802	GOL	O1-C1-C2-C3
2	A	803	GOL	O1-C1-C2-O2
2	E	802	GOL	O1-C1-C2-C3
2	E	802	GOL	C1-C2-C3-O3
2	I	801	GOL	C1-C2-C3-O3
2	G	801	GOL	O2-C2-C3-O3
2	A	801	GOL	C1-C2-C3-O3
2	A	803	GOL	O1-C1-C2-C3
2	A	804	GOL	O1-C1-C2-C3
2	B	801	GOL	O1-C1-C2-C3
2	B	802	GOL	O1-C1-C2-C3
2	D	801	GOL	C1-C2-C3-O3
2	E	803	GOL	O1-C1-C2-C3
2	G	801	GOL	C1-C2-C3-O3
2	A	804	GOL	O1-C1-C2-O2
2	E	802	GOL	O2-C2-C3-O3
2	E	803	GOL	O1-C1-C2-O2
2	A	802	GOL	O1-C1-C2-O2
2	E	802	GOL	O1-C1-C2-O2
2	A	801	GOL	O2-C2-C3-O3
2	A	803	GOL	O2-C2-C3-O3
2	A	804	GOL	O2-C2-C3-O3
2	E	801	GOL	O2-C2-C3-O3
2	D	801	GOL	O2-C2-C3-O3
2	I	801	GOL	O2-C2-C3-O3
2	B	801	GOL	O1-C1-C2-O2
2	B	802	GOL	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	802	GOL	1	0
2	B	801	GOL	1	0
2	E	801	GOL	1	0

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	348/387 (89%)	-0.37	9 (2%) 57 47	27, 60, 109, 201	5 (1%)
1	B	347/387 (89%)	0.02	14 (4%) 42 33	36, 77, 154, 225	0
1	C	342/387 (88%)	0.32	15 (4%) 39 31	48, 104, 185, 241	0
1	D	343/387 (88%)	-0.11	10 (2%) 53 43	27, 75, 149, 224	2 (0%)
1	E	343/387 (88%)	-0.04	10 (2%) 53 43	44, 92, 166, 238	0
1	F	340/387 (87%)	0.32	20 (5%) 28 21	30, 106, 182, 239	1 (0%)
1	G	343/387 (88%)	0.12	18 (5%) 33 25	40, 89, 148, 233	1 (0%)
1	H	320/387 (82%)	0.43	17 (5%) 32 24	62, 117, 189, 247	0
1	I	334/387 (86%)	0.18	13 (3%) 43 34	37, 94, 171, 211	3 (0%)
All	All	3060/3483 (87%)	0.09	126 (4%) 41 33	27, 88, 171, 247	12 (0%)

All (126) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	660	GLY	8.0
1	F	659	ALA	6.8
1	C	709	ILE	6.7
1	A	625	GLY	5.4
1	A	363	PRO	5.3
1	B	408	GLN	5.1
1	B	411	GLY	5.1
1	I	411	GLY	5.0
1	I	640	PRO	4.8
1	F	661	GLY	4.7
1	D	411	GLY	4.6
1	I	626	ALA	4.6
1	F	664	PRO	4.5
1	E	658	PRO	4.3
1	I	410	ILE	4.1

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Mol	Chain	Res	Type	RSRZ
1	C	541	VAL	4.0
1	B	542	THR	4.0
1	F	462	TYR	3.9
1	G	652	ASN	3.7
1	I	363	PRO	3.7
1	B	634	SER	3.6
1	D	632	ALA	3.5
1	H	462	TYR	3.5
1	G	658	PRO	3.5
1	G	411	GLY	3.5
1	H	643	VAL	3.5
1	H	464	LEU	3.4
1	G	414	TYR	3.4
1	C	459	LEU	3.4
1	G	420	ARG	3.4
1	F	635	TYR	3.3
1	G	364	ASP	3.3
1	D	363	PRO	3.3
1	H	426	LEU	3.3
1	B	409	ASP	3.3
1	B	541	VAL	3.3
1	H	463	ASN	3.2
1	A	384	ILE	3.2
1	A	386[A]	HIS	3.2
1	F	658	PRO	3.2
1	F	464	LEU	3.2
1	I	362	ASN	3.1
1	B	632	ALA	3.1
1	H	609	ALA	3.1
1	H	642	SER	3.1
1	H	663	PRO	3.1
1	I	364	ASP	3.1
1	C	617	PRO	3.1
1	F	650	MET	3.1
1	D	627	ILE	3.0
1	B	363	PRO	3.0
1	C	627	ILE	3.0
1	G	412	ILE	3.0
1	C	645	LEU	3.0
1	G	661	GLY	2.9
1	G	413	ARG	2.9
1	I	407	TYR	2.9

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Mol	Chain	Res	Type	RSRZ
1	I	627	ILE	2.9
1	E	638	LEU	2.9
1	C	656	PRO	2.8
1	H	640	PRO	2.8
1	B	659	ALA	2.8
1	H	416	GLY	2.8
1	H	646	ILE	2.7
1	E	494	LYS	2.7
1	G	426	LEU	2.7
1	D	658	PRO	2.6
1	G	678	LYS	2.6
1	A	365	ASP	2.6
1	C	632	ALA	2.6
1	F	472	ALA	2.6
1	H	675	TYR	2.6
1	E	487	ILE	2.6
1	F	623	ILE	2.6
1	G	659	ALA	2.5
1	F	663	PRO	2.5
1	A	623	ILE	2.5
1	H	509	PHE	2.5
1	E	630	THR	2.5
1	D	410	ILE	2.4
1	C	625	GLY	2.4
1	G	660	GLY	2.4
1	A	632	ALA	2.4
1	D	635	TYR	2.4
1	D	630	THR	2.4
1	C	661	GLY	2.4
1	E	541	VAL	2.4
1	I	494	LYS	2.3
1	H	440	TYR	2.3
1	I	645	LEU	2.3
1	B	624	ARG	2.3
1	C	654	ILE	2.3
1	E	364	ASP	2.3
1	B	390	SER	2.3
1	A	626	ALA	2.2
1	E	464	LEU	2.2
1	C	468	THR	2.2
1	F	363	PRO	2.2
1	B	407	TYR	2.2

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Mol	Chain	Res	Type	RSRZ
1	F	440	TYR	2.2
1	H	684	PHE	2.2
1	F	383	LEU	2.2
1	G	650	MET	2.2
1	C	494	LYS	2.2
1	B	365	ASP	2.2
1	E	542	THR	2.2
1	B	466	PHE	2.2
1	A	627	ILE	2.1
1	F	465	PRO	2.1
1	D	633	GLY	2.1
1	C	636	LEU	2.1
1	F	615	LEU	2.1
1	I	541	VAL	2.1
1	D	365	ASP	2.1
1	H	645	LEU	2.1
1	G	635	TYR	2.1
1	F	646	ILE	2.1
1	G	627	ILE	2.1
1	C	657	THR	2.1
1	G	608	SER	2.1
1	E	635	TYR	2.1
1	G	662	GLN	2.0
1	H	427	GLU	2.0
1	I	607	GLN	2.0
1	F	636	LEU	2.0
1	F	471	ASN	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	GOL	C	801	6/6	0.48	0.20	155,158,162,165	0
2	GOL	I	801	6/6	0.66	0.18	105,120,127,131	0
2	GOL	G	801	6/6	0.78	0.30	110,114,116,117	0
2	GOL	E	801	6/6	0.83	0.23	79,94,112,117	0
2	GOL	E	803	6/6	0.83	0.18	117,121,136,139	0
2	GOL	D	801	6/6	0.84	0.12	75,98,101,102	0
2	GOL	A	804	6/6	0.85	0.20	66,79,82,82	0
2	GOL	B	802	6/6	0.87	0.24	99,109,112,113	0
2	GOL	E	802	6/6	0.89	0.20	65,75,84,90	0
2	GOL	A	802	6/6	0.89	0.21	75,101,102,105	0
2	GOL	A	801	6/6	0.90	0.17	85,96,112,118	0
2	GOL	B	801	6/6	0.90	0.18	68,75,85,90	0
2	GOL	A	803	6/6	0.91	0.10	87,98,105,109	0

6.5 Other polymers [i](#)

There are no such residues in this entry.