



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 08:17 AM UTC

PDB ID : 4IHC / pdb_00004ihc
Title : Crystal structure of probable mannonate dehydratase Dd703_0947 (target EFI-502222) from Dickeya dadantii Ech703
Authors : Patskovsky, Y.; Toro, R.; Bhosle, R.; Hillerich, B.; Seidel, R.D.; Washington, E.; Scott Glenn, A.; Chowdhury, S.; Evans, B.; Hammonds, J.; Zencheck, W.D.; Imker, H.J.; Al Obaidi, N.F.; Stead, M.; Love, J.; Gerlt, J.A.; Almo, S.C.; Enzyme Function Initiative (EFI)
Deposited on : 2012-12-18
Resolution : 2.00 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:

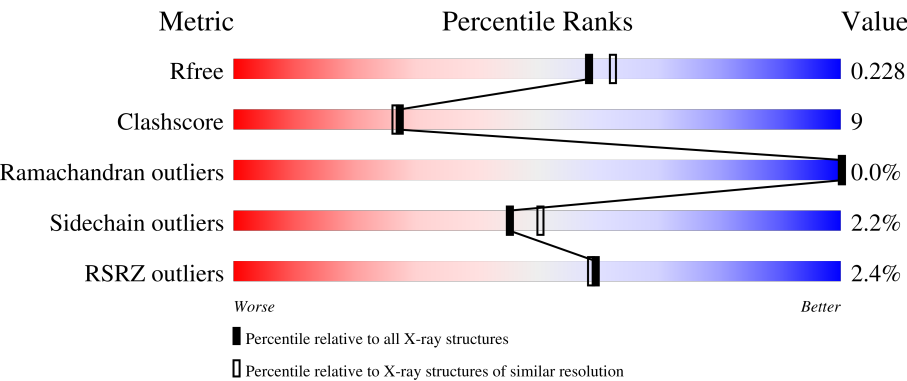
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	440	77%12%• 10%
1	B	440	3% 72%16%• 10%
1	C	440	3% 73%16%• 10%
1	D	440	2% 73%16%• 10%

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Mol	Chain	Length	Quality of chain
1	E	440	 % 72% 17% 10%
1	F	440	 2% 76% 13% 10%
1	G	440	 4% 75% 15% 10%
1	H	440	 % 72% 16% 10%

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	FMT	B	505	-	-	X	-
4	FMT	G	505	-	-	X	-
5	IOD	A	506	-	-	X	-
5	IOD	C	506	-	-	X	-
5	IOD	E	508	-	-	X	-
5	IOD	F	509	-	-	X	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 26763 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 Mandelate racemase/muconate lactonizing protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	396	Total	C	N	O	S	0	0	0
			3084	1964	522	577	21			
1	B	395	Total	C	N	O	S	0	2	0
			3086	1967	521	576	22			
1	C	395	Total	C	N	O	S	0	2	0
			3072	1958	520	573	21			
1	D	395	Total	C	N	O	S	0	1	0
			3077	1960	520	576	21			
1	E	395	Total	C	N	O	S	0	1	0
			3081	1963	521	576	21			
1	F	395	Total	C	N	O	S	0	2	0
			3083	1964	520	578	21			
1	G	396	Total	C	N	O	S	0	1	0
			3079	1962	518	577	22			
1	H	395	Total	C	N	O	S	0	0	0
			3075	1959	520	575	21			

There are 184 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-22	MET	-	expression tag	UNP C6CBG9
A	-21	HIS	-	expression tag	UNP C6CBG9
A	-20	HIS	-	expression tag	UNP C6CBG9
A	-19	HIS	-	expression tag	UNP C6CBG9
A	-18	HIS	-	expression tag	UNP C6CBG9
A	-17	HIS	-	expression tag	UNP C6CBG9
A	-16	HIS	-	expression tag	UNP C6CBG9
A	-15	SER	-	expression tag	UNP C6CBG9
A	-14	SER	-	expression tag	UNP C6CBG9
A	-13	GLY	-	expression tag	UNP C6CBG9
A	-12	VAL	-	expression tag	UNP C6CBG9
A	-11	ASP	-	expression tag	UNP C6CBG9
A	-10	LEU	-	expression tag	UNP C6CBG9

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-9	GLY	-	expression tag	UNP C6CBG9
A	-8	THR	-	expression tag	UNP C6CBG9
A	-7	GLU	-	expression tag	UNP C6CBG9
A	-6	ASN	-	expression tag	UNP C6CBG9
A	-5	LEU	-	expression tag	UNP C6CBG9
A	-4	TYR	-	expression tag	UNP C6CBG9
A	-3	PHE	-	expression tag	UNP C6CBG9
A	-2	GLN	-	expression tag	UNP C6CBG9
A	-1	SER	-	expression tag	UNP C6CBG9
A	0	MET	-	expression tag	UNP C6CBG9
B	-22	MET	-	expression tag	UNP C6CBG9
B	-21	HIS	-	expression tag	UNP C6CBG9
B	-20	HIS	-	expression tag	UNP C6CBG9
B	-19	HIS	-	expression tag	UNP C6CBG9
B	-18	HIS	-	expression tag	UNP C6CBG9
B	-17	HIS	-	expression tag	UNP C6CBG9
B	-16	HIS	-	expression tag	UNP C6CBG9
B	-15	SER	-	expression tag	UNP C6CBG9
B	-14	SER	-	expression tag	UNP C6CBG9
B	-13	GLY	-	expression tag	UNP C6CBG9
B	-12	VAL	-	expression tag	UNP C6CBG9
B	-11	ASP	-	expression tag	UNP C6CBG9
B	-10	LEU	-	expression tag	UNP C6CBG9
B	-9	GLY	-	expression tag	UNP C6CBG9
B	-8	THR	-	expression tag	UNP C6CBG9
B	-7	GLU	-	expression tag	UNP C6CBG9
B	-6	ASN	-	expression tag	UNP C6CBG9
B	-5	LEU	-	expression tag	UNP C6CBG9
B	-4	TYR	-	expression tag	UNP C6CBG9
B	-3	PHE	-	expression tag	UNP C6CBG9
B	-2	GLN	-	expression tag	UNP C6CBG9
B	-1	SER	-	expression tag	UNP C6CBG9
B	0	MET	-	expression tag	UNP C6CBG9
C	-22	MET	-	expression tag	UNP C6CBG9
C	-21	HIS	-	expression tag	UNP C6CBG9
C	-20	HIS	-	expression tag	UNP C6CBG9
C	-19	HIS	-	expression tag	UNP C6CBG9
C	-18	HIS	-	expression tag	UNP C6CBG9
C	-17	HIS	-	expression tag	UNP C6CBG9
C	-16	HIS	-	expression tag	UNP C6CBG9
C	-15	SER	-	expression tag	UNP C6CBG9
C	-14	SER	-	expression tag	UNP C6CBG9

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-13	GLY	-	expression tag	UNP C6CBG9
C	-12	VAL	-	expression tag	UNP C6CBG9
C	-11	ASP	-	expression tag	UNP C6CBG9
C	-10	LEU	-	expression tag	UNP C6CBG9
C	-9	GLY	-	expression tag	UNP C6CBG9
C	-8	THR	-	expression tag	UNP C6CBG9
C	-7	GLU	-	expression tag	UNP C6CBG9
C	-6	ASN	-	expression tag	UNP C6CBG9
C	-5	LEU	-	expression tag	UNP C6CBG9
C	-4	TYR	-	expression tag	UNP C6CBG9
C	-3	PHE	-	expression tag	UNP C6CBG9
C	-2	GLN	-	expression tag	UNP C6CBG9
C	-1	SER	-	expression tag	UNP C6CBG9
C	0	MET	-	expression tag	UNP C6CBG9
D	-22	MET	-	expression tag	UNP C6CBG9
D	-21	HIS	-	expression tag	UNP C6CBG9
D	-20	HIS	-	expression tag	UNP C6CBG9
D	-19	HIS	-	expression tag	UNP C6CBG9
D	-18	HIS	-	expression tag	UNP C6CBG9
D	-17	HIS	-	expression tag	UNP C6CBG9
D	-16	HIS	-	expression tag	UNP C6CBG9
D	-15	SER	-	expression tag	UNP C6CBG9
D	-14	SER	-	expression tag	UNP C6CBG9
D	-13	GLY	-	expression tag	UNP C6CBG9
D	-12	VAL	-	expression tag	UNP C6CBG9
D	-11	ASP	-	expression tag	UNP C6CBG9
D	-10	LEU	-	expression tag	UNP C6CBG9
D	-9	GLY	-	expression tag	UNP C6CBG9
D	-8	THR	-	expression tag	UNP C6CBG9
D	-7	GLU	-	expression tag	UNP C6CBG9
D	-6	ASN	-	expression tag	UNP C6CBG9
D	-5	LEU	-	expression tag	UNP C6CBG9
D	-4	TYR	-	expression tag	UNP C6CBG9
D	-3	PHE	-	expression tag	UNP C6CBG9
D	-2	GLN	-	expression tag	UNP C6CBG9
D	-1	SER	-	expression tag	UNP C6CBG9
D	0	MET	-	expression tag	UNP C6CBG9
E	-22	MET	-	expression tag	UNP C6CBG9
E	-21	HIS	-	expression tag	UNP C6CBG9
E	-20	HIS	-	expression tag	UNP C6CBG9
E	-19	HIS	-	expression tag	UNP C6CBG9
E	-18	HIS	-	expression tag	UNP C6CBG9

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-17	HIS	-	expression tag	UNP C6CBG9
E	-16	HIS	-	expression tag	UNP C6CBG9
E	-15	SER	-	expression tag	UNP C6CBG9
E	-14	SER	-	expression tag	UNP C6CBG9
E	-13	GLY	-	expression tag	UNP C6CBG9
E	-12	VAL	-	expression tag	UNP C6CBG9
E	-11	ASP	-	expression tag	UNP C6CBG9
E	-10	LEU	-	expression tag	UNP C6CBG9
E	-9	GLY	-	expression tag	UNP C6CBG9
E	-8	THR	-	expression tag	UNP C6CBG9
E	-7	GLU	-	expression tag	UNP C6CBG9
E	-6	ASN	-	expression tag	UNP C6CBG9
E	-5	LEU	-	expression tag	UNP C6CBG9
E	-4	TYR	-	expression tag	UNP C6CBG9
E	-3	PHE	-	expression tag	UNP C6CBG9
E	-2	GLN	-	expression tag	UNP C6CBG9
E	-1	SER	-	expression tag	UNP C6CBG9
E	0	MET	-	expression tag	UNP C6CBG9
F	-22	MET	-	expression tag	UNP C6CBG9
F	-21	HIS	-	expression tag	UNP C6CBG9
F	-20	HIS	-	expression tag	UNP C6CBG9
F	-19	HIS	-	expression tag	UNP C6CBG9
F	-18	HIS	-	expression tag	UNP C6CBG9
F	-17	HIS	-	expression tag	UNP C6CBG9
F	-16	HIS	-	expression tag	UNP C6CBG9
F	-15	SER	-	expression tag	UNP C6CBG9
F	-14	SER	-	expression tag	UNP C6CBG9
F	-13	GLY	-	expression tag	UNP C6CBG9
F	-12	VAL	-	expression tag	UNP C6CBG9
F	-11	ASP	-	expression tag	UNP C6CBG9
F	-10	LEU	-	expression tag	UNP C6CBG9
F	-9	GLY	-	expression tag	UNP C6CBG9
F	-8	THR	-	expression tag	UNP C6CBG9
F	-7	GLU	-	expression tag	UNP C6CBG9
F	-6	ASN	-	expression tag	UNP C6CBG9
F	-5	LEU	-	expression tag	UNP C6CBG9
F	-4	TYR	-	expression tag	UNP C6CBG9
F	-3	PHE	-	expression tag	UNP C6CBG9
F	-2	GLN	-	expression tag	UNP C6CBG9
F	-1	SER	-	expression tag	UNP C6CBG9
F	0	MET	-	expression tag	UNP C6CBG9
G	-22	MET	-	expression tag	UNP C6CBG9

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Chain	Residue	Modelled	Actual	Comment	Reference
G	-21	HIS	-	expression tag	UNP C6CBG9
G	-20	HIS	-	expression tag	UNP C6CBG9
G	-19	HIS	-	expression tag	UNP C6CBG9
G	-18	HIS	-	expression tag	UNP C6CBG9
G	-17	HIS	-	expression tag	UNP C6CBG9
G	-16	HIS	-	expression tag	UNP C6CBG9
G	-15	SER	-	expression tag	UNP C6CBG9
G	-14	SER	-	expression tag	UNP C6CBG9
G	-13	GLY	-	expression tag	UNP C6CBG9
G	-12	VAL	-	expression tag	UNP C6CBG9
G	-11	ASP	-	expression tag	UNP C6CBG9
G	-10	LEU	-	expression tag	UNP C6CBG9
G	-9	GLY	-	expression tag	UNP C6CBG9
G	-8	THR	-	expression tag	UNP C6CBG9
G	-7	GLU	-	expression tag	UNP C6CBG9
G	-6	ASN	-	expression tag	UNP C6CBG9
G	-5	LEU	-	expression tag	UNP C6CBG9
G	-4	TYR	-	expression tag	UNP C6CBG9
G	-3	PHE	-	expression tag	UNP C6CBG9
G	-2	GLN	-	expression tag	UNP C6CBG9
G	-1	SER	-	expression tag	UNP C6CBG9
G	0	MET	-	expression tag	UNP C6CBG9
H	-22	MET	-	expression tag	UNP C6CBG9
H	-21	HIS	-	expression tag	UNP C6CBG9
H	-20	HIS	-	expression tag	UNP C6CBG9
H	-19	HIS	-	expression tag	UNP C6CBG9
H	-18	HIS	-	expression tag	UNP C6CBG9
H	-17	HIS	-	expression tag	UNP C6CBG9
H	-16	HIS	-	expression tag	UNP C6CBG9
H	-15	SER	-	expression tag	UNP C6CBG9
H	-14	SER	-	expression tag	UNP C6CBG9
H	-13	GLY	-	expression tag	UNP C6CBG9
H	-12	VAL	-	expression tag	UNP C6CBG9
H	-11	ASP	-	expression tag	UNP C6CBG9
H	-10	LEU	-	expression tag	UNP C6CBG9
H	-9	GLY	-	expression tag	UNP C6CBG9
H	-8	THR	-	expression tag	UNP C6CBG9
H	-7	GLU	-	expression tag	UNP C6CBG9
H	-6	ASN	-	expression tag	UNP C6CBG9
H	-5	LEU	-	expression tag	UNP C6CBG9
H	-4	TYR	-	expression tag	UNP C6CBG9
H	-3	PHE	-	expression tag	UNP C6CBG9

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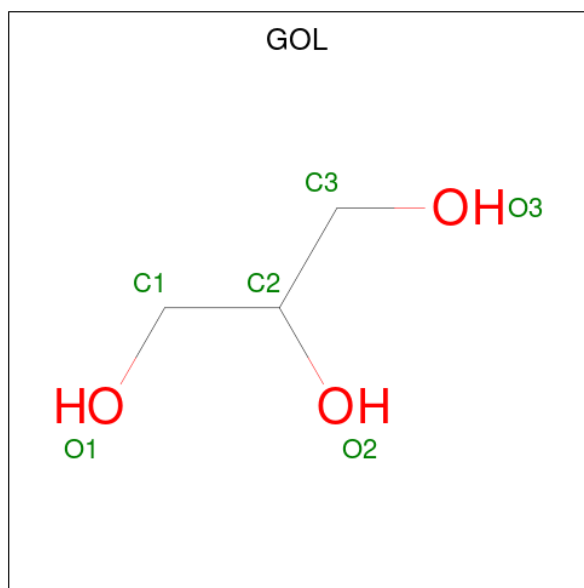
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Chain	Residue	Modelled	Actual	Comment	Reference
H	-2	GLN	-	expression tag	UNP C6CBG9
H	-1	SER	-	expression tag	UNP C6CBG9
H	0	MET	-	expression tag	UNP C6CBG9

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

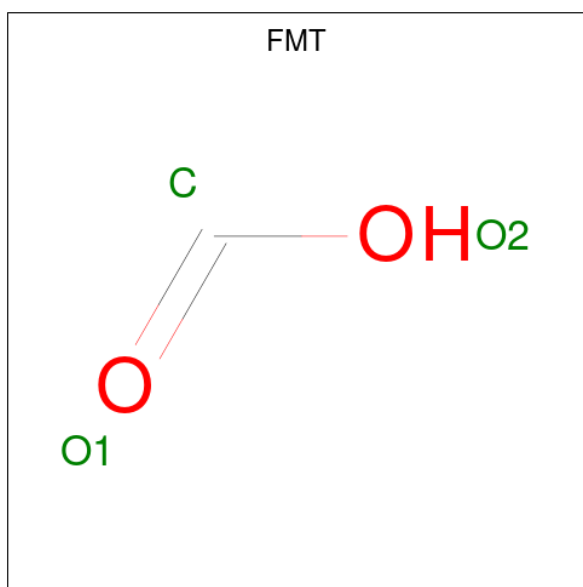
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Mg 1 1	0	0
2	B	1	Total Mg 1 1	0	0
2	C	1	Total Mg 1 1	0	0
2	D	1	Total Mg 1 1	0	0
2	E	1	Total Mg 1 1	0	0
2	F	1	Total Mg 1 1	0	0
2	G	1	Total Mg 1 1	0	0
2	H	1	Total Mg 1 1	0	0

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 6 3 3	0	0
3	A	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0
3	C	1	Total C O 6 3 3	0	0
3	C	1	Total C O 6 3 3	0	0
3	D	1	Total C O 6 3 3	0	0
3	D	1	Total C O 6 3 3	0	0
3	E	1	Total C O 6 3 3	0	0
3	F	1	Total C O 6 3 3	0	0
3	F	1	Total C O 6 3 3	0	0
3	G	1	Total C O 6 3 3	0	0
3	H	1	Total C O 6 3 3	0	0
3	H	1	Total C O 6 3 3	0	0

- Molecule 4 is FORMIC ACID (CCD ID: FMT) (formula: CH₂O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 3 1 2	0	0
4	B	1	Total C O 3 1 2	0	0
4	C	1	Total C O 3 1 2	0	0
4	D	1	Total C O 3 1 2	0	0
4	E	1	Total C O 3 1 2	0	0
4	F	1	Total C O 3 1 2	0	0
4	F	1	Total C O 3 1 2	0	0
4	G	1	Total C O 3 1 2	0	0
4	H	1	Total C O 3 1 2	0	0

- Molecule 5 is IODIDE ION (CCD ID: IOD) (formula: I).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	2	Total I 2 2	0	0
5	B	3	Total I 3 3	0	0
5	C	3	Total I 3 3	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	D	3	Total I 3 3	0	0
5	E	3	Total I 3 3	0	0
5	F	3	Total I 3 3	0	0
5	H	2	Total I 2 2	0	0

- Molecule 6 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

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

- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	244	Total O 244 244	0	0
7	B	255	Total O 255 255	0	0
7	C	237	Total O 237 237	0	0
7	D	237	Total O 237 237	0	0
7	E	275	Total O 276 276	0	1
7	F	239	Total O 239 239	0	0

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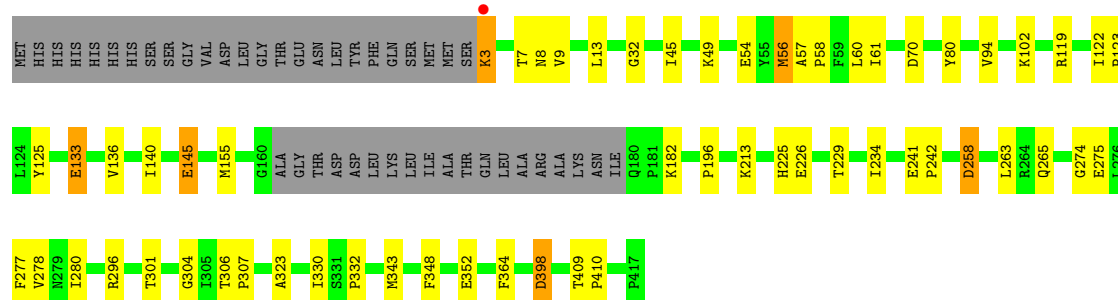
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	G	234	Total 234	O 234	0	0
7	H	257	Total 257	O 257	0	0

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.

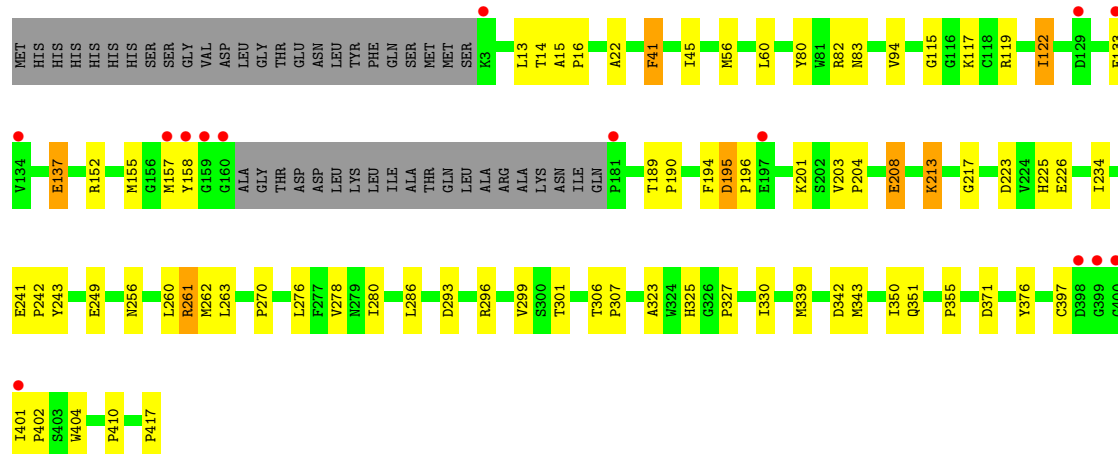
- Molecule 1: Mandelate racemase/muconate lactonizing protein

Chain A: 



- Molecule 1: Mandelate racemase/muconate lactonizing protein

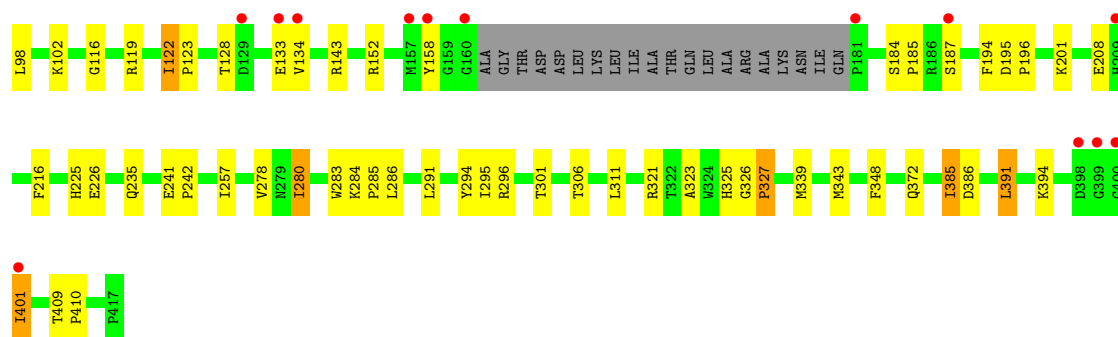
Chain B: 



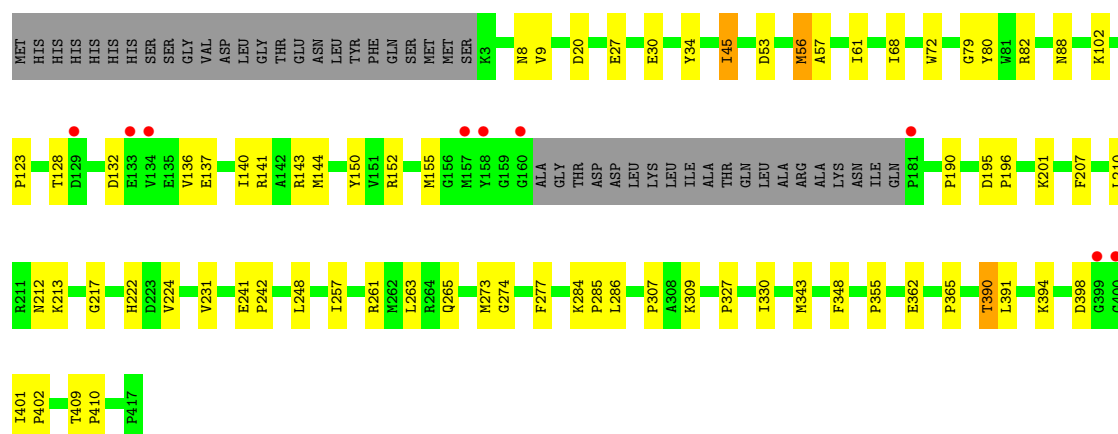
- Molecule 1: Mandelate racemase/muconate lactonizing protein

Chain C: 

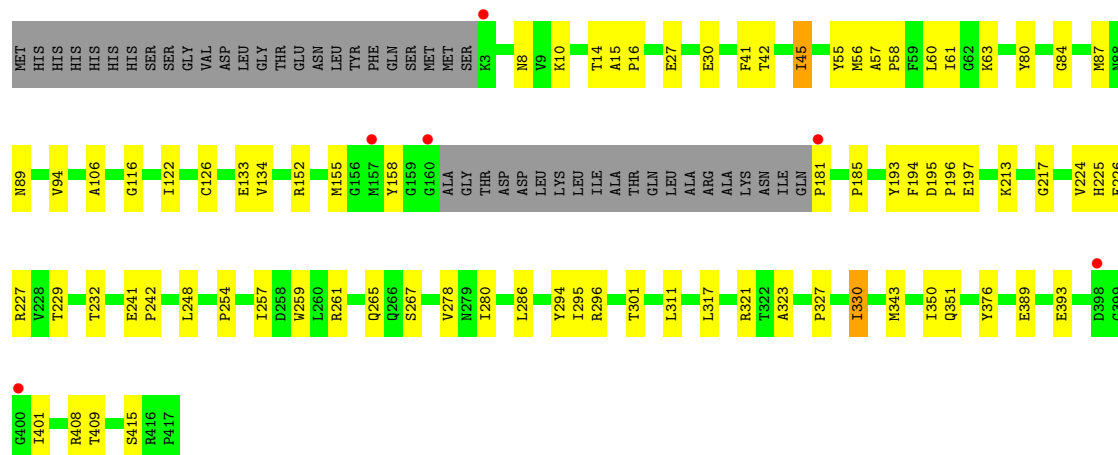
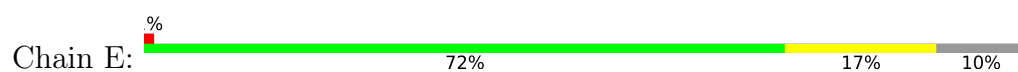




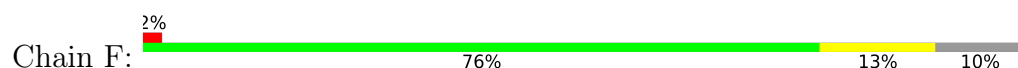
- Molecule 1: Mandelate racemase/muconate lactonizing protein

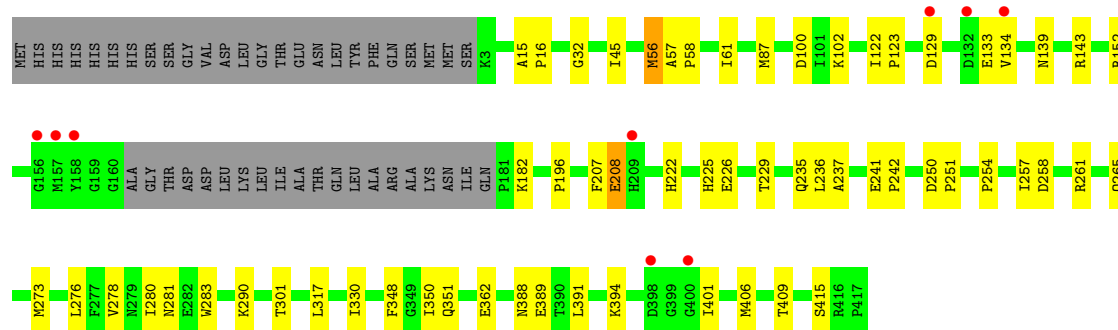


- Molecule 1: Mandelate racemase/muconate lactonizing protein

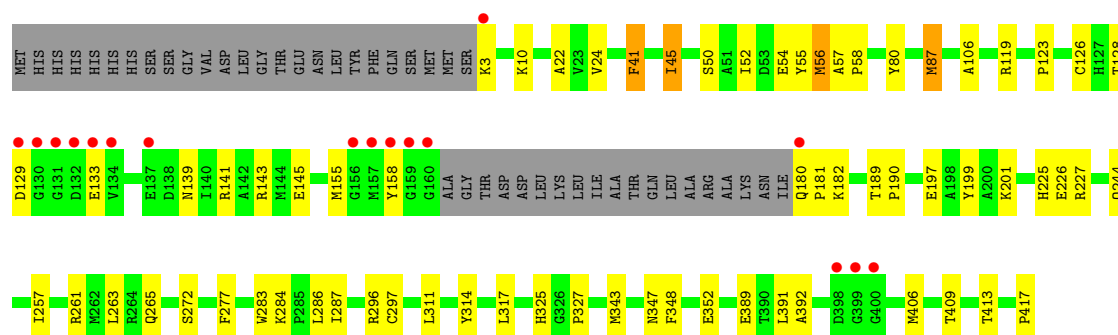


- Molecule 1: Mandelate racemase/muconate lactonizing protein

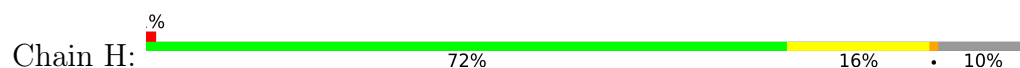




• Molecule 1: Mandelate racemase/muconate lactonizing protein



• Molecule 1: Mandelate racemase/muconate lactonizing protein



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	224.33Å 224.35Å 87.08Å 90.00° 97.96° 90.00°	Depositor
Resolution (Å)	40.00 – 2.00 40.00 – 2.00	Depositor EDS
% Data completeness (in resolution range)	88.0 (40.00-2.00) 88.0 (40.00-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.19	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.26 (at 1.94Å)	Xtriage
Refinement program	REFMAC 5.7.0025	Depositor
R, R_{free}	0.173 , 0.228 (Not available) , 0.228	Depositor DCC
R_{free} test set	8445 reflections (2.69%)	wwPDB-VP
Wilson B-factor (Å ²)	15.6	Xtriage
Anisotropy	0.448	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 40.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	26763	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.72% 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, CL, IOD, FMT, MG

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	0.72	1/3162 (0.0%)	0.91	1/4296 (0.0%)
1	B	0.73	1/3170 (0.0%)	0.95	4/4305 (0.1%)
1	C	0.70	1/3156 (0.0%)	0.97	5/4289 (0.1%)
1	D	0.74	2/3158 (0.1%)	0.95	2/4291 (0.0%)
1	E	0.77	2/3162 (0.1%)	0.98	2/4295 (0.0%)
1	F	0.72	1/3167 (0.0%)	0.95	2/4303 (0.0%)
1	G	0.74	2/3160 (0.1%)	0.96	3/4295 (0.1%)
1	H	0.72	0/3153	0.97	6/4283 (0.1%)
All	All	0.73	10/25288 (0.0%)	0.96	25/34357 (0.1%)

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	195	ASP	C-O	-8.38	1.16	1.24
1	A	70	ASP	C-O	-6.80	1.16	1.24
1	G	197	GLU	C-O	-6.04	1.17	1.24
1	D	88	ASN	C-O	-5.80	1.17	1.24
1	D	53	ASP	C-O	-5.52	1.17	1.24
1	B	261	ARG	C-O	-5.46	1.17	1.24
1	E	197	GLU	C-O	-5.44	1.17	1.24
1	F	208	GLU	C-O	-5.31	1.18	1.24
1	G	201	LYS	C-O	-5.24	1.17	1.24
1	C	235	GLN	C-O	-5.01	1.18	1.24

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	330	ILE	N-CA-C	-7.26	98.77	108.06
1	B	195	ASP	CA-C-N	6.98	126.95	119.28
1	B	195	ASP	C-N-CA	6.98	126.95	119.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	122	ILE	N-CA-C	6.71	114.19	107.55
1	F	100	ASP	N-CA-C	-6.13	104.52	111.14
1	E	330	ILE	N-CA-C	-6.06	100.35	108.35
1	G	227	ARG	N-CA-C	5.92	120.34	113.18
1	G	389	GLU	N-CA-C	5.61	117.07	111.07
1	H	326	GLY	CA-C-N	5.50	126.71	119.84
1	H	326	GLY	C-N-CA	5.50	126.71	119.84
1	B	122	ILE	N-CA-C	5.40	113.38	107.60
1	G	297	CYS	N-CA-C	5.39	117.24	110.24
1	C	122	ILE	N-CA-C	5.25	113.22	107.60
1	C	326	GLY	CA-C-N	5.19	126.32	119.84
1	C	326	GLY	C-N-CA	5.19	126.32	119.84
1	D	195	ASP	CA-C-N	5.16	124.78	119.05
1	D	195	ASP	C-N-CA	5.16	124.78	119.05
1	C	306	THR	CA-C-N	-5.14	113.74	119.19
1	C	306	THR	C-N-CA	-5.14	113.74	119.19
1	H	189	THR	CA-C-N	5.12	125.06	119.78
1	H	189	THR	C-N-CA	5.12	125.06	119.78
1	E	63	LYS	N-CA-C	5.10	117.44	110.55
1	F	122	ILE	N-CA-C	5.08	112.58	107.55
1	B	299	VAL	N-CA-C	5.03	115.64	110.36
1	H	40	THR	N-CA-C	5.02	117.46	108.17

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	3084	0	3036	53	0
1	B	3086	0	3046	66	0
1	C	3072	0	3023	68	0
1	D	3077	0	3026	59	0
1	E	3081	0	3037	63	0
1	F	3083	0	3032	43	0
1	G	3079	0	3023	53	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	H	3075	0	3029	61	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	12	0	16	3	0
3	B	12	0	16	1	0
3	C	12	0	16	1	0
3	D	12	0	16	2	0
3	E	6	0	8	0	0
3	F	12	0	16	2	0
3	G	6	0	8	0	0
3	H	12	0	16	1	0
4	A	3	0	1	1	0
4	B	3	0	1	3	0
4	C	3	0	1	0	0
4	D	3	0	1	0	0
4	E	3	0	1	0	0
4	F	6	0	2	1	0
4	G	3	0	1	3	0
4	H	3	0	1	1	0
5	A	2	0	0	2	0
5	B	3	0	0	1	0
5	C	3	0	0	4	0
5	D	3	0	0	2	0
5	E	3	0	0	3	0
5	F	3	0	0	4	0
5	H	2	0	0	1	0
6	B	1	0	0	0	0
6	C	1	0	0	0	0
6	D	1	0	0	0	0
6	E	2	0	0	0	0
6	F	1	0	0	0	0
6	G	2	0	0	0	0
6	H	1	0	0	0	0
7	A	244	0	0	6	0
7	B	255	0	0	12	0
7	C	237	0	0	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	D	237	0	0	10	0
7	E	276	0	0	13	0
7	F	239	0	0	8	0
7	G	234	0	0	8	0
7	H	257	0	0	13	0
All	All	26763	0	24373	448	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 9.

All (448) 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:122:ILE:HD11	1:E:343:MET:CE	1.83	1.08
7:E:874:HOH:O	1:F:257:ILE:HD11	1.63	0.98
1:E:122:ILE:HD11	1:E:343:MET:HE2	1.46	0.97
1:C:280:ILE:HD13	1:C:280:ILE:O	1.67	0.94
1:H:416:ARG:HD2	7:H:846:HOH:O	1.66	0.94
1:D:201:LYS:HE2	7:F:828:HOH:O	1.69	0.92
1:E:126:CYS:HB2	7:E:744:HOH:O	1.69	0.92
1:E:122:ILE:HD11	1:E:343:MET:HE1	1.56	0.88
1:E:122:ILE:CD1	1:E:343:MET:HE1	2.07	0.85
1:A:8:ASN:HA	1:A:61:ILE:HD11	1.61	0.83
1:H:57:ALA:HB3	1:H:58:PRO:HD3	1.61	0.83
1:A:145:GLU:HA	1:A:145:GLU:OE1	1.77	0.82
1:C:187:SER:HB2	7:E:852:HOH:O	1.80	0.82
1:B:325:HIS:CD2	4:B:505:FMT:H	2.16	0.80
1:A:182:LYS:HE2	1:D:150:TYR:CZ	2.17	0.79
1:G:352:GLU:OE2	4:G:505:FMT:H	1.83	0.79
7:A:752:HOH:O	5:C:506:IOD:I	2.71	0.79
1:C:280:ILE:HD11	1:C:284:LYS:HD3	1.65	0.78
1:H:388:ASN:CG	1:H:391:LEU:HD12	2.09	0.77
1:D:137:GLU:HB3	1:D:141:ARG:NH2	2.00	0.76
1:G:181:PRO:HB2	7:G:765:HOH:O	1.84	0.76
1:A:13:LEU:HD11	1:A:49:LYS:HD3	1.68	0.75
1:F:182:LYS:HE3	7:G:766:HOH:O	1.86	0.75
1:G:56:MET:HE3	1:G:56:MET:HA	1.70	0.74
1:H:196:PRO:HD2	7:H:805:HOH:O	1.86	0.74
1:B:325:HIS:NE2	4:B:505:FMT:H	2.03	0.74
7:A:620:HOH:O	1:C:284:LYS:HE3	1.89	0.73
7:A:794:HOH:O	1:B:401:ILE:HD13	1.88	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:56:MET:HE3	1:H:56:MET:HA	1.70	0.72
1:F:57:ALA:O	1:F:61:ILE:HG12	1.89	0.71
1:G:24:VAL:CG2	1:G:52:ILE:HD13	2.21	0.71
1:E:229:THR:HG23	5:E:508:IOD:I	2.59	0.71
1:C:280:ILE:HD13	1:C:280:ILE:C	2.16	0.70
1:E:232:THR:HG23	7:E:864:HOH:O	1.92	0.70
1:C:4:LEU:HD13	1:C:30:GLU:HG3	1.74	0.70
1:H:145:GLU:HG2	7:H:699:HOH:O	1.91	0.70
1:H:196:PRO:HB2	5:H:507:IOD:I	2.61	0.69
1:C:280:ILE:CD1	1:C:284:LYS:HB2	2.23	0.69
1:G:391:LEU:HD11	7:G:825:HOH:O	1.92	0.69
1:F:56:MET:HE3	1:F:56:MET:HA	1.74	0.69
1:E:57:ALA:O	1:E:61:ILE:HG12	1.92	0.69
1:D:391:LEU:HD12	1:D:394:LYS:HE2	1.75	0.68
1:B:152:ARG:HD2	7:B:710:HOH:O	1.94	0.67
1:A:56:MET:HE3	1:A:56:MET:HA	1.77	0.67
1:A:57:ALA:HB3	1:A:58:PRO:HD3	1.77	0.67
1:F:196:PRO:HB2	5:F:509:IOD:I	2.66	0.66
1:B:56:MET:HE3	1:B:56:MET:HA	1.78	0.66
1:A:32:GLY:H	3:A:505:GOL:H2	1.60	0.66
1:A:122:ILE:HD11	1:A:343:MET:HG2	1.77	0.65
1:C:280:ILE:HD11	1:C:284:LYS:HB2	1.77	0.65
1:A:234:ILE:HG12	1:A:263:LEU:HB2	1.80	0.64
1:B:117:LYS:HG2	1:B:343[A]:MET:SD	2.37	0.64
1:A:8:ASN:HD22	1:A:9:VAL:H	1.44	0.64
1:C:184:SER:HB3	1:C:185:PRO:HD2	1.80	0.63
1:D:391:LEU:CD1	1:D:394:LYS:HE2	2.29	0.63
1:H:263:LEU:C	1:H:263:LEU:HD23	2.23	0.63
1:E:278:VAL:HG12	1:E:301:THR:HG22	1.79	0.63
1:E:122:ILE:CD1	1:E:343:MET:CE	2.65	0.62
1:D:137:GLU:HB3	1:D:141:ARG:HH21	1.64	0.62
1:A:196:PRO:HB2	5:A:506:IOD:I	2.70	0.62
1:F:278:VAL:HG12	1:F:301:THR:HG22	1.82	0.62
1:F:251:PRO:HD2	1:F:273:MET:SD	2.40	0.62
1:H:145:GLU:CG	7:H:699:HOH:O	2.48	0.62
1:A:258:ASP:HB2	7:A:773:HOH:O	1.98	0.61
1:A:7:THR:C	1:A:61:ILE:HD12	2.25	0.61
1:B:119:ARG:O	1:B:343[B]:MET:HE1	2.01	0.61
1:A:7:THR:O	1:A:61:ILE:HD12	2.00	0.61
1:C:391:LEU:HD13	1:C:394:LYS:HD3	1.82	0.60
1:B:343[B]:MET:HE3	1:B:343[B]:MET:HA	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:388:ASN:CB	1:H:391:LEU:HD12	2.31	0.60
1:A:182:LYS:HE2	1:D:150:TYR:OH	2.00	0.60
1:A:225:HIS:O	1:A:226:GLU:HG2	2.01	0.60
1:A:145:GLU:OE1	1:A:145:GLU:CA	2.46	0.60
1:D:401:ILE:HD12	7:H:849:HOH:O	2.02	0.60
1:C:284:LYS:N	1:C:285:PRO:HD2	2.17	0.59
7:A:801:HOH:O	1:G:182:LYS:HE3	2.02	0.59
1:B:152:ARG:NH2	7:B:613:HOH:O	2.34	0.59
1:G:119:ARG:O	1:G:343[A]:MET:HE1	2.03	0.58
1:D:190:PRO:HA	7:D:738:HOH:O	2.03	0.58
1:G:24:VAL:HG22	1:G:52:ILE:HD13	1.83	0.58
1:C:391:LEU:CD1	1:C:394:LYS:HD3	2.34	0.58
1:D:56:MET:HE3	1:D:56:MET:HA	1.84	0.58
1:C:82:ARG:HA	1:C:87:MET:CE	2.32	0.58
1:D:132:ASP:O	1:D:136:VAL:HG23	2.03	0.58
1:B:137:GLU:CD	1:B:213:LYS:HE2	2.28	0.58
1:A:7:THR:C	1:A:61:ILE:CD1	2.77	0.58
1:B:261:ARG:HD2	7:B:769:HOH:O	2.04	0.57
1:F:235:GLN:CD	7:F:756:HOH:O	2.47	0.57
1:E:389:GLU:O	1:E:393:GLU:HG3	2.05	0.57
1:F:152:ARG:HD3	7:F:822:HOH:O	2.05	0.57
1:E:122:ILE:CG1	1:E:343:MET:HE1	2.35	0.57
1:D:327:PRO:O	1:D:330:ILE:HG22	2.04	0.57
1:E:152:ARG:HD3	7:E:667:HOH:O	2.05	0.56
1:F:406:MET:O	1:F:406:MET:HG3	2.03	0.56
1:G:257:ILE:HB	1:G:286:LEU:HD21	1.86	0.56
1:C:4:LEU:HD13	1:C:30:GLU:CG	2.35	0.56
1:C:60:LEU:HD11	1:C:94:VAL:HG21	1.88	0.56
1:D:152:ARG:HD3	7:D:798:HOH:O	2.04	0.56
1:A:123:PRO:HD2	1:A:348:PHE:CE2	2.40	0.56
1:C:391:LEU:HD12	1:C:391:LEU:O	2.05	0.56
1:E:261:ARG:O	1:E:265[B]:GLN:HG3	2.05	0.56
1:C:401:ILE:HD13	1:C:401:ILE:N	2.21	0.56
1:G:343[A]:MET:HE3	1:G:343[A]:MET:HA	1.88	0.55
1:C:278:VAL:HG12	1:C:301:THR:HG22	1.88	0.55
1:E:181:PRO:HB2	7:E:771:HOH:O	2.07	0.55
1:C:122:ILE:HD11	1:C:343:MET:HG2	1.88	0.55
1:B:196:PRO:HB2	5:B:508:IOD:I	2.77	0.55
1:D:274:GLY:HA2	1:D:277:PHE:CD2	2.42	0.55
1:A:133:GLU:OE1	1:A:155:MET:SD	2.65	0.55
1:B:225:HIS:ND1	1:B:417:PRO:OXT	2.39	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:82:ARG:HA	1:C:87:MET:HE3	1.89	0.55
1:D:68:ILE:HD12	1:D:102:LYS:HE3	1.89	0.55
1:D:284:LYS:HB3	1:D:285:PRO:HD3	1.89	0.54
1:A:398:ASP:HB2	7:A:756:HOH:O	2.07	0.54
1:E:196:PRO:HB2	5:E:508:IOD:I	2.76	0.54
1:G:158:TYR:HD1	1:G:199:TYR:CE1	2.26	0.54
1:H:133:GLU:C	1:H:133:GLU:OE1	2.50	0.54
1:A:136:VAL:O	1:A:140:ILE:HG13	2.07	0.54
1:C:201:LYS:HE2	7:E:861:HOH:O	2.07	0.54
1:E:295:ILE:HG12	1:E:311:LEU:HD21	1.89	0.54
1:B:208:GLU:HG3	1:B:243:TYR:OH	2.08	0.54
1:A:60:LEU:HD11	1:A:94:VAL:HG21	1.90	0.54
1:D:231:VAL:HG12	7:D:696:HOH:O	2.08	0.54
1:D:241:GLU:N	1:D:242:PRO:CD	2.70	0.54
1:D:20:ASP:HB3	1:D:45:ILE:HD13	1.89	0.54
1:E:57:ALA:HB3	1:E:58:PRO:HD3	1.90	0.54
1:F:15:ALA:N	1:F:16:PRO:HD3	2.23	0.54
1:G:141:ARG:O	1:G:145:GLU:HG2	2.08	0.53
5:D:505:IOD:I	1:H:84:GLY:HA3	2.79	0.53
1:G:284:LYS:NZ	7:G:688:HOH:O	2.40	0.53
1:B:278:VAL:HG12	1:B:301:THR:HG22	1.90	0.53
1:C:196:PRO:HB2	5:C:508:IOD:I	2.78	0.53
1:H:365:PRO:O	1:H:385:ILE:HD12	2.08	0.53
1:E:280:ILE:HG22	7:E:831:HOH:O	2.09	0.53
1:H:13:LEU:HD11	1:H:49:LYS:HD2	1.90	0.53
1:C:301:THR:HG22	1:C:301:THR:O	2.09	0.53
1:A:275:GLU:OE2	1:B:80:TYR:OH	2.25	0.53
1:C:68:ILE:HD11	1:C:102:LYS:HE2	1.90	0.53
1:B:15:ALA:N	1:B:16:PRO:HD3	2.24	0.53
1:B:325:HIS:NE2	4:B:505:FMT:C	2.70	0.53
1:E:8:ASN:OD1	1:E:27:GLU:OE1	2.27	0.53
1:E:317:LEU:HD13	1:G:283:TRP:HH2	1.73	0.52
1:C:257:ILE:HB	1:C:286:LEU:HD21	1.90	0.52
1:C:280:ILE:C	1:C:280:ILE:CD1	2.82	0.52
1:B:122:ILE:HD13	1:B:339:MET:HG2	1.91	0.52
1:G:55:TYR:C	1:G:58:PRO:HD2	2.35	0.52
1:A:409:THR:HB	1:A:410:PRO:CD	2.40	0.52
1:C:325:HIS:CE1	1:C:327:PRO:HG3	2.45	0.52
1:E:327:PRO:O	1:E:330:ILE:HG22	2.09	0.52
1:F:317:LEU:HD13	1:H:283:TRP:HH2	1.75	0.52
5:F:507:IOD:I	1:H:116:GLY:HA2	2.80	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:502:GOL:H12	1:B:80:TYR:CD2	2.45	0.51
1:B:155:MET:HE2	7:B:671:HOH:O	2.10	0.51
1:E:106:ALA:HB1	1:G:106:ALA:HB1	1.91	0.51
1:F:229:THR:HG23	5:F:509:IOD:I	2.81	0.51
1:C:123:PRO:HG2	1:C:348:PHE:HE2	1.76	0.51
1:H:260:LEU:HD12	1:H:286:LEU:HD21	1.91	0.51
1:C:184:SER:HB3	1:C:185:PRO:CD	2.40	0.51
1:G:406:MET:O	1:G:406:MET:HG3	2.11	0.51
1:H:388:ASN:CG	1:H:391:LEU:CD1	2.80	0.51
1:A:125:TYR:CZ	1:A:352:GLU:HB2	2.45	0.51
1:C:280:ILE:CD1	1:C:284:LYS:HD3	2.39	0.51
1:E:133:GLU:HG3	1:E:155:MET:SD	2.50	0.51
1:A:265:GLN:HE22	1:D:265[B]:GLN:HG2	1.75	0.51
1:A:3:LYS:O	1:A:3:LYS:HG3	2.09	0.51
1:E:224:VAL:HG23	1:E:248:LEU:HD11	1.92	0.51
1:H:20:ASP:HB3	1:H:45:ILE:HD13	1.92	0.51
1:A:241:GLU:N	1:A:242:PRO:CD	2.73	0.51
1:G:284:LYS:HE2	1:G:314:TYR:OH	2.11	0.51
1:D:224:VAL:HG23	1:D:248:LEU:HD11	1.93	0.51
1:A:409:THR:HB	1:A:410:PRO:HD2	1.92	0.50
1:D:355:PRO:HD2	7:D:757:HOH:O	2.11	0.50
1:F:57:ALA:HB3	1:F:58:PRO:HD3	1.94	0.50
1:A:32:GLY:N	3:A:505:GOL:H2	2.26	0.50
1:E:15:ALA:N	1:E:16:PRO:HD3	2.26	0.50
1:C:158:TYR:O	1:C:194:PHE:HB2	2.12	0.50
1:D:140:ILE:O	1:D:144:MET:HG3	2.11	0.50
1:D:343:MET:HE3	7:D:780:HOH:O	2.12	0.50
1:A:296:ARG:HD2	1:A:323:ALA:O	2.11	0.50
1:H:309:LYS:NZ	7:H:731:HOH:O	2.45	0.50
1:B:201:LYS:HE2	1:C:216:PHE:O	2.12	0.50
1:C:152:ARG:HD3	7:C:687:HOH:O	2.12	0.50
1:E:217:GLY:N	7:E:773:HOH:O	2.44	0.50
1:F:362[A]:GLU:HG3	7:F:675:HOH:O	2.11	0.50
1:F:290:LYS:NZ	7:F:677:HOH:O	2.45	0.50
1:H:123:PRO:HG2	1:H:348:PHE:HE2	1.77	0.50
1:C:280:ILE:HD11	1:C:284:LYS:CD	2.37	0.49
1:G:123:PRO:HG2	1:G:348:PHE:HE2	1.76	0.49
1:H:388:ASN:HB3	1:H:391:LEU:HD12	1.94	0.49
1:D:68:ILE:CD1	1:D:102:LYS:HE3	2.41	0.49
1:H:283:TRP:NE1	7:H:757:HOH:O	2.38	0.49
1:A:304:GLY:O	1:A:307:PRO:HD2	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:295:ILE:HG12	1:C:311:LEU:HD21	1.95	0.49
1:D:30:GLU:OE1	1:D:30:GLU:HA	2.12	0.49
1:D:241:GLU:N	1:D:242:PRO:HD2	2.27	0.49
1:D:207:PHE:CZ	1:D:222:HIS:HB2	2.48	0.49
1:F:283:TRP:HH2	1:H:317:LEU:HD13	1.77	0.49
1:C:15:ALA:N	1:C:16:PRO:HD3	2.28	0.49
1:B:22:ALA:HB2	1:B:45:ILE:HD11	1.95	0.49
1:E:259:TRP:CE2	1:H:290:LYS:HD2	2.47	0.49
1:G:343[A]:MET:HE3	1:G:343[A]:MET:CA	2.42	0.49
1:C:116:GLY:HA2	5:C:506:IOD:I	2.83	0.49
1:B:343[B]:MET:HE3	1:B:343[B]:MET:CA	2.43	0.48
1:C:123:PRO:HG2	1:C:348:PHE:CE2	2.48	0.48
1:E:241:GLU:N	1:E:242:PRO:CD	2.76	0.48
1:G:133:GLU:HG3	1:G:155:MET:SD	2.53	0.48
1:D:137:GLU:HG3	1:D:210:LEU:CD2	2.44	0.48
1:H:384:GLY:HA2	7:H:736:HOH:O	2.12	0.48
1:A:278:VAL:HG12	1:A:301:THR:HG22	1.95	0.48
1:E:224:VAL:CG2	1:E:248:LEU:HD11	2.44	0.48
1:D:390:THR:HB	7:D:828:HOH:O	2.13	0.48
1:E:14:THR:C	1:E:16:PRO:HD3	2.38	0.48
1:E:122:ILE:HG13	1:E:343:MET:HE1	1.94	0.48
1:E:257:ILE:HD11	1:F:281:ASN:OD1	2.14	0.48
1:F:250:ASP:HA	1:F:273:MET:SD	2.53	0.48
1:B:137:GLU:OE2	1:B:213:LYS:HE2	2.12	0.48
1:G:277:PHE:CD2	1:G:283:TRP:HB3	2.49	0.48
1:E:257:ILE:HB	1:E:286:LEU:HD21	1.95	0.48
1:G:263:LEU:HD23	1:G:263:LEU:C	2.38	0.48
1:A:304:GLY:C	1:A:307:PRO:HD2	2.39	0.48
1:C:68:ILE:CD1	1:C:102:LYS:HE2	2.44	0.48
1:G:158:TYR:CD1	1:G:199:TYR:CE1	3.02	0.48
1:B:217:GLY:N	7:B:821:HOH:O	2.38	0.48
1:C:280:ILE:HD13	1:C:284:LYS:HB2	1.92	0.48
1:E:350:ILE:HG13	1:E:351:GLN:N	2.29	0.48
1:G:189:THR:CG2	1:G:190:PRO:HD2	2.43	0.48
1:B:241:GLU:N	1:B:242:PRO:HD2	2.29	0.47
1:B:201:LYS:HE2	7:C:764:HOH:O	2.14	0.47
1:E:376:TYR:HA	7:E:823:HOH:O	2.14	0.47
1:D:132:ASP:HA	1:D:155:MET:SD	2.54	0.47
1:E:60:LEU:HD11	1:E:94:VAL:HG21	1.96	0.47
1:B:234:ILE:HG12	1:B:263:LEU:HB2	1.95	0.47
1:D:8:ASN:ND2	1:D:9:VAL:H	2.12	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:13:LEU:HB3	1:B:397:CYS:SG	2.54	0.47
1:E:294:TYR:HA	1:E:321:ARG:O	2.15	0.47
1:C:119:ARG:O	1:C:343:MET:HE1	2.15	0.47
1:D:34:TYR:HB3	3:D:503:GOL:H31	1.97	0.47
1:E:254:PRO:HD3	7:E:637:HOH:O	2.15	0.47
1:F:207:PHE:CZ	1:F:222:HIS:HB2	2.50	0.47
1:G:272:SER:HB2	1:G:296:ARG:CG	2.45	0.47
1:H:128:THR:CG2	1:H:143:ARG:HG3	2.45	0.47
1:A:332:PRO:HG2	1:A:364:PHE:CD2	2.49	0.47
1:B:350:ILE:HG13	1:B:351:GLN:N	2.29	0.47
1:D:402:PRO:HD3	7:D:660:HOH:O	2.14	0.47
1:H:284:LYS:HB3	1:H:285:PRO:HD3	1.97	0.47
1:A:263:LEU:C	1:A:263:LEU:HD23	2.40	0.47
1:E:42:THR:HA	1:E:45:ILE:HG23	1.95	0.47
1:G:283:TRP:CE3	1:G:287:ILE:HD12	2.50	0.47
1:C:4:LEU:CD1	1:C:30:GLU:CG	2.93	0.46
1:H:128:THR:HG22	1:H:143:ARG:HG3	1.97	0.46
1:D:196:PRO:HB2	5:D:508:IOD:I	2.84	0.46
1:D:261:ARG:O	1:D:265[A]:GLN:HG3	2.15	0.46
1:E:55:TYR:O	1:F:401:ILE:HD11	2.14	0.46
1:E:261:ARG:O	1:E:265[A]:GLN:HG3	2.14	0.46
1:A:296:ARG:NH2	4:A:503:FMT:H	2.31	0.46
1:C:401:ILE:HD13	1:C:401:ILE:H	1.79	0.46
1:F:32:GLY:H	3:F:508:GOL:H2	1.79	0.46
1:B:158:TYR:O	1:B:194:PHE:HB2	2.15	0.46
1:D:263:LEU:HD23	1:D:263:LEU:C	2.40	0.46
1:E:56:MET:CG	1:E:94:VAL:HG21	2.44	0.46
1:A:265:GLN:HE22	1:D:265[A]:GLN:HG2	1.80	0.46
1:G:189:THR:HG23	1:G:190:PRO:HD2	1.97	0.46
1:E:80:TYR:CZ	1:F:276:LEU:HD21	2.50	0.46
1:F:389:GLU:HG3	7:F:779:HOH:O	2.14	0.46
1:F:254:PRO:O	1:F:257:ILE:HG23	2.16	0.46
1:F:261:ARG:O	1:F:265[B]:GLN:HG3	2.16	0.46
1:H:193:TYR:HA	1:H:408:ARG:O	2.15	0.46
1:H:260:LEU:HD12	1:H:286:LEU:CD2	2.46	0.46
1:B:226:GLU:OE1	1:B:417:PRO:O	2.34	0.46
1:B:260:LEU:HD12	1:B:286:LEU:CD2	2.46	0.46
1:B:355:PRO:HD2	7:B:712:HOH:O	2.16	0.46
1:C:225:HIS:CD2	7:C:724:HOH:O	2.68	0.46
1:D:123:PRO:HG2	1:D:348:PHE:HE2	1.80	0.46
1:A:8:ASN:ND2	1:A:9:VAL:H	2.13	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:339:MET:HA	1:B:342:ASP:HB2	1.97	0.45
1:E:193:TYR:HA	1:E:408:ARG:O	2.16	0.45
1:D:261:ARG:O	1:D:265[B]:GLN:HG3	2.14	0.45
1:G:409:THR:OG1	1:G:413:THR:HB	2.15	0.45
1:A:8:ASN:CA	1:A:61:ILE:HD11	2.37	0.45
1:G:126:CYS:SG	1:G:143:ARG:HD3	2.57	0.45
1:G:352:GLU:OE2	4:G:505:FMT:C	2.58	0.45
1:B:152:ARG:NH1	7:B:748:HOH:O	2.49	0.45
1:B:223:ASP:HA	1:B:249:GLU:HB3	1.99	0.45
1:H:57:ALA:CB	1:H:58:PRO:HD3	2.41	0.45
1:H:409:THR:HB	1:H:410:PRO:CD	2.47	0.45
1:D:72:TRP:CZ2	1:D:307:PRO:HG3	2.52	0.45
1:E:225:HIS:O	1:E:226:GLU:HG3	2.17	0.45
1:G:139:ASN:O	1:G:143:ARG:HG2	2.16	0.45
1:G:352:GLU:CD	4:G:505:FMT:H	2.41	0.45
1:H:278:VAL:HG12	1:H:301:THR:HG22	1.98	0.45
1:C:225:HIS:O	1:C:226:GLU:HG2	2.15	0.45
1:A:8:ASN:HD22	1:A:9:VAL:N	2.13	0.45
1:C:4:LEU:CD1	1:C:30:GLU:HG2	2.46	0.45
1:E:10:LYS:HB2	1:E:10:LYS:HE3	1.70	0.45
1:F:56:MET:HE3	1:F:56:MET:CA	2.45	0.45
1:H:157:MET:HE2	7:H:694:HOH:O	2.15	0.45
1:B:82:ARG:O	1:B:83:ASN:HB2	2.16	0.45
1:B:280:ILE:HG12	1:B:280:ILE:O	2.16	0.45
1:E:226:GLU:HA	7:E:838:HOH:O	2.17	0.45
1:F:350:ILE:HG13	1:F:351:GLN:N	2.32	0.45
1:H:122:ILE:HD13	1:H:339:MET:HG2	1.99	0.45
1:A:229:THR:HG23	5:A:506:IOD:I	2.87	0.45
7:B:822:HOH:O	1:H:201:LYS:HE2	2.16	0.45
1:C:225:HIS:O	1:C:226:GLU:CG	2.64	0.45
1:D:224:VAL:CG2	1:D:248:LEU:HD11	2.46	0.45
1:F:261:ARG:O	1:F:265[A]:GLN:HG3	2.17	0.45
1:H:158:TYR:CD1	1:H:227:ARG:HG3	2.51	0.45
1:H:226:GLU:OE1	1:H:417:PRO:O	2.35	0.45
1:A:13:LEU:HD11	1:A:49:LYS:CD	2.42	0.44
1:B:189:THR:HG23	1:B:190:PRO:HD2	1.98	0.44
1:C:409:THR:HB	1:C:410:PRO:HD2	1.99	0.44
1:G:87:MET:HE2	7:G:603:HOH:O	2.16	0.44
1:B:261:ARG:CZ	1:B:261:ARG:HB3	2.47	0.44
1:D:257:ILE:HB	1:D:286:LEU:HD21	2.00	0.44
1:E:196:PRO:HD2	7:E:862:HOH:O	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:258:ASP:HB3	7:G:802:HOH:O	2.17	0.44
1:B:371:ASP:HB3	1:B:376:TYR:HE2	1.82	0.44
1:C:128:THR:HG22	1:C:143:ARG:HG3	1.98	0.44
1:C:280:ILE:HG12	1:C:283:TRP:CZ2	2.52	0.44
1:D:82:ARG:HD2	3:H:504:GOL:H31	1.99	0.44
1:B:115:GLY:HA2	7:B:798:HOH:O	2.17	0.44
1:B:225:HIS:O	1:B:226:GLU:HG2	2.17	0.44
1:C:385:ILE:HG13	1:C:386:ASP:N	2.32	0.44
1:D:273:MET:HE3	1:D:273:MET:HB3	1.88	0.44
1:B:263:LEU:HD23	1:B:263:LEU:C	2.43	0.44
1:H:119:ARG:O	1:H:343:MET:HE1	2.18	0.44
1:C:296:ARG:HD2	1:C:323:ALA:O	2.18	0.44
1:F:362[B]:GLU:HG2	1:F:391:LEU:CD1	2.47	0.44
1:G:22:ALA:HB2	1:G:45:ILE:HG22	1.99	0.44
1:H:273:MET:HE3	1:H:273:MET:HB3	1.87	0.44
1:A:213:LYS:HA	1:A:213:LYS:HD3	1.89	0.44
1:B:260:LEU:HD12	1:B:286:LEU:HD21	1.98	0.44
1:C:64:ASP:OD1	1:C:64:ASP:C	2.61	0.44
1:D:80:TYR:CZ	1:H:276:LEU:HD21	2.52	0.44
1:D:217:GLY:N	7:D:762:HOH:O	2.47	0.44
1:H:280:ILE:HG22	7:H:718:HOH:O	2.17	0.44
1:G:50:SER:O	1:G:54:GLU:HB2	2.18	0.44
1:B:296:ARG:HD2	1:B:323:ALA:O	2.18	0.43
1:B:194:PHE:O	1:B:410:PRO:HD3	2.18	0.43
1:B:261:ARG:CZ	1:B:261:ARG:CB	2.95	0.43
1:G:244:GLN:HA	7:G:801:HOH:O	2.17	0.43
1:B:56:MET:HE3	1:B:56:MET:CA	2.46	0.43
1:F:388:ASN:OD1	1:F:388:ASN:C	2.61	0.43
1:G:128:THR:HG22	1:G:143:ARG:HG3	2.00	0.43
1:H:258:ASP:OD1	1:H:261:ARG:HD3	2.18	0.43
1:C:24:VAL:HG22	1:C:52:ILE:HD13	1.99	0.43
1:D:409:THR:HB	1:D:410:PRO:CD	2.48	0.43
1:H:60:LEU:HD11	1:H:94:VAL:HG11	2.00	0.43
1:E:84:GLY:HA3	5:F:505:IOD:I	2.89	0.43
1:G:347:ASN:O	1:G:348:PHE:C	2.62	0.43
1:H:15:ALA:N	1:H:16:PRO:HD3	2.33	0.43
1:H:57:ALA:HB3	1:H:58:PRO:CD	2.42	0.43
1:C:294:TYR:HA	1:C:321:ARG:O	2.18	0.43
1:D:362:GLU:O	1:D:365:PRO:HD3	2.19	0.43
1:E:41:PHE:CD1	1:E:41:PHE:C	2.95	0.43
1:E:158:TYR:O	1:E:194:PHE:HB2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:261:ARG:O	1:G:265:GLN:HG3	2.19	0.43
1:B:261:ARG:HB3	1:B:261:ARG:NH2	2.34	0.43
1:F:409:THR:HG23	1:F:415:SER:OG	2.19	0.43
1:H:184:SER:HB3	1:H:185:PRO:HD2	2.01	0.43
1:H:409:THR:HB	1:H:410:PRO:HD2	2.01	0.43
1:B:306:THR:HB	1:B:307:PRO:CD	2.49	0.43
1:C:57:ALA:O	1:C:61:ILE:HG12	2.19	0.43
1:C:241:GLU:N	1:C:242:PRO:CD	2.82	0.43
1:D:343:MET:CE	7:D:780:HOH:O	2.67	0.43
1:G:225:HIS:O	1:G:226:GLU:HG3	2.19	0.43
1:B:83:ASN:OD1	3:B:504:GOL:H11	2.19	0.42
1:B:256:ASN:ND2	7:B:639:HOH:O	2.30	0.42
1:C:57:ALA:HB3	1:C:58:PRO:HD3	2.01	0.42
1:E:185:PRO:HG3	1:H:374:TYR:CE1	2.54	0.42
1:E:241:GLU:N	1:E:242:PRO:HD2	2.34	0.42
1:G:317:LEU:HD23	1:G:317:LEU:HA	1.83	0.42
1:B:14:THR:C	1:B:16:PRO:HD3	2.45	0.42
1:B:203:VAL:HB	1:B:204:PRO:HD3	2.01	0.42
1:E:41:PHE:HB2	1:E:89:ASN:ND2	2.35	0.42
1:E:116:GLY:HA2	5:E:506:IOD:I	2.89	0.42
1:F:241:GLU:N	1:F:242:PRO:CD	2.82	0.42
1:D:309:LYS:HB3	7:D:827:HOH:O	2.18	0.42
1:E:296:ARG:HD2	1:E:323:ALA:O	2.20	0.42
1:A:102:LYS:HE3	1:A:102:LYS:HB2	1.91	0.42
1:E:409:THR:HG23	1:E:415:SER:OG	2.19	0.42
1:D:79:GLY:O	1:D:80:TYR:HB3	2.20	0.42
1:E:56:MET:HG2	1:E:94:VAL:HG21	2.02	0.42
1:G:41:PHE:CD1	1:G:41:PHE:C	2.97	0.42
1:G:225:HIS:ND1	1:G:417:PRO:OXT	2.51	0.42
1:C:280:ILE:CD1	1:C:284:LYS:CB	2.93	0.42
1:H:60:LEU:HD11	1:H:94:VAL:HG21	2.02	0.42
1:H:155:MET:HE2	7:H:830:HOH:O	2.20	0.42
1:G:55:TYR:O	1:G:58:PRO:HD2	2.19	0.41
1:G:325:HIS:CE1	1:G:327:PRO:HG3	2.55	0.41
1:C:8:ASN:ND2	1:C:27:GLU:OE1	2.45	0.41
1:C:291:LEU:HD23	1:C:291:LEU:HA	1.85	0.41
1:G:391:LEU:O	1:G:392:ALA:C	2.62	0.41
4:F:506:FMT:H	7:F:800:HOH:O	2.18	0.41
1:G:283:TRP:CE2	1:G:311:LEU:HD13	2.55	0.41
1:B:195:ASP:HA	1:B:196:PRO:HD3	1.84	0.41
1:D:27:GLU:HB3	3:D:503:GOL:H32	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:87:MET:CE	7:G:603:HOH:O	2.69	0.41
1:H:56:MET:HE3	1:H:56:MET:CA	2.43	0.41
1:B:41:PHE:C	1:B:41:PHE:CD1	2.98	0.41
1:B:115:GLY:CA	7:B:798:HOH:O	2.69	0.41
1:C:94:VAL:O	1:C:98:LEU:HG	2.21	0.41
1:E:80:TYR:CE1	1:F:276:LEU:HD21	2.54	0.41
1:A:265:GLN:NE2	1:D:265[B]:GLN:HG2	2.35	0.41
1:B:402:PRO:HB2	1:B:404:TRP:CD1	2.56	0.41
1:F:265[B]:GLN:NE2	1:G:265:GLN:HA	2.35	0.41
1:H:376:TYR:HA	7:H:754:HOH:O	2.20	0.41
1:A:119:ARG:O	1:A:343:MET:HE1	2.20	0.41
1:B:262:MET:HB2	7:B:763:HOH:O	2.19	0.41
1:D:409:THR:HB	1:D:410:PRO:HD2	2.01	0.41
1:F:139:ASN:O	1:F:143:ARG:HG2	2.21	0.41
1:F:330:ILE:O	1:F:330:ILE:HG23	2.21	0.41
1:H:13:LEU:CD1	1:H:49:LYS:HD2	2.50	0.41
3:C:502:GOL:H12	1:G:80:TYR:CD2	2.56	0.41
1:D:137:GLU:HG3	1:D:210:LEU:HD23	2.03	0.41
1:E:80:TYR:CD2	3:F:502:GOL:H12	2.56	0.41
1:G:57:ALA:HB3	1:G:58:PRO:HD3	2.02	0.41
1:B:327:PRO:O	1:B:330:ILE:HG22	2.21	0.41
1:C:56:MET:CG	1:C:94:VAL:HG21	2.50	0.41
1:C:122:ILE:HA	1:C:123:PRO:HD2	1.90	0.41
1:D:57:ALA:O	1:D:61:ILE:HG12	2.20	0.41
1:E:227:ARG:HD3	1:E:227:ARG:HA	1.94	0.41
1:F:123:PRO:HG2	1:F:348:PHE:HE2	1.86	0.41
1:F:236:LEU:O	1:F:237:ALA:C	2.63	0.41
1:A:274:GLY:HA2	1:A:277:PHE:CD2	2.57	0.41
1:C:123:PRO:CG	1:C:348:PHE:HE2	2.34	0.41
1:D:274:GLY:HA2	1:D:277:PHE:CE2	2.56	0.41
1:H:57:ALA:O	1:H:61:ILE:HG12	2.21	0.41
1:C:122:ILE:HD13	1:C:339:MET:HG2	2.03	0.40
1:F:225:HIS:O	1:F:226:GLU:CG	2.69	0.40
1:F:394:LYS:HB2	1:F:394:LYS:HE3	1.81	0.40
1:A:80:TYR:CZ	1:B:276:LEU:HD21	2.56	0.40
1:A:280:ILE:HD13	7:C:613:HOH:O	2.20	0.40
1:A:343:MET:HE3	1:A:343:MET:CA	2.52	0.40
1:H:262:MET:HA	1:H:265:GLN:HE21	1.85	0.40
1:H:296:ARG:HD2	1:H:323:ALA:O	2.21	0.40
1:H:296:ARG:NH2	4:H:505:FMT:H	2.36	0.40
1:B:60:LEU:HD11	1:B:94:VAL:HG11	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:87:MET:HG2	5:C:505:IOD:I	2.91	0.40
1:F:280:ILE:HD13	7:H:673:HOH:O	2.21	0.40
1:A:306:THR:HB	1:A:307:PRO:HD3	2.04	0.40
1:C:195:ASP:HA	1:C:196:PRO:HD2	1.95	0.40
1:E:401:ILE:HD13	7:F:821:HOH:O	2.20	0.40
1:G:272:SER:HB2	1:G:296:ARG:HG2	2.02	0.40
1:B:270:PRO:HA	1:B:293:ASP:OD2	2.21	0.40
1:D:128:THR:HG22	1:D:143:ARG:HG3	2.04	0.40
1:H:92:SER:OG	1:H:96:MET:HE3	2.21	0.40
1:H:158:TYR:OH	1:H:225:HIS:HB2	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	392/440 (89%)	384 (98%)	8 (2%)	0	100	100
1	B	393/440 (89%)	379 (96%)	14 (4%)	0	100	100
1	C	393/440 (89%)	382 (97%)	10 (2%)	1 (0%)	36	35
1	D	392/440 (89%)	380 (97%)	12 (3%)	0	100	100
1	E	392/440 (89%)	380 (97%)	12 (3%)	0	100	100
1	F	393/440 (89%)	378 (96%)	15 (4%)	0	100	100
1	G	393/440 (89%)	383 (98%)	10 (2%)	0	100	100
1	H	391/440 (89%)	381 (97%)	10 (3%)	0	100	100
All	All	3139/3520 (89%)	3047 (97%)	91 (3%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	327	PRO

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	331/369 (90%)	323 (98%)	8 (2%)	43	47
1	B	332/369 (90%)	326 (98%)	6 (2%)	51	58
1	C	329/369 (89%)	320 (97%)	9 (3%)	39	42
1	D	330/369 (89%)	324 (98%)	6 (2%)	51	58
1	E	331/369 (90%)	325 (98%)	6 (2%)	51	58
1	F	331/369 (90%)	323 (98%)	8 (2%)	43	47
1	G	330/369 (89%)	322 (98%)	8 (2%)	43	47
1	H	330/369 (89%)	322 (98%)	8 (2%)	43	47
All	All	2644/2952 (90%)	2585 (98%)	59 (2%)	45	50

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LYS
1	A	45	ILE
1	A	54	GLU
1	A	56	MET
1	A	133	GLU
1	A	145	GLU
1	A	258	ASP
1	A	398	ASP
1	B	41	PHE
1	B	133	GLU
1	B	137	GLU
1	B	157	MET
1	B	208	GLU
1	B	213	LYS
1	C	45	ILE

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Mol	Chain	Res	Type
1	C	133	GLU
1	C	134	VAL
1	C	208	GLU
1	C	280	ILE
1	C	372	GLN
1	C	385	ILE
1	C	391	LEU
1	C	401	ILE
1	D	45	ILE
1	D	56	MET
1	D	212	ASN
1	D	213	LYS
1	D	390	THR
1	D	398	ASP
1	E	30	GLU
1	E	45	ILE
1	E	87	MET
1	E	134	VAL
1	E	213	LYS
1	E	267	SER
1	F	45	ILE
1	F	56	MET
1	F	87	MET
1	F	102	LYS
1	F	129	ASP
1	F	133	GLU
1	F	134	VAL
1	F	208	GLU
1	G	3	LYS
1	G	10	LYS
1	G	41	PHE
1	G	45	ILE
1	G	56	MET
1	G	87	MET
1	G	129	ASP
1	G	180	GLN
1	H	3	LYS
1	H	19	ILE
1	H	56	MET
1	H	117	LYS
1	H	133	GLU
1	H	145	GLU

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Mol	Chain	Res	Type
1	H	157	MET
1	H	205	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	ASN
1	A	256	ASN
1	A	265	GLN
1	B	8	ASN
1	B	212	ASN
1	C	209	HIS
1	C	256	ASN
1	C	265	GLN
1	D	8	ASN
1	D	244	GLN
1	D	256	ASN
1	E	8	ASN
1	E	212	ASN
1	E	256	ASN
1	F	8	ASN
1	F	372	GLN
1	G	8	ASN
1	G	89	ASN
1	G	180	GLN
1	G	212	ASN
1	H	256	ASN
1	H	265	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 59 ligands modelled in this entry, 36 are monoatomic - leaving 23 for Mogul analysis.

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
3	GOL	E	502	-	5,5,5	0.36	0	5,5,5	0.77	0
4	FMT	G	505	2	2,2,2	0.75	0	1,1,1	0.44	0
3	GOL	H	502	-	5,5,5	0.39	0	5,5,5	0.58	0
3	GOL	D	502	-	5,5,5	0.54	0	5,5,5	0.43	0
3	GOL	B	504	-	5,5,5	0.39	0	5,5,5	0.44	0
3	GOL	F	508	-	5,5,5	0.18	0	5,5,5	0.61	0
3	GOL	C	502	-	5,5,5	0.41	0	5,5,5	0.36	0
4	FMT	F	503	2	2,2,2	0.39	0	1,1,1	0.47	0
4	FMT	D	506	2	2,2,2	0.67	0	1,1,1	0.61	0
3	GOL	A	505	-	5,5,5	0.26	0	5,5,5	0.53	0
4	FMT	B	505	2	2,2,2	0.81	0	1,1,1	0.43	0
4	FMT	C	503	2	2,2,2	0.68	0	1,1,1	0.61	0
3	GOL	G	502	-	5,5,5	0.36	0	5,5,5	0.19	0
3	GOL	A	502	-	5,5,5	0.34	0	5,5,5	0.36	0
4	FMT	F	506	-	2,2,2	0.93	0	1,1,1	0.41	0
3	GOL	B	502	-	5,5,5	0.55	0	5,5,5	0.43	0
3	GOL	F	502	-	5,5,5	0.50	0	5,5,5	0.45	0
4	FMT	E	505	2	2,2,2	0.46	0	1,1,1	0.47	0
3	GOL	C	507	-	5,5,5	0.30	0	5,5,5	0.57	0
4	FMT	A	503	2	2,2,2	0.36	0	1,1,1	0.63	0
4	FMT	H	505	2	2,2,2	0.62	0	1,1,1	0.62	0
3	GOL	H	504	-	5,5,5	0.27	0	5,5,5	0.56	0
3	GOL	D	503	-	5,5,5	0.20	0	5,5,5	0.59	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
3	GOL	E	502	-	-	2/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	B	504	-	-	2/4/4/4	-
3	GOL	F	508	-	-	3/4/4/4	-
3	GOL	G	502	-	-	3/4/4/4	-
3	GOL	A	502	-	-	4/4/4/4	-
3	GOL	C	502	-	-	2/4/4/4	-
3	GOL	B	502	-	-	2/4/4/4	-
3	GOL	H	504	-	-	2/4/4/4	-
3	GOL	A	505	-	-	0/4/4/4	-
3	GOL	F	502	-	-	4/4/4/4	-
3	GOL	D	502	-	-	0/4/4/4	-
3	GOL	D	503	-	-	2/4/4/4	-
3	GOL	H	502	-	-	2/4/4/4	-
3	GOL	C	507	-	-	1/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (29) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	502	GOL	O1-C1-C2-C3
3	B	504	GOL	O1-C1-C2-C3
3	C	502	GOL	O1-C1-C2-C3
3	F	502	GOL	O1-C1-C2-C3
3	H	502	GOL	O1-C1-C2-C3
3	H	504	GOL	C1-C2-C3-O3
3	H	504	GOL	O2-C2-C3-O3
3	F	508	GOL	O2-C2-C3-O3
3	H	502	GOL	O1-C1-C2-O2
3	A	502	GOL	C1-C2-C3-O3
3	B	502	GOL	C1-C2-C3-O3
3	D	503	GOL	C1-C2-C3-O3
3	E	502	GOL	O1-C1-C2-C3
3	F	502	GOL	C1-C2-C3-O3
3	F	508	GOL	C1-C2-C3-O3
3	G	502	GOL	C1-C2-C3-O3
3	B	504	GOL	O1-C1-C2-O2
3	D	503	GOL	O2-C2-C3-O3
3	E	502	GOL	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
3	A	502	GOL	O1-C1-C2-O2
3	C	502	GOL	O1-C1-C2-O2
3	F	502	GOL	O1-C1-C2-O2
3	B	502	GOL	O2-C2-C3-O3
3	F	502	GOL	O2-C2-C3-O3
3	G	502	GOL	O2-C2-C3-O3
3	G	502	GOL	O1-C1-C2-O2
3	F	508	GOL	O1-C1-C2-C3
3	A	502	GOL	O2-C2-C3-O3
3	C	507	GOL	O1-C1-C2-O2

There are no ring outliers.

13 monomers are involved in 19 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	G	505	FMT	3	0
3	B	504	GOL	1	0
3	F	508	GOL	1	0
3	C	502	GOL	1	0
3	A	505	GOL	2	0
4	B	505	FMT	3	0
3	A	502	GOL	1	0
4	F	506	FMT	1	0
3	F	502	GOL	1	0
4	A	503	FMT	1	0
4	H	505	FMT	1	0
3	H	504	GOL	1	0
3	D	503	GOL	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	396/440 (90%)	-0.36	1 (0%) 90 89	8, 17, 40, 69	0
1	B	395/440 (89%)	-0.32	14 (3%) 47 46	7, 16, 45, 89	2 (0%)
1	C	395/440 (89%)	-0.32	13 (3%) 49 48	7, 16, 40, 75	2 (0%)
1	D	395/440 (89%)	-0.33	9 (2%) 61 60	7, 16, 41, 65	1 (0%)
1	E	395/440 (89%)	-0.39	6 (1%) 72 71	6, 15, 39, 77	1 (0%)
1	F	395/440 (89%)	-0.34	9 (2%) 61 60	7, 17, 43, 85	2 (0%)
1	G	396/440 (90%)	-0.30	17 (4%) 40 39	7, 17, 47, 100	1 (0%)
1	H	395/440 (89%)	-0.42	6 (1%) 72 71	6, 15, 39, 70	0
All	All	3162/3520 (89%)	-0.35	75 (2%) 59 59	6, 16, 42, 100	9 (0%)

All (75) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	160	GLY	6.1
1	B	157	MET	5.9
1	F	400	GLY	4.9
1	D	181	PRO	4.8
1	C	157	MET	4.5
1	D	157	MET	4.5
1	G	131	GLY	4.5
1	E	157	MET	4.5
1	G	130	GLY	4.5
1	G	157	MET	4.2
1	E	160	GLY	4.1
1	H	160	GLY	3.9
1	F	157	MET	3.8
1	G	180	GLN	3.8
1	B	160	GLY	3.6
1	G	156	GLY	3.6

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Mol	Chain	Res	Type	RSRZ
1	H	157	MET	3.5
1	B	398	ASP	3.4
1	C	134	VAL	3.4
1	B	399	GLY	3.4
1	B	400	GLY	3.3
1	C	158	TYR	3.2
1	B	3	LYS	3.2
1	G	398	ASP	3.2
1	D	400	GLY	3.2
1	G	134	VAL	3.1
1	G	399	GLY	3.1
1	G	400	GLY	3.1
1	C	160	GLY	3.1
1	D	399	GLY	3.0
1	B	134	VAL	2.9
1	C	133	GLU	2.8
1	G	158	TYR	2.8
1	B	129	ASP	2.8
1	G	137	GLU	2.8
1	G	133	GLU	2.8
1	E	400	GLY	2.8
1	E	398	ASP	2.7
1	C	399	GLY	2.7
1	C	401	ILE	2.7
1	C	129	ASP	2.7
1	F	398	ASP	2.6
1	G	129	ASP	2.6
1	H	400	GLY	2.6
1	F	134	VAL	2.5
1	H	181	PRO	2.5
1	G	3	LYS	2.5
1	B	133	GLU	2.5
1	D	134	VAL	2.4
1	B	197	GLU	2.4
1	F	156	GLY	2.4
1	D	158	TYR	2.4
1	F	132	ASP	2.3
1	F	209	HIS	2.3
1	F	129	ASP	2.3
1	H	398	ASP	2.3
1	B	181	PRO	2.3
1	G	159	GLY	2.3

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Mol	Chain	Res	Type	RSRZ
1	E	3	LYS	2.2
1	B	159	GLY	2.2
1	C	209	HIS	2.2
1	B	401	ILE	2.2
1	B	158	TYR	2.2
1	C	181	PRO	2.2
1	E	181	PRO	2.2
1	C	187	SER	2.2
1	D	129	ASP	2.2
1	G	132	ASP	2.2
1	D	133	GLU	2.1
1	C	400	GLY	2.1
1	A	3	LYS	2.1
1	F	158	TYR	2.1
1	C	398	ASP	2.1
1	H	209	HIS	2.0
1	D	160	GLY	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
3	GOL	A	505	6/6	0.88	0.14	31,43,56,56	0
3	GOL	B	502	6/6	0.89	0.14	17,22,37,39	0
3	GOL	B	504	6/6	0.91	0.12	24,40,40,46	0
3	GOL	F	508	6/6	0.91	0.12	26,29,55,56	0
4	FMT	E	505	3/3	0.91	0.12	23,23,29,36	0
4	FMT	F	503	3/3	0.91	0.10	26,26,31,31	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	FMT	F	506	3/3	0.91	0.12	30,30,31,46	0
3	GOL	D	503	6/6	0.92	0.10	33,34,41,47	0
3	GOL	E	502	6/6	0.92	0.11	19,23,25,46	0
3	GOL	D	502	6/6	0.92	0.11	12,22,26,28	0
4	FMT	A	503	3/3	0.93	0.10	30,30,32,42	0
3	GOL	F	502	6/6	0.93	0.11	17,23,29,30	0
3	GOL	C	502	6/6	0.93	0.12	17,22,31,40	0
3	GOL	H	504	6/6	0.93	0.10	20,31,39,41	0
3	GOL	H	502	6/6	0.94	0.09	9,19,25,34	0
3	GOL	A	502	6/6	0.94	0.10	13,22,29,44	0
5	IOD	A	504	1/1	0.94	0.07	38,38,38,38	1
5	IOD	E	506	1/1	0.94	0.06	34,34,34,34	1
6	CL	C	504	1/1	0.94	0.07	38,38,38,38	0
4	FMT	G	505	3/3	0.95	0.11	19,19,28,37	0
3	GOL	G	502	6/6	0.95	0.09	16,19,27,38	0
3	GOL	C	507	6/6	0.95	0.10	26,27,33,43	0
4	FMT	D	506	3/3	0.95	0.11	28,28,43,53	0
6	CL	D	504	1/1	0.95	0.07	27,27,27,27	0
5	IOD	D	507	1/1	0.96	0.06	33,33,33,33	1
4	FMT	B	505	3/3	0.96	0.11	28,28,29,49	0
5	IOD	E	507	1/1	0.96	0.05	31,31,31,31	1
6	CL	B	503	1/1	0.96	0.07	33,33,33,33	0
5	IOD	B	507	1/1	0.96	0.06	32,32,32,32	1
5	IOD	C	506	1/1	0.96	0.06	33,33,33,33	1
6	CL	E	503	1/1	0.96	0.06	33,33,33,33	0
6	CL	F	504	1/1	0.96	0.06	34,34,34,34	0
6	CL	G	503	1/1	0.96	0.09	30,30,30,30	0
6	CL	G	504	1/1	0.96	0.06	35,35,35,35	0
4	FMT	H	505	3/3	0.97	0.06	17,17,25,29	0
5	IOD	F	507	1/1	0.97	0.05	35,35,35,35	1
6	CL	E	504	1/1	0.97	0.06	35,35,35,35	0
5	IOD	H	506	1/1	0.97	0.05	30,30,30,30	1
5	IOD	D	508	1/1	0.97	0.09	38,38,38,38	1
4	FMT	C	503	3/3	0.97	0.06	26,26,27,35	0
6	CL	H	503	1/1	0.97	0.06	37,37,37,37	0
2	MG	B	501	1/1	0.98	0.06	14,14,14,14	0
5	IOD	B	508	1/1	0.98	0.06	33,33,33,33	1
2	MG	E	501	1/1	0.98	0.03	17,17,17,17	0
5	IOD	E	508	1/1	0.98	0.08	32,32,32,32	1
2	MG	H	501	1/1	0.98	0.03	14,14,14,14	0
5	IOD	H	507	1/1	0.99	0.06	30,30,30,30	1
5	IOD	D	505	1/1	0.99	0.02	14,14,14,14	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	MG	C	501	1/1	0.99	0.02	16,16,16,16	0
5	IOD	A	506	1/1	0.99	0.07	31,31,31,31	1
5	IOD	B	506	1/1	0.99	0.02	15,15,15,15	0
2	MG	D	501	1/1	0.99	0.03	14,14,14,14	0
2	MG	A	501	1/1	0.99	0.02	11,11,11,11	0
2	MG	F	501	1/1	0.99	0.02	20,20,20,20	0
5	IOD	F	509	1/1	0.99	0.08	35,35,35,35	1
5	IOD	C	508	1/1	0.99	0.07	33,33,33,33	1
5	IOD	C	505	1/1	1.00	0.02	15,15,15,15	0
5	IOD	F	505	1/1	1.00	0.01	15,15,15,15	0
2	MG	G	501	1/1	1.00	0.02	15,15,15,15	0

6.5 Other polymers [i](#)

There are no such residues in this entry.