



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:07 PM UTC

PDB ID : 4HN8 / pdb_00004hn8
Title : Crystal structure of a putative D-glucarate dehydratase from *Pseudomonas mendocina* ymp
Authors : Hegde, R.P.; Toro, R.; Burley, S.K.; Almo, S.C.; Ramagopal, U.A.; New York Structural Genomics Research Consortium (NYSGRG)
Deposited on : 2012-10-19
Resolution : 2.20 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

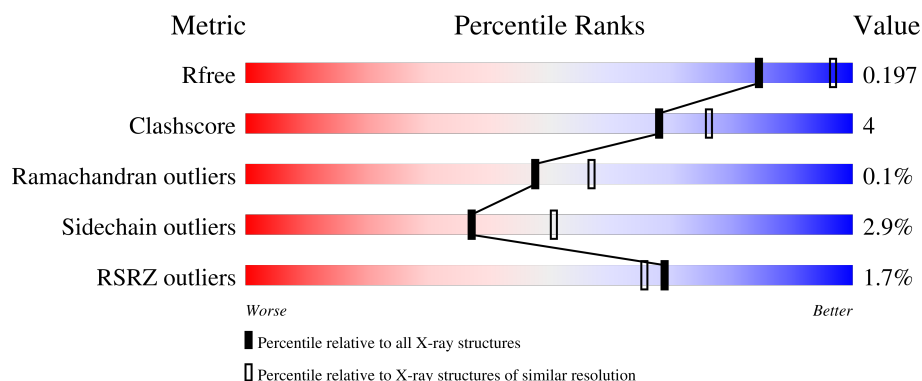
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	473	0.1% 83% 7% 10%
1	B	473	0.1% 82% 7% 10%
1	C	473	2% 82% 8% 10%
1	D	473	2% 83% 7% 10%
1	E	473	0.1% 81% 8% 10%

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Mol	Chain	Length	Quality of chain
1	F	473	 % 83% 7% 10%
1	G	473	 % 80% 10% 10%
1	H	473	 2% 82% 7% 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
2	GOL	G	501	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 27164 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 D-glucarate dehydratase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	427	Total	C	N	O	S	Se	0	4	0
			3328	2089	617	602	6	14			
1	B	426	Total	C	N	O	S	Se	0	1	0
			3301	2071	611	599	6	14			
1	C	427	Total	C	N	O	S	Se	0	7	0
			3348	2102	620	606	6	14			
1	D	427	Total	C	N	O	S	Se	0	6	0
			3333	2095	612	606	6	14			
1	E	426	Total	C	N	O	S	Se	0	3	0
			3312	2079	614	599	6	14			
1	F	426	Total	C	N	O	S	Se	0	3	0
			3319	2082	619	598	6	14			
1	G	426	Total	C	N	O	S	Se	0	3	0
			3314	2079	615	600	6	14			
1	H	427	Total	C	N	O	S	Se	0	2	0
			3313	2081	612	600	6	14			

There are 88 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MSE	-	expression tag	UNP A4XRL3
A	0	SER	-	expression tag	UNP A4XRL3
A	1	LEU	-	expression tag	UNP A4XRL3
A	464	GLU	-	expression tag	UNP A4XRL3
A	465	GLY	-	expression tag	UNP A4XRL3
A	466	HIS	-	expression tag	UNP A4XRL3
A	467	HIS	-	expression tag	UNP A4XRL3
A	468	HIS	-	expression tag	UNP A4XRL3
A	469	HIS	-	expression tag	UNP A4XRL3
A	470	HIS	-	expression tag	UNP A4XRL3
A	471	HIS	-	expression tag	UNP A4XRL3
B	-1	MSE	-	expression tag	UNP A4XRL3
B	0	SER	-	expression tag	UNP A4XRL3

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Chain	Residue	Modelled	Actual	Comment	Reference
B	1	LEU	-	expression tag	UNP A4XRL3
B	464	GLU	-	expression tag	UNP A4XRL3
B	465	GLY	-	expression tag	UNP A4XRL3
B	466	HIS	-	expression tag	UNP A4XRL3
B	467	HIS	-	expression tag	UNP A4XRL3
B	468	HIS	-	expression tag	UNP A4XRL3
B	469	HIS	-	expression tag	UNP A4XRL3
B	470	HIS	-	expression tag	UNP A4XRL3
B	471	HIS	-	expression tag	UNP A4XRL3
C	-1	MSE	-	expression tag	UNP A4XRL3
C	0	SER	-	expression tag	UNP A4XRL3
C	1	LEU	-	expression tag	UNP A4XRL3
C	464	GLU	-	expression tag	UNP A4XRL3
C	465	GLY	-	expression tag	UNP A4XRL3
C	466	HIS	-	expression tag	UNP A4XRL3
C	467	HIS	-	expression tag	UNP A4XRL3
C	468	HIS	-	expression tag	UNP A4XRL3
C	469	HIS	-	expression tag	UNP A4XRL3
C	470	HIS	-	expression tag	UNP A4XRL3
C	471	HIS	-	expression tag	UNP A4XRL3
D	-1	MSE	-	expression tag	UNP A4XRL3
D	0	SER	-	expression tag	UNP A4XRL3
D	1	LEU	-	expression tag	UNP A4XRL3
D	464	GLU	-	expression tag	UNP A4XRL3
D	465	GLY	-	expression tag	UNP A4XRL3
D	466	HIS	-	expression tag	UNP A4XRL3
D	467	HIS	-	expression tag	UNP A4XRL3
D	468	HIS	-	expression tag	UNP A4XRL3
D	469	HIS	-	expression tag	UNP A4XRL3
D	470	HIS	-	expression tag	UNP A4XRL3
D	471	HIS	-	expression tag	UNP A4XRL3
E	-1	MSE	-	expression tag	UNP A4XRL3
E	0	SER	-	expression tag	UNP A4XRL3
E	1	LEU	-	expression tag	UNP A4XRL3
E	464	GLU	-	expression tag	UNP A4XRL3
E	465	GLY	-	expression tag	UNP A4XRL3
E	466	HIS	-	expression tag	UNP A4XRL3
E	467	HIS	-	expression tag	UNP A4XRL3
E	468	HIS	-	expression tag	UNP A4XRL3
E	469	HIS	-	expression tag	UNP A4XRL3
E	470	HIS	-	expression tag	UNP A4XRL3
E	471	HIS	-	expression tag	UNP A4XRL3

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Chain	Residue	Modelled	Actual	Comment	Reference
F	-1	MSE	-	expression tag	UNP A4XRL3
F	0	SER	-	expression tag	UNP A4XRL3
F	1	LEU	-	expression tag	UNP A4XRL3
F	464	GLU	-	expression tag	UNP A4XRL3
F	465	GLY	-	expression tag	UNP A4XRL3
F	466	HIS	-	expression tag	UNP A4XRL3
F	467	HIS	-	expression tag	UNP A4XRL3
F	468	HIS	-	expression tag	UNP A4XRL3
F	469	HIS	-	expression tag	UNP A4XRL3
F	470	HIS	-	expression tag	UNP A4XRL3
F	471	HIS	-	expression tag	UNP A4XRL3
G	-1	MSE	-	expression tag	UNP A4XRL3
G	0	SER	-	expression tag	UNP A4XRL3
G	1	LEU	-	expression tag	UNP A4XRL3
G	464	GLU	-	expression tag	UNP A4XRL3
G	465	GLY	-	expression tag	UNP A4XRL3
G	466	HIS	-	expression tag	UNP A4XRL3
G	467	HIS	-	expression tag	UNP A4XRL3
G	468	HIS	-	expression tag	UNP A4XRL3
G	469	HIS	-	expression tag	UNP A4XRL3
G	470	HIS	-	expression tag	UNP A4XRL3
G	471	HIS	-	expression tag	UNP A4XRL3
H	-1	MSE	-	expression tag	UNP A4XRL3
H	0	SER	-	expression tag	UNP A4XRL3
H	1	LEU	-	expression tag	UNP A4XRL3
H	464	GLU	-	expression tag	UNP A4XRL3
H	465	GLY	-	expression tag	UNP A4XRL3
H	466	HIS	-	expression tag	UNP A4XRL3
H	467	HIS	-	expression tag	UNP A4XRL3
H	468	HIS	-	expression tag	UNP A4XRL3
H	469	HIS	-	expression tag	UNP A4XRL3
H	470	HIS	-	expression tag	UNP A4XRL3
H	471	HIS	-	expression tag	UNP A4XRL3

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



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

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	103	Total	O	0	0
			103	103		
3	B	84	Total	O	0	0
			84	84		
3	C	71	Total	O	0	0
			71	71		

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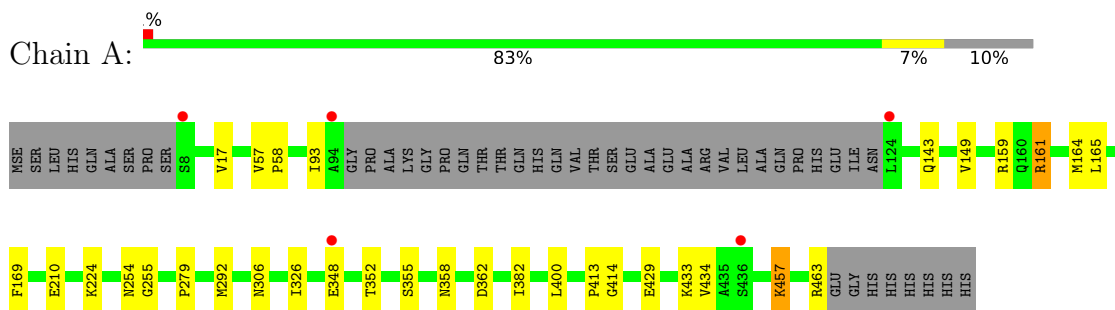
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	64	Total 64	O 64	0	0
3	E	61	Total 61	O 61	0	0
3	F	32	Total 32	O 32	0	0
3	G	53	Total 53	O 53	0	0
3	H	74	Total 74	O 74	0	0

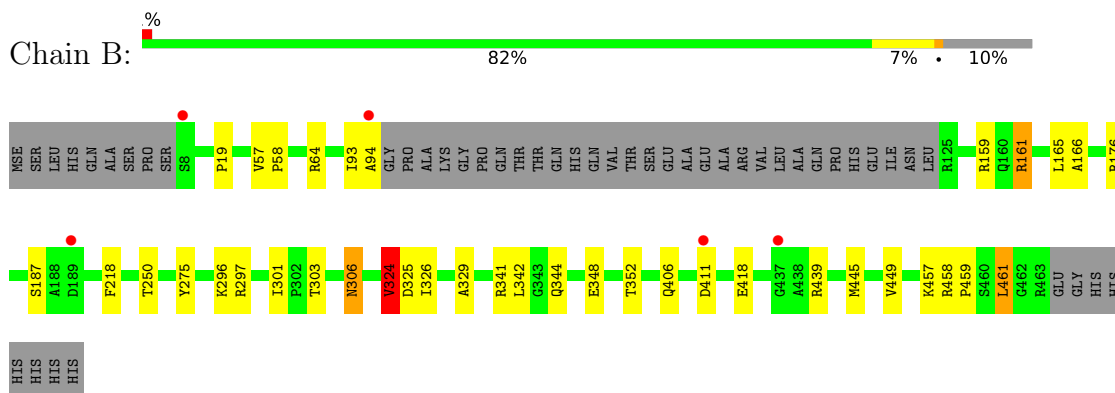
3 Residue-property plots [i](#)

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

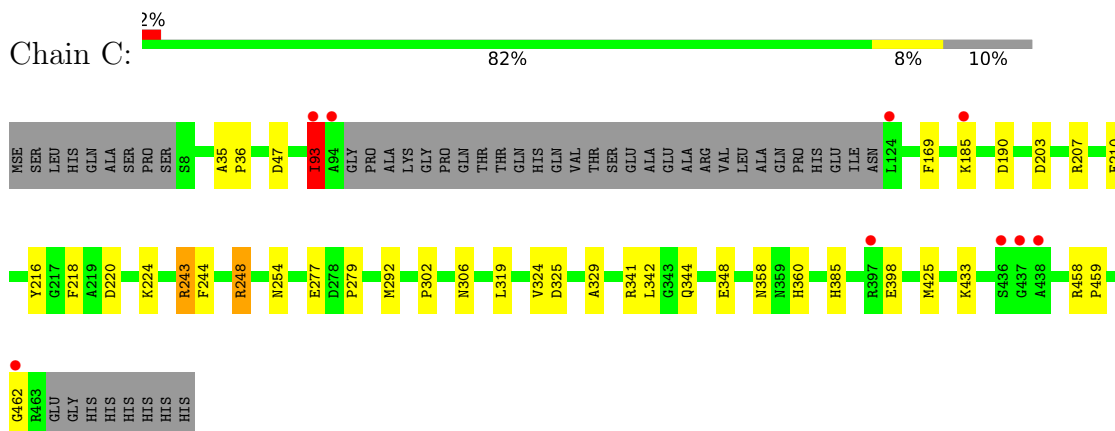
- Molecule 1: D-glucarate dehydratase



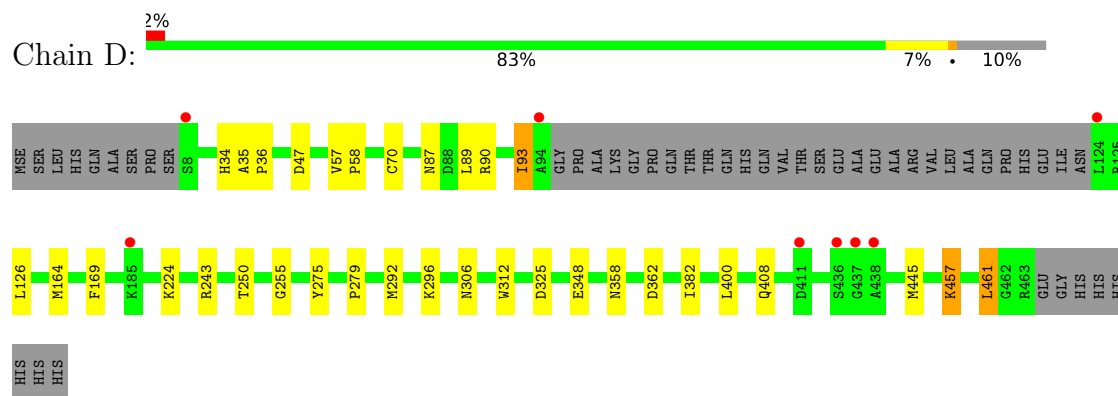
- Molecule 1: D-glucarate dehydratase



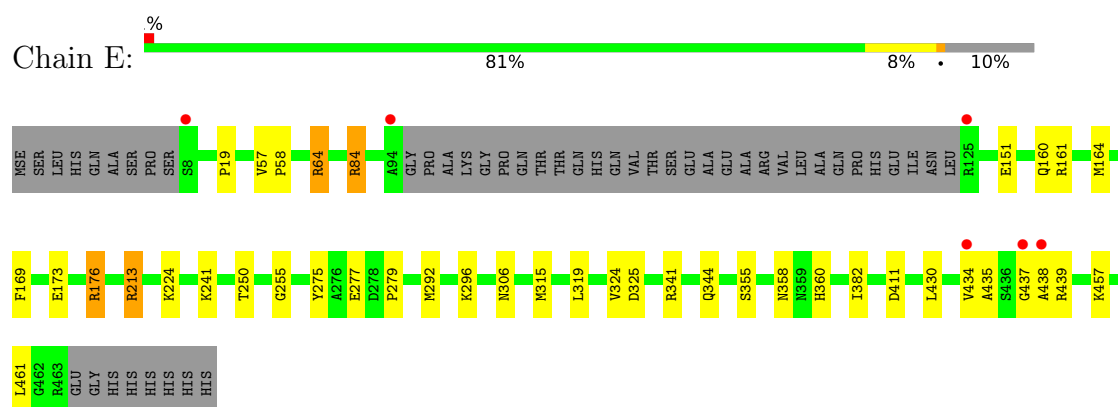
- Molecule 1: D-glucarate dehydratase



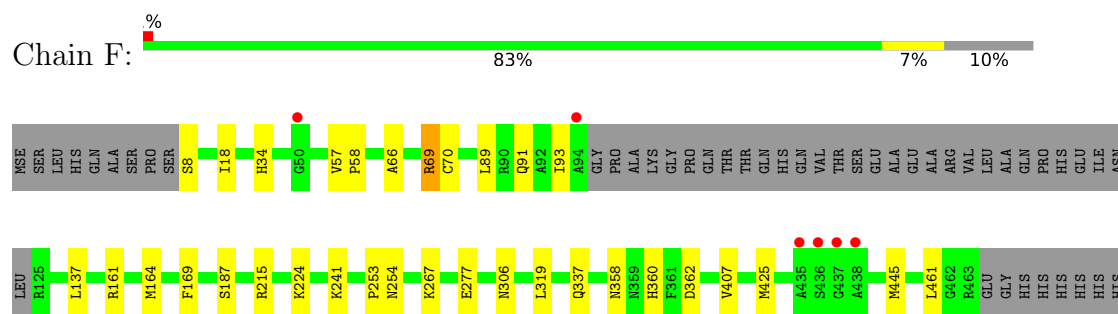
- Molecule 1: D-glucarate dehydratase



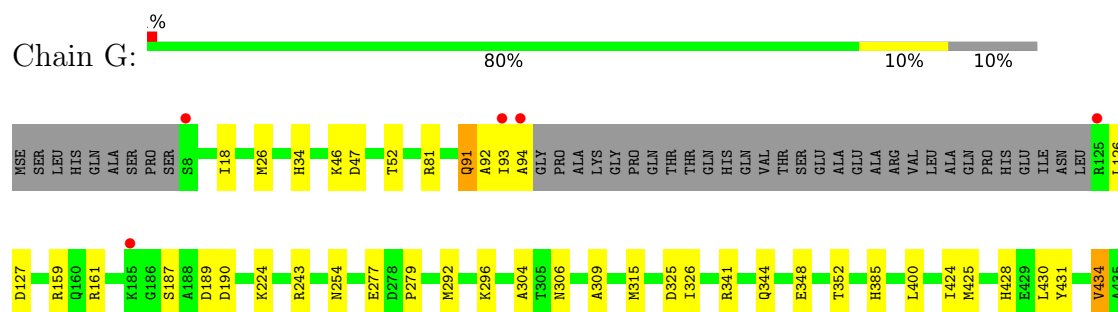
- Molecule 1: D-glucarate dehydratase

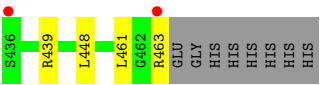


- Molecule 1: D-glucarate dehydratase

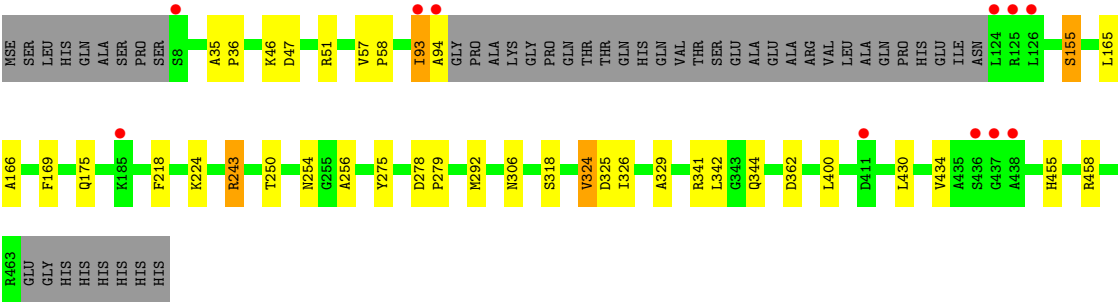
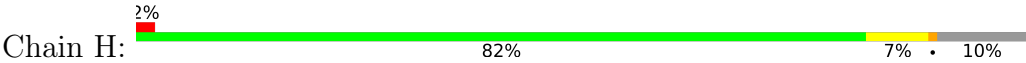


- Molecule 1: D-glucarate dehydratase





● Molecule 1: D-glucarate dehydratase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	126.36Å 148.84Å 198.46Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.18 – 2.20 46.18 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.7 (46.18-2.20) 99.7 (46.18-2.20)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.68 (at 2.20Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.176 , 0.222 (Not available) , 0.197	Depositor DCC
R_{free} test set	9478 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	24.4	Xtriage
Anisotropy	0.016	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 24.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	27164	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 16.18% 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 [i](#)

5.1 Standard geometry [i](#)

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	0.81	0/3394	0.86	1/4574 (0.0%)
1	B	0.81	0/3361	0.88	2/4530 (0.0%)
1	C	0.77	0/3423	0.87	5/4613 (0.1%)
1	D	0.77	0/3409	0.88	0/4597
1	E	0.75	0/3379	0.86	0/4555
1	F	0.74	0/3386	0.87	1/4563 (0.0%)
1	G	0.75	0/3377	0.87	0/4551
1	H	0.77	1/3377 (0.0%)	0.88	1/4553 (0.0%)
All	All	0.77	1/27106 (0.0%)	0.87	10/36536 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	278	ASP	C-O	-6.25	1.20	1.23

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	18	ILE	N-CA-C	6.32	114.41	108.15
1	B	301	ILE	CB-CA-C	-6.05	105.31	110.94
1	H	256	ALA	N-CA-C	5.93	118.70	111.82
1	B	324	VAL	N-CA-C	5.51	116.21	108.17
1	C	398	GLU	CA-C-N	5.49	125.50	119.90
1	C	398	GLU	C-N-CA	5.49	125.50	119.90
1	A	434	VAL	CB-CA-C	-5.37	104.82	110.62
1	C	433	LYS	N-CA-C	5.14	116.97	111.36
1	C	244	PHE	CA-C-N	5.05	124.71	119.56
1	C	244	PHE	C-N-CA	5.05	124.71	119.56

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	3328	0	3295	20	0
1	B	3301	0	3262	35	0
1	C	3348	0	3320	17	0
1	D	3333	0	3299	22	0
1	E	3312	0	3279	33	0
1	F	3319	0	3293	16	0
1	G	3314	0	3276	29	0
1	H	3313	0	3279	22	0
2	A	12	0	16	3	0
2	B	6	0	8	3	0
2	C	6	0	8	1	0
2	D	6	0	8	3	0
2	E	6	0	8	1	0
2	F	6	0	8	2	0
2	G	6	0	8	4	0
2	H	6	0	8	0	0
3	A	103	0	0	2	0
3	B	84	0	0	1	0
3	C	71	0	0	0	0
3	D	64	0	0	0	0
3	E	61	0	0	3	0
3	F	32	0	0	0	0
3	G	53	0	0	1	0
3	H	74	0	0	1	0
All	All	27164	0	26375	192	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 4.

All (192) 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:213[A]:ARG:HG2	1:E:213[A]:ARG:HH11	1.08	1.15

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:161:ARG:HH11	1:B:161:ARG:HG3	1.27	1.00
1:D:445:MSE:HE1	1:D:461:LEU:CD1	1.94	0.97
1:E:84:ARG:HH11	1:E:84:ARG:HG3	1.26	0.96
1:B:445:MSE:HE1	1:B:461:LEU:HD11	1.46	0.96
1:F:445:MSE:HE1	1:F:461:LEU:HD21	1.52	0.92
1:C:93:ILE:HG22	1:C:93:ILE:O	1.73	0.86
1:D:445:MSE:HE1	1:D:461:LEU:HD11	1.58	0.86
1:E:213[A]:ARG:HG2	1:E:213[A]:ARG:NH1	1.83	0.85
1:F:445:MSE:HE1	1:F:461:LEU:CD2	2.07	0.83
1:H:243:ARG:HH11	1:H:243:ARG:HG2	1.46	0.80
1:B:445:MSE:HE1	1:B:461:LEU:CD1	2.12	0.80
1:B:161:ARG:HH11	1:B:161:ARG:CG	2.01	0.74
1:D:358:ASN:HD22	2:D:501:GOL:C1	2.03	0.71
1:C:93:ILE:O	1:C:93:ILE:CG2	2.39	0.70
1:B:445:MSE:HE3	1:B:449:VAL:HG23	1.75	0.69
1:B:176:ARG:HH11	1:B:176:ARG:HG2	1.59	0.68
1:G:34:HIS:HE1	2:G:501:GOL:H31	1.60	0.67
1:G:424:ILE:HG22	1:G:425:MSE:HE2	1.76	0.66
1:H:155:SER:OG	3:H:661:HOH:O	2.13	0.65
1:G:309:ALA:HB1	1:G:315:MSE:CE	2.26	0.65
1:G:309:ALA:HB1	1:G:315:MSE:HE2	1.77	0.65
1:E:439:ARG:HH22	2:E:501:GOL:H12	1.61	0.65
1:G:224:LYS:HD2	3:G:635:HOH:O	1.98	0.64
1:E:57:VAL:HB	1:E:58:PRO:HD2	1.80	0.64
1:D:358:ASN:HD22	2:D:501:GOL:H12	1.61	0.63
1:E:84:ARG:HG3	1:E:84:ARG:NH1	2.04	0.63
1:G:34:HIS:CE1	2:G:501:GOL:H31	2.34	0.62
1:G:424:ILE:HG22	1:G:425:MSE:CE	2.30	0.62
1:H:455:HIS:CG	1:H:458:ARG:HD2	2.35	0.62
1:B:19:PRO:HG3	1:B:64:ARG:HD3	1.80	0.62
1:H:243:ARG:HH11	1:H:243:ARG:CG	2.12	0.61
1:F:34:HIS:CE1	2:F:501:GOL:H11	2.36	0.60
1:E:255:GLY:O	1:E:457:LYS:HE3	2.01	0.60
1:E:84:ARG:HH11	1:E:84:ARG:CG	2.06	0.60
1:C:210:GLU:OE1	1:C:243:ARG:NH1	2.35	0.60
1:G:296:LYS:HE3	1:G:325:ASP:OD2	2.03	0.59
1:G:91:GLN:HG3	1:G:92:ALA:N	2.18	0.58
1:F:93:ILE:HD12	1:F:93:ILE:N	2.18	0.58
1:E:250:THR:HG21	1:E:275:TYR:CZ	2.39	0.57
1:A:429:GLU:O	1:A:433:LYS:HD2	2.04	0.57
1:D:93:ILE:HG21	1:D:126:LEU:HD13	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:445:MSE:HE1	1:D:461:LEU:HD13	1.82	0.57
1:B:445:MSE:HE3	1:B:449:VAL:CG2	2.33	0.57
1:D:164:MSE:HE2	1:D:382:ILE:HD11	1.87	0.57
1:G:93:ILE:CG2	1:G:93:ILE:O	2.53	0.57
1:A:164:MSE:HE2	1:A:382:ILE:HD11	1.85	0.56
1:H:169:PHE:CE2	1:H:224:LYS:HG2	2.41	0.56
1:A:358:ASN:OD1	2:A:502:GOL:C1	2.54	0.55
1:E:57:VAL:HB	1:E:58:PRO:CD	2.35	0.54
1:G:385:HIS:NE2	2:G:501:GOL:H32	2.22	0.54
1:H:318:SER:OG	1:H:324:VAL:HG13	2.07	0.54
1:E:277:GLU:OE2	1:E:306:ASN:ND2	2.42	0.53
1:E:250:THR:HG21	1:E:275:TYR:CE2	2.43	0.53
1:A:169:PHE:CE2	1:A:224:LYS:HG2	2.43	0.53
1:B:445:MSE:CE	1:B:461:LEU:CD1	2.85	0.53
1:F:358:ASN:O	1:F:360:HIS:HD2	1.92	0.53
1:F:89:LEU:O	1:F:93:ILE:HD13	2.08	0.53
1:A:17:VAL:O	2:A:501:GOL:H31	2.08	0.52
1:B:166:ALA:HB2	1:B:218:PHE:CD2	2.45	0.52
1:F:445:MSE:HE1	1:F:461:LEU:HD23	1.91	0.52
1:C:341:ARG:HA	1:C:344:GLN:HE21	1.73	0.52
1:D:35:ALA:HB1	1:D:36:PRO:HD2	1.91	0.52
1:B:161:ARG:HG3	1:B:161:ARG:NH1	2.08	0.52
1:F:57:VAL:HB	1:F:58:PRO:HD2	1.92	0.52
1:B:176:ARG:HG2	1:B:176:ARG:NH1	2.25	0.51
1:B:296:LYS:HE3	1:B:325:ASP:OD1	2.11	0.51
1:E:213[A]:ARG:HH11	1:E:213[A]:ARG:CG	1.98	0.51
1:B:348:GLU:HG2	1:D:312:TRP:HB2	1.93	0.51
1:D:93:ILE:HG21	1:D:126:LEU:CD1	2.41	0.51
1:H:93:ILE:O	1:H:94:ALA:C	2.54	0.51
1:E:341:ARG:HA	1:E:344:GLN:HE21	1.77	0.50
1:A:355:SER:HB3	1:A:382:ILE:HB	1.93	0.50
1:F:70:CYS:SG	1:F:93:ILE:HD11	2.52	0.50
1:G:18:ILE:HG21	1:G:425:MSE:HE1	1.94	0.49
1:H:243:ARG:HG2	1:H:243:ARG:NH1	2.17	0.49
1:B:165:LEU:C	1:B:165:LEU:HD12	2.37	0.49
1:G:93:ILE:HG22	1:G:126:LEU:HD13	1.93	0.49
1:E:161:ARG:NH1	3:E:616:HOH:O	2.37	0.49
1:F:34:HIS:HE1	2:F:501:GOL:H11	1.76	0.49
1:G:93:ILE:O	1:G:94:ALA:HB2	2.12	0.49
1:A:93:ILE:HD12	1:A:93:ILE:N	2.27	0.49
1:E:296:LYS:HE2	1:E:325:ASP:OD1	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:279:PRO:HD2	1:E:292:MSE:SE	2.63	0.49
1:E:213[A]:ARG:NH1	1:E:213[A]:ARG:CG	2.64	0.49
1:H:279:PRO:HD2	1:H:292:MSE:SE	2.63	0.48
1:D:296:LYS:HE3	1:D:325:ASP:OD1	2.12	0.48
1:D:70:CYS:SG	1:D:89:LEU:HD22	2.53	0.48
1:D:169:PHE:CE2	1:D:224:LYS:HG2	2.49	0.48
1:D:279:PRO:HD2	1:D:292:MSE:SE	2.64	0.48
1:B:306[A]:ASN:ND2	3:B:612:HOH:O	2.47	0.48
1:D:93:ILE:CG2	1:D:126:LEU:HD13	2.43	0.48
1:B:57:VAL:HB	1:B:58:PRO:HD2	1.95	0.47
1:F:66:ALA:O	1:F:69[B]:ARG:HB2	2.14	0.47
1:C:277:GLU:OE2	1:C:306[A]:ASN:ND2	2.48	0.47
1:G:439:ARG:HH22	2:G:501:GOL:H2	1.78	0.47
1:B:439:ARG:HH22	2:B:501:GOL:H32	1.80	0.47
1:H:47:ASP:OD1	1:H:47:ASP:C	2.57	0.47
1:H:165:LEU:HD12	1:H:165:LEU:C	2.40	0.47
1:B:161:ARG:CG	1:B:161:ARG:NH1	2.68	0.47
1:C:302:PRO:HA	1:C:325:ASP:OD2	2.15	0.47
1:A:93:ILE:N	1:A:93:ILE:CD1	2.78	0.46
1:B:445:MSE:CE	1:B:449:VAL:HG23	2.43	0.46
1:C:35:ALA:HB1	1:C:36:PRO:HD2	1.97	0.46
1:E:19:PRO:HG3	1:E:64:ARG:HG3	1.97	0.46
1:E:169:PHE:CE2	1:E:224:LYS:HG2	2.50	0.46
1:D:34:HIS:CE1	2:D:501:GOL:H2	2.50	0.46
1:H:166:ALA:HB2	1:H:218:PHE:CD2	2.50	0.46
1:A:224:LYS:HD2	3:A:643:HOH:O	2.15	0.46
1:B:445:MSE:CE	1:B:449:VAL:CG2	2.93	0.46
1:B:93:ILE:O	1:B:93:ILE:HG22	2.15	0.46
1:C:279:PRO:HD2	1:C:292:MSE:SE	2.66	0.46
1:H:250:THR:HG21	1:H:275:TYR:CZ	2.50	0.46
1:H:329:ALA:HB2	1:H:342:LEU:HD23	1.97	0.46
1:H:341:ARG:HA	1:H:344:GLN:HE21	1.81	0.46
1:A:165:LEU:C	1:A:165:LEU:HD12	2.41	0.46
1:A:358:ASN:OD1	2:A:502:GOL:H12	2.16	0.46
1:E:411:ASP:C	1:E:411:ASP:OD1	2.57	0.46
1:G:93:ILE:O	1:G:93:ILE:HG23	2.16	0.45
1:B:93:ILE:O	1:B:94:ALA:HB3	2.17	0.45
1:B:19:PRO:HG3	1:B:64:ARG:CD	2.46	0.45
1:E:84:ARG:NH1	1:E:84:ARG:CG	2.69	0.45
1:B:341:ARG:HA	1:B:344:GLN:HE21	1.81	0.45
1:A:159[A]:ARG:HD3	1:A:161:ARG:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:224:LYS:HD2	3:E:647:HOH:O	2.16	0.45
1:E:344:GLN:NE2	3:E:661:HOH:O	2.45	0.45
1:G:279:PRO:HD2	1:G:292:MSE:SE	2.67	0.45
1:E:430:LEU:O	1:E:434:VAL:HG22	2.17	0.44
1:A:326:ILE:HA	1:A:352:THR:O	2.18	0.44
1:E:173:GLU:HB3	1:E:176:ARG:HG3	1.98	0.44
1:B:296:LYS:HE3	1:B:325:ASP:CG	2.42	0.44
1:C:458:ARG:O	1:C:459:PRO:C	2.61	0.44
1:H:35:ALA:HB1	1:H:36:PRO:HD2	1.99	0.44
1:C:220:ASP:OD1	1:C:248:ARG:HD2	2.18	0.44
1:G:326:ILE:HA	1:G:352:THR:O	2.18	0.43
1:C:93:ILE:CD1	1:C:93:ILE:N	2.81	0.43
1:C:358:ASN:O	1:C:360:HIS:HD2	2.01	0.43
1:F:93:ILE:N	1:F:93:ILE:CD1	2.81	0.43
1:E:358:ASN:O	1:E:360:HIS:HD2	2.02	0.43
1:D:87:ASN:OD1	1:D:90:ARG:NH1	2.52	0.43
1:F:253:PRO:HD2	1:F:277:GLU:O	2.18	0.43
1:D:362:ASP:OD1	1:D:362:ASP:N	2.48	0.43
1:G:159[A]:ARG:HD3	1:G:161:ARG:O	2.18	0.43
1:H:155:SER:OG	1:H:155:SER:O	2.36	0.43
1:A:306[A]:ASN:ND2	3:A:635:HOH:O	2.51	0.43
1:C:216:TYR:HB2	1:C:218:PHE:CE2	2.54	0.43
1:D:250:THR:HG21	1:D:275:TYR:CE2	2.53	0.43
1:C:47:ASP:OD1	1:C:47:ASP:C	2.61	0.43
1:A:143:GLN:HA	1:A:413:PRO:HB2	2.00	0.42
1:G:26:MSE:HE2	1:G:448:LEU:HD11	2.01	0.42
1:E:164:MSE:HE2	1:E:382:ILE:HD11	2.00	0.42
1:B:329:ALA:HB2	1:B:342:LEU:HD23	2.01	0.42
1:D:255:GLY:O	1:D:457:LYS:HE3	2.19	0.42
1:F:169:PHE:CE2	1:F:224:LYS:HG2	2.53	0.42
1:H:46:LYS:HA	1:H:51:ARG:O	2.19	0.42
1:B:458:ARG:O	1:B:459:PRO:C	2.62	0.42
1:H:250:THR:HG21	1:H:275:TYR:CE2	2.54	0.42
1:H:57:VAL:HB	1:H:58:PRO:HD2	2.01	0.42
1:D:47:ASP:OD1	1:D:47:ASP:C	2.62	0.42
1:G:47:ASP:OD1	1:G:47:ASP:C	2.61	0.42
1:G:430:LEU:O	1:G:434:VAL:HB	2.20	0.42
1:A:279:PRO:HD2	1:A:292:MSE:SE	2.69	0.42
1:B:326:ILE:HA	1:B:352:THR:O	2.20	0.42
1:C:329:ALA:HB2	1:C:342:LEU:HD23	2.02	0.42
1:E:435:ALA:C	1:E:437:GLY:H	2.27	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:439:ARG:HH12	2:B:501:GOL:C3	2.33	0.41
1:E:315:MSE:HE2	1:E:319:LEU:HG	2.03	0.41
1:A:255:GLY:O	1:A:457:LYS:HE3	2.21	0.41
1:G:277:GLU:OE2	1:G:306[A]:ASN:ND2	2.52	0.41
1:A:149:VAL:HG23	1:A:414:GLY:C	2.44	0.41
1:E:151:GLU:HG3	1:G:81:ARG:NH1	2.35	0.41
1:F:137:LEU:HD23	1:F:137:LEU:HA	1.92	0.41
1:A:57:VAL:HB	1:A:58:PRO:HD2	2.01	0.41
1:B:411:ASP:OD1	1:B:411:ASP:C	2.60	0.41
1:E:169:PHE:CD2	1:E:224:LYS:HG2	2.55	0.41
1:B:439:ARG:NH2	2:B:501:GOL:H32	2.36	0.41
1:B:159:ARG:HH11	1:B:159:ARG:HD3	1.73	0.41
1:F:164:MSE:HE3	1:F:407:VAL:CG2	2.50	0.41
1:G:309:ALA:HB1	1:G:315:MSE:HE1	2.02	0.41
1:G:341:ARG:HA	1:G:344:GLN:HE21	1.86	0.41
1:G:428:HIS:O	1:G:431:TYR:HB3	2.21	0.41
1:H:430:LEU:O	1:H:434:VAL:HG22	2.21	0.41
1:B:250:THR:HG21	1:B:275:TYR:CE2	2.56	0.41
1:C:169:PHE:CE2	1:C:224:LYS:HG2	2.55	0.41
1:B:303:THR:OG1	1:B:324:VAL:HA	2.21	0.40
1:D:57:VAL:HB	1:D:58:PRO:HD2	2.04	0.40
1:G:304:ALA:HA	1:G:326:ILE:O	2.21	0.40
1:H:325:ASP:C	1:H:326:ILE:HG13	2.44	0.40
1:C:385:HIS:NE2	2:C:501:GOL:O2	2.43	0.40
1:G:46:LYS:HD2	1:G:52:THR:OG1	2.22	0.40
1:E:355:SER:HB3	1:E:382:ILE:HB	2.03	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	427/473 (90%)	417 (98%)	10 (2%)	0	100	100
1	B	423/473 (89%)	414 (98%)	9 (2%)	0	100	100
1	C	430/473 (91%)	420 (98%)	8 (2%)	2 (0%)	24	27
1	D	429/473 (91%)	419 (98%)	10 (2%)	0	100	100
1	E	425/473 (90%)	414 (97%)	10 (2%)	1 (0%)	43	51
1	F	425/473 (90%)	408 (96%)	14 (3%)	3 (1%)	18	19
1	G	425/473 (90%)	415 (98%)	10 (2%)	0	100	100
1	H	425/473 (90%)	419 (99%)	6 (1%)	0	100	100
All	All	3409/3784 (90%)	3326 (98%)	77 (2%)	6 (0%)	48	51

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	462	GLY
1	E	438	ALA
1	F	187	SER
1	C	93	ILE
1	F	69[A]	ARG
1	F	69[B]	ARG

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	335/356 (94%)	327 (98%)	8 (2%)	43	58
1	B	332/356 (93%)	322 (97%)	10 (3%)	36	49
1	C	338/356 (95%)	325 (96%)	13 (4%)	29	40
1	D	336/356 (94%)	327 (97%)	9 (3%)	39	53
1	E	333/356 (94%)	324 (97%)	9 (3%)	39	53
1	F	334/356 (94%)	322 (96%)	12 (4%)	31	42
1	G	333/356 (94%)	321 (96%)	12 (4%)	31	42

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	H	333/356 (94%)	324 (97%)	9 (3%)	39 53
All	All	2674/2848 (94%)	2592 (97%)	82 (3%)	37 48

All (82) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	161	ARG
1	A	210	GLU
1	A	254	ASN
1	A	348	GLU
1	A	362	ASP
1	A	400	LEU
1	A	457	LYS
1	A	463	ARG
1	B	161	ARG
1	B	187	SER
1	B	297	ARG
1	B	306[A]	ASN
1	B	306[B]	ASN
1	B	324	VAL
1	B	406	GLN
1	B	418	GLU
1	B	457	LYS
1	B	461	LEU
1	C	93	ILE
1	C	185	LYS
1	C	190	ASP
1	C	203	ASP
1	C	207	ARG
1	C	243	ARG
1	C	248	ARG
1	C	254	ASN
1	C	319	LEU
1	C	324	VAL
1	C	348[A]	GLU
1	C	348[B]	GLU
1	C	425	MSE
1	D	93	ILE
1	D	243	ARG
1	D	306	ASN
1	D	348[A]	GLU
1	D	348[B]	GLU

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Mol	Chain	Res	Type
1	D	400	LEU
1	D	408	GLN
1	D	457	LYS
1	D	461	LEU
1	E	64	ARG
1	E	84	ARG
1	E	160	GLN
1	E	176	ARG
1	E	213[A]	ARG
1	E	213[B]	ARG
1	E	241	LYS
1	E	324	VAL
1	E	461	LEU
1	F	8	SER
1	F	91	GLN
1	F	161	ARG
1	F	215	ARG
1	F	241	LYS
1	F	254	ASN
1	F	267	LYS
1	F	306	ASN
1	F	319	LEU
1	F	337	GLN
1	F	362	ASP
1	F	425	MSE
1	G	91	GLN
1	G	127	ASP
1	G	187	SER
1	G	189	ASP
1	G	190	ASP
1	G	243	ARG
1	G	254	ASN
1	G	348	GLU
1	G	400	LEU
1	G	434	VAL
1	G	461	LEU
1	G	463	ARG
1	H	93	ILE
1	H	155	SER
1	H	175	GLN
1	H	243	ARG
1	H	254	ASN

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Mol	Chain	Res	Type
1	H	306	ASN
1	H	324	VAL
1	H	362	ASP
1	H	400	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (35) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	91	GLN
1	A	128	ASN
1	A	406	GLN
1	B	41	ASN
1	B	77	GLN
1	B	128	ASN
1	B	344	GLN
1	B	408	GLN
1	C	143	GLN
1	C	284	ASN
1	C	344	GLN
1	C	422	GLN
1	D	77	GLN
1	D	83	ASN
1	D	254	ASN
1	D	358	ASN
1	D	453	GLN
1	E	16	GLN
1	E	344	GLN
1	E	358	ASN
1	F	128	ASN
1	F	143	GLN
1	F	144	HIS
1	F	269	GLN
1	F	306	ASN
1	F	337	GLN
1	F	408	GLN
1	G	128	ASN
1	G	193	HIS
1	G	344	GLN
1	H	65	GLN
1	H	128	ASN
1	H	144	HIS
1	H	193	HIS

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Mol	Chain	Res	Type
1	H	344	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 ⓘ

9 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	E	501	-	5,5,5	0.62	0	5,5,5	1.47	1 (20%)
2	GOL	B	501	-	5,5,5	0.78	0	5,5,5	0.90	0
2	GOL	F	501	-	5,5,5	0.62	0	5,5,5	0.87	0
2	GOL	C	501	-	5,5,5	0.67	0	5,5,5	0.71	0
2	GOL	H	501	-	5,5,5	0.95	0	5,5,5	1.38	2 (40%)
2	GOL	G	501	-	5,5,5	0.67	0	5,5,5	0.79	0
2	GOL	D	501	-	5,5,5	0.64	0	5,5,5	1.23	0
2	GOL	A	501	-	5,5,5	0.23	0	5,5,5	0.32	0
2	GOL	A	502	-	5,5,5	0.70	0	5,5,5	1.36	1 (20%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the

Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns.
'-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	E	501	-	-	4/4/4/4	-
2	GOL	B	501	-	-	2/4/4/4	-
2	GOL	F	501	-	-	4/4/4/4	-
2	GOL	C	501	-	-	2/4/4/4	-
2	GOL	H	501	-	-	0/4/4/4	-
2	GOL	G	501	-	-	3/4/4/4	-
2	GOL	D	501	-	-	4/4/4/4	-
2	GOL	A	501	-	-	2/4/4/4	-
2	GOL	A	502	-	-	2/4/4/4	-

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	501	GOL	O2-C2-C1	-2.51	98.80	109.18
2	H	501	GOL	O2-C2-C1	-2.17	100.21	109.18
2	A	502	GOL	O2-C2-C3	-2.14	100.32	109.18
2	H	501	GOL	O3-C3-C2	2.02	119.46	110.38

There are no chirality outliers.

All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	GOL	O1-C1-C2-O2
2	A	501	GOL	O1-C1-C2-C3
2	A	502	GOL	O1-C1-C2-O2
2	A	502	GOL	O1-C1-C2-C3
2	C	501	GOL	C1-C2-C3-O3
2	D	501	GOL	O1-C1-C2-C3
2	D	501	GOL	C1-C2-C3-O3
2	E	501	GOL	C1-C2-C3-O3
2	F	501	GOL	O1-C1-C2-O2
2	F	501	GOL	O1-C1-C2-C3
2	G	501	GOL	O1-C1-C2-O2
2	G	501	GOL	O1-C1-C2-C3
2	B	501	GOL	O1-C1-C2-C3
2	E	501	GOL	O1-C1-C2-C3
2	F	501	GOL	C1-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
2	C	501	GOL	O2-C2-C3-O3
2	D	501	GOL	O1-C1-C2-O2
2	D	501	GOL	O2-C2-C3-O3
2	E	501	GOL	O2-C2-C3-O3
2	E	501	GOL	O1-C1-C2-O2
2	F	501	GOL	O2-C2-C3-O3
2	B	501	GOL	O1-C1-C2-O2
2	G	501	GOL	O2-C2-C3-O3

There are no ring outliers.

8 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	501	GOL	1	0
2	B	501	GOL	3	0
2	F	501	GOL	2	0
2	C	501	GOL	1	0
2	G	501	GOL	4	0
2	D	501	GOL	3	0
2	A	501	GOL	1	0
2	A	502	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	413/473 (87%)	-0.48	5 (1%) 76 74	7, 18, 39, 70	4 (0%)
1	B	412/473 (87%)	-0.42	5 (1%) 76 74	10, 20, 41, 66	1 (0%)
1	C	413/473 (87%)	-0.30	9 (2%) 62 59	9, 21, 45, 88	7 (1%)
1	D	413/473 (87%)	-0.20	8 (1%) 66 63	10, 22, 47, 82	6 (1%)
1	E	412/473 (87%)	-0.19	6 (1%) 72 69	14, 25, 46, 74	3 (0%)
1	F	412/473 (87%)	-0.05	6 (1%) 72 69	14, 28, 50, 77	3 (0%)
1	G	412/473 (87%)	-0.24	7 (1%) 69 66	10, 24, 47, 72	3 (0%)
1	H	413/473 (87%)	-0.24	11 (2%) 56 53	10, 22, 48, 91	2 (0%)
All	All	3300/3784 (87%)	-0.27	57 (1%) 69 66	7, 23, 46, 91	29 (0%)

All (57) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	94	ALA	6.4
1	B	94	ALA	5.3
1	G	94	ALA	5.3
1	C	94	ALA	4.9
1	A	94	ALA	4.2
1	D	437	GLY	4.1
1	H	93	ILE	4.0
1	H	436	SER	3.9
1	H	124	LEU	3.9
1	C	397[A]	ARG	3.9
1	D	436	SER	3.8
1	A	124	LEU	3.8
1	E	437	GLY	3.6
1	C	93	ILE	3.4
1	H	437	GLY	3.4
1	D	124	LEU	3.4

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Mol	Chain	Res	Type	RSRZ
1	C	436	SER	3.3
1	E	94	ALA	3.3
1	A	436	SER	3.3
1	C	438	ALA	3.2
1	A	8	SER	3.2
1	D	94	ALA	3.1
1	H	438	ALA	3.1
1	E	438	ALA	3.0
1	F	437	GLY	3.0
1	F	94	ALA	3.0
1	D	185	LYS	3.0
1	C	437	GLY	2.9
1	F	438	ALA	2.9
1	H	125	ARG	2.8
1	E	125	ARG	2.7
1	G	125	ARG	2.7
1	B	411	ASP	2.6
1	C	462	GLY	2.6
1	F	50	GLY	2.5
1	B	8	SER	2.5
1	B	437	GLY	2.5
1	E	434	VAL	2.5
1	D	8	SER	2.4
1	E	8	SER	2.4
1	C	124	LEU	2.4
1	H	126	LEU	2.4
1	G	185	LYS	2.4
1	D	438	ALA	2.4
1	F	436	SER	2.4
1	G	463	ARG	2.3
1	H	411	ASP	2.3
1	B	189	ASP	2.3
1	F	435	ALA	2.3
1	H	8	SER	2.3
1	A	348	GLU	2.2
1	H	185	LYS	2.2
1	G	8	SER	2.1
1	G	436	SER	2.1
1	D	411	ASP	2.1
1	G	93	ILE	2.1
1	C	185	LYS	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($\text{\AA}^2$)	Q<0.9
2	GOL	F	501	6/6	0.84	0.18	29,37,45,46	0
2	GOL	G	501	6/6	0.84	0.17	23,32,37,52	0
2	GOL	C	501	6/6	0.85	0.16	24,32,34,43	0
2	GOL	D	501	6/6	0.87	0.16	29,32,41,42	0
2	GOL	A	502	6/6	0.89	0.12	20,26,28,37	0
2	GOL	A	501	6/6	0.90	0.11	35,36,44,45	0
2	GOL	H	501	6/6	0.90	0.15	21,26,32,37	0
2	GOL	B	501	6/6	0.92	0.10	21,27,28,36	0
2	GOL	E	501	6/6	0.93	0.12	30,32,38,44	0

6.5 Other polymers [i](#)

There are no such residues in this entry.