



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 06:29 AM UTC

PDB ID : 4J5I / pdb_00004j5i
Title : Crystal structure of an alpha-ketoglutarate-dependent taurine dioxygenase from *Mycobacterium smegmatis*
Authors : Seattle Structural Genomics Center for Infectious Disease (SSGCID)
Deposited on : 2013-02-08
Resolution : 2.60 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:

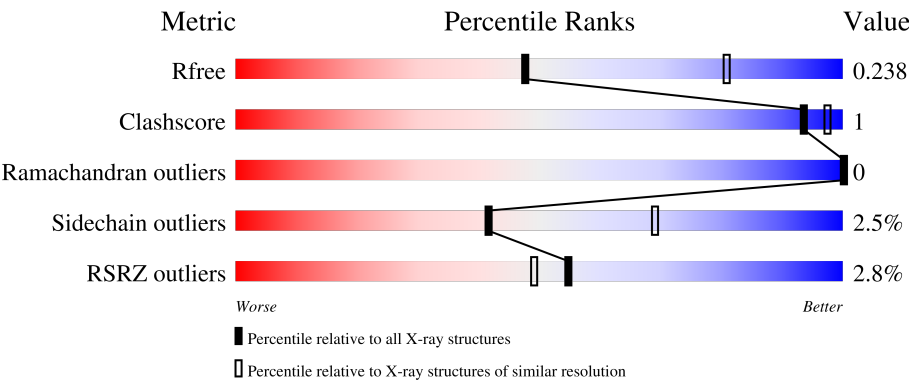
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

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

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	304	3% 78% 18%
1	B	304	% 89% 5% 6%
1	C	304	4% 86% 11%
1	D	304	2% 75% 21%
1	E	304	2% 78% 18%

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Mol	Chain	Length	Quality of chain
1	F	304	 % 79% 18% 6%
1	G	304	 6% 83% 13%
1	H	304	 % 78% 18% 6%

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
3	EDO	A	402	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 16323 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 Alpha-ketoglutarate-dependent taurine dioxygenase.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	248	Total	C	N	O	0	0	0
			1911	1217	339	355			
1	B	285	Total	C	N	O	0	0	0
			2186	1385	385	416			
1	C	271	Total	C	N	O	0	0	0
			2067	1311	363	393			
1	D	239	Total	C	N	O	0	0	0
			1839	1171	325	343			
1	E	250	Total	C	N	O	0	1	0
			1925	1225	338	362			
1	F	250	Total	C	N	O	0	0	0
			1929	1224	339	366			
1	G	264	Total	C	N	O	0	0	0
			2026	1289	356	381			
1	H	249	Total	C	N	O	0	0	0
			1918	1219	340	359			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	expression tag	UNP A0QZ23
A	-2	PRO	-	expression tag	UNP A0QZ23
A	-1	GLY	-	expression tag	UNP A0QZ23
A	0	SER	-	expression tag	UNP A0QZ23
A	1	MET	-	expression tag	UNP A0QZ23
A	2	VAL	-	expression tag	UNP A0QZ23
B	-3	GLY	-	expression tag	UNP A0QZ23
B	-2	PRO	-	expression tag	UNP A0QZ23
B	-1	GLY	-	expression tag	UNP A0QZ23
B	0	SER	-	expression tag	UNP A0QZ23
B	1	MET	-	expression tag	UNP A0QZ23
B	2	VAL	-	expression tag	UNP A0QZ23
C	-3	GLY	-	expression tag	UNP A0QZ23

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	PRO	-	expression tag	UNP A0QZ23
C	-1	GLY	-	expression tag	UNP A0QZ23
C	0	SER	-	expression tag	UNP A0QZ23
C	1	MET	-	expression tag	UNP A0QZ23
C	2	VAL	-	expression tag	UNP A0QZ23
D	-3	GLY	-	expression tag	UNP A0QZ23
D	-2	PRO	-	expression tag	UNP A0QZ23
D	-1	GLY	-	expression tag	UNP A0QZ23
D	0	SER	-	expression tag	UNP A0QZ23
D	1	MET	-	expression tag	UNP A0QZ23
D	2	VAL	-	expression tag	UNP A0QZ23
E	-3	GLY	-	expression tag	UNP A0QZ23
E	-2	PRO	-	expression tag	UNP A0QZ23
E	-1	GLY	-	expression tag	UNP A0QZ23
E	0	SER	-	expression tag	UNP A0QZ23
E	1	MET	-	expression tag	UNP A0QZ23
E	2	VAL	-	expression tag	UNP A0QZ23
F	-3	GLY	-	expression tag	UNP A0QZ23
F	-2	PRO	-	expression tag	UNP A0QZ23
F	-1	GLY	-	expression tag	UNP A0QZ23
F	0	SER	-	expression tag	UNP A0QZ23
F	1	MET	-	expression tag	UNP A0QZ23
F	2	VAL	-	expression tag	UNP A0QZ23
G	-3	GLY	-	expression tag	UNP A0QZ23
G	-2	PRO	-	expression tag	UNP A0QZ23
G	-1	GLY	-	expression tag	UNP A0QZ23
G	0	SER	-	expression tag	UNP A0QZ23
G	1	MET	-	expression tag	UNP A0QZ23
G	2	VAL	-	expression tag	UNP A0QZ23
H	-3	GLY	-	expression tag	UNP A0QZ23
H	-2	PRO	-	expression tag	UNP A0QZ23
H	-1	GLY	-	expression tag	UNP A0QZ23
H	0	SER	-	expression tag	UNP A0QZ23
H	1	MET	-	expression tag	UNP A0QZ23
H	2	VAL	-	expression tag	UNP A0QZ23

- Molecule 2 is FE (III) ION (CCD ID: FE) (formula: Fe).

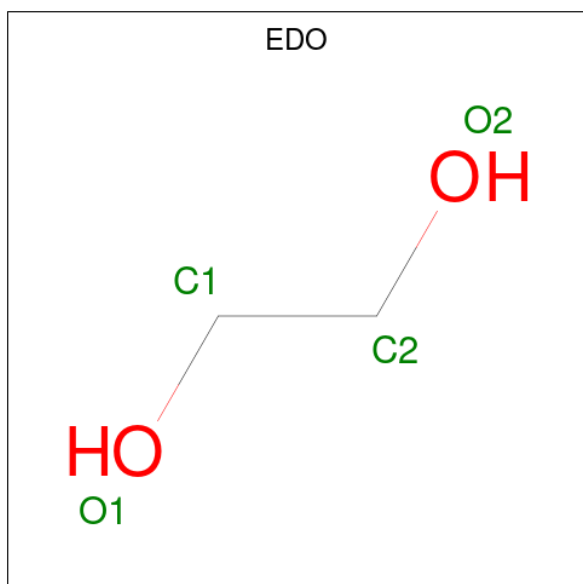
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Fe 1 1	0	0
2	B	1	Total Fe 1 1	0	0

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

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	F	1	Total	Mg	0	0
			1	1		

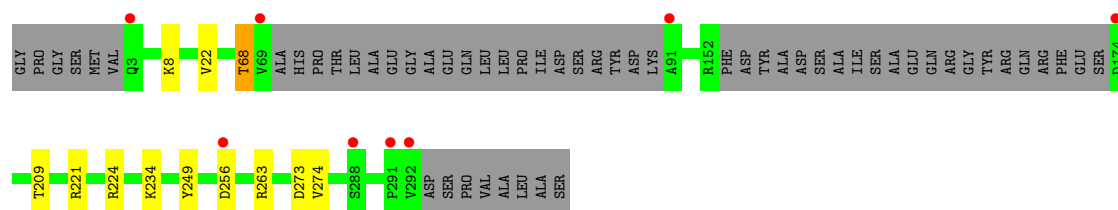
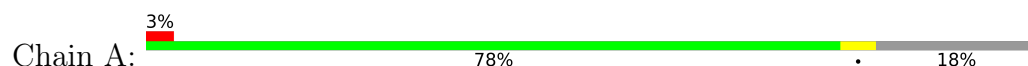
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	60	Total 60	O 60	0	0
5	B	88	Total 88	O 88	0	0
5	C	63	Total 63	O 63	0	0
5	D	44	Total 44	O 44	0	0
5	E	67	Total 67	O 67	0	0
5	F	67	Total 67	O 67	0	0
5	G	63	Total 63	O 63	0	0
5	H	57	Total 57	O 57	0	0

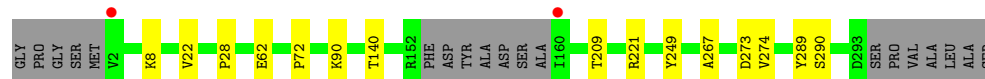
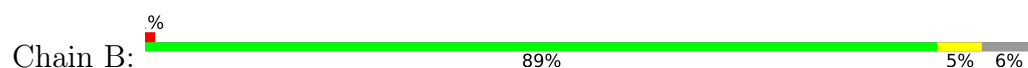
3 Residue-property plots [i](#)

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.

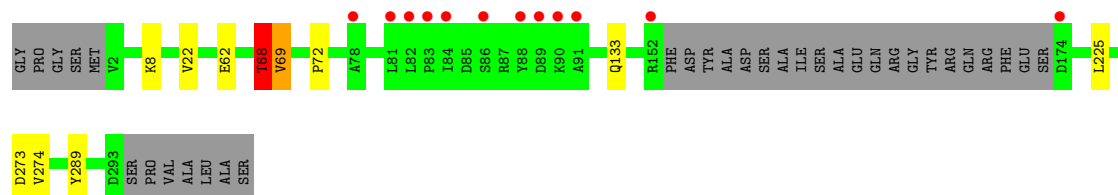
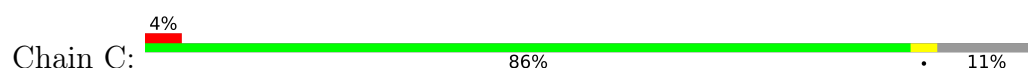
- Molecule 1: Alpha-ketoglutarate-dependent taurine dioxygenase



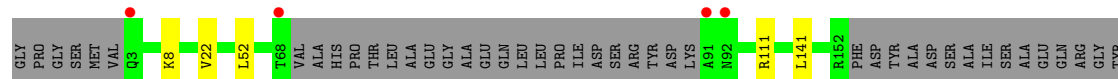
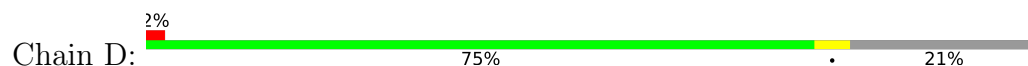
- Molecule 1: Alpha-ketoglutarate-dependent taurine dioxygenase



- Molecule 1: Alpha-ketoglutarate-dependent taurine dioxygenase

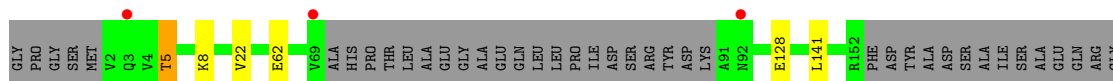
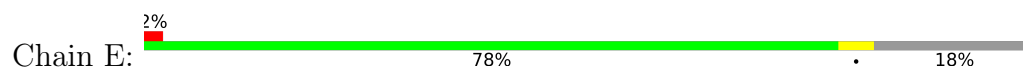


- Molecule 1: Alpha-ketoglutarate-dependent taurine dioxygenase

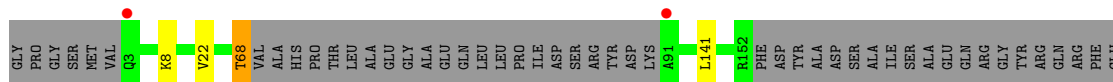
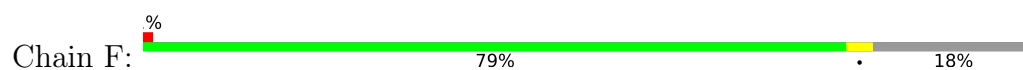




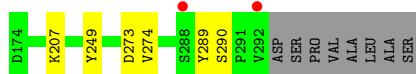
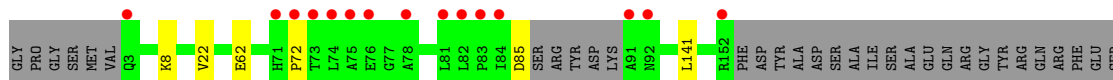
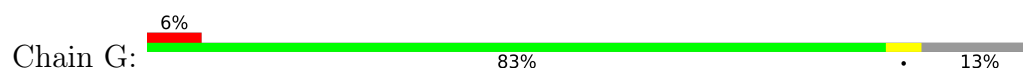
- Molecule 1: Alpha-ketoglutarate-dependent taurine dioxygenase



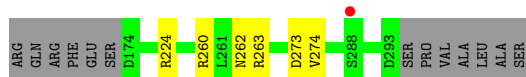
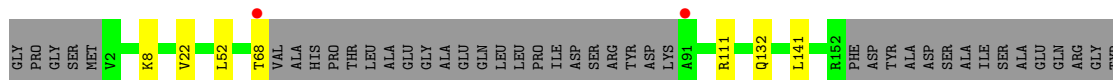
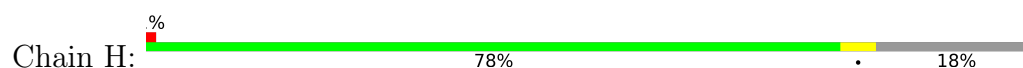
- Molecule 1: Alpha-ketoglutarate-dependent taurine dioxygenase



- Molecule 1: Alpha-ketoglutarate-dependent taurine dioxygenase



- Molecule 1: Alpha-ketoglutarate-dependent taurine dioxygenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	77.01Å 166.68Å 99.72Å 90.00° 90.57° 90.00°	Depositor
Resolution (Å)	49.40 – 2.60 49.40 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.0 (49.40-2.60) 99.0 (49.40-2.60)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.81 (at 2.61Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.211 , 0.234 0.215 , 0.238	Depositor DCC
R_{free} test set	3833 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	26.6	Xtriage
Anisotropy	0.043	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 26.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.045 for h,-k,-l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	16323	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.17% 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: FE, EDO, 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	0.67	0/1957	0.83	4/2675 (0.1%)
1	B	0.73	1/2239 (0.0%)	0.82	1/3063 (0.0%)
1	C	0.65	0/2117	0.77	1/2901 (0.0%)
1	D	0.64	0/1883	0.77	1/2576 (0.0%)
1	E	0.64	1/1974 (0.1%)	0.78	4/2702 (0.1%)
1	F	0.66	0/1976	0.76	1/2704 (0.0%)
1	G	0.63	0/2075	0.78	1/2840 (0.0%)
1	H	0.64	0/1964	0.77	0/2688
All	All	0.66	2/16185 (0.0%)	0.79	13/22149 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	5	THR	CA-CB	6.08	1.62	1.53
1	B	140	THR	CA-CB	5.20	1.61	1.53

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	263	ARG	NE-CZ-NH2	8.52	126.87	119.20
1	A	263	ARG	NE-CZ-NH1	-6.78	114.72	121.50
1	A	263	ARG	CD-NE-CZ	6.59	133.62	124.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	68	THR	N-CA-CB	-5.79	101.20	110.46
1	E	278	ARG	CA-CB-CG	5.78	125.66	114.10
1	E	234	LYS	CD-CE-NZ	5.66	130.02	111.90
1	D	234	LYS	CD-CE-NZ	5.57	129.71	111.90
1	F	68	THR	N-CA-CB	-5.39	102.34	111.50
1	B	290	SER	N-CA-C	5.38	112.97	108.13
1	E	278	ARG	CB-CG-CD	-5.23	99.27	111.30
1	A	68	THR	N-CA-CB	-5.16	102.45	110.29
1	E	290	SER	N-CA-C	5.15	112.76	108.13
1	G	290	SER	N-CA-C	5.01	112.64	108.13

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	256	ASP	Peptide

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1911	0	1849	8	0
1	B	2186	0	2078	6	0
1	C	2067	0	1972	11	0
1	D	1839	0	1764	4	0
1	E	1925	0	1840	4	0
1	F	1929	0	1849	5	0
1	G	2026	0	1952	5	0
1	H	1918	0	1839	6	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

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	4	0	6	4	0
4	F	1	0	0	0	0
5	A	60	0	0	1	0
5	B	88	0	0	3	0
5	C	63	0	0	1	0
5	D	44	0	0	0	0
5	E	67	0	0	1	0
5	F	67	0	0	1	0
5	G	63	0	0	0	0
5	H	57	0	0	1	0
All	All	16323	0	15149	42	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 1.

All (42) 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:C:68:THR:HG22	1:C:69:VAL:O	1.51	1.09
1:B:221:ARG:HD3	5:B:520:HOH:O	1.96	0.66
1:A:224:ARG:HD2	3:A:402:EDO:O2	1.98	0.64
1:F:263:ARG:NH1	5:F:561:HOH:O	2.28	0.59
1:F:269:ASP:HB3	1:F:293:ASP:O	2.03	0.59
1:C:68:THR:CG2	1:C:69:VAL:O	2.41	0.56
1:A:224:ARG:HD2	3:A:402:EDO:C2	2.37	0.54
1:E:263:ARG:NH1	5:E:564:HOH:O	2.24	0.54
1:H:260:ARG:HH11	1:H:262:ASN:CG	2.15	0.53
1:A:234:LYS:HE3	5:A:508:HOH:O	2.09	0.51
1:A:249:TYR:CD2	1:C:225:LEU:HD12	2.46	0.51
1:C:68:THR:CG2	1:C:289:TYR:CE1	2.95	0.50
1:C:69:VAL:HG12	1:C:289:TYR:HD1	1.78	0.49
1:C:72:PRO:HD3	1:C:289:TYR:CE1	2.48	0.49
1:C:133:GLN:HB3	5:C:518:HOH:O	2.12	0.49
1:E:225:LEU:HD12	1:G:249:TYR:CD2	2.46	0.49
1:A:221:ARG:HA	3:A:402:EDO:H12	1.92	0.49
1:B:249:TYR:CD2	1:D:225:LEU:HD12	2.48	0.49
1:H:263:ARG:NH1	5:H:553:HOH:O	2.21	0.49
1:G:72:PRO:HD3	1:G:289:TYR:CE1	2.48	0.48
1:E:128:GLU:OE2	1:E:191[A]:GLU:CD	2.57	0.48
1:E:141:LEU:HD22	1:F:141:LEU:HD22	1.96	0.48
1:B:72:PRO:HD3	1:B:289:TYR:CE1	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:141:LEU:HD22	1:H:141:LEU:HD22	1.96	0.46
1:C:69:VAL:HG12	1:C:289:TYR:CD1	2.49	0.46
1:F:273:ASP:C	1:F:273:ASP:OD1	2.58	0.46
1:B:28:PRO:HA	5:B:585:HOH:O	2.16	0.45
1:D:141:LEU:HD21	1:D:214:LEU:HD13	1.99	0.44
1:C:68:THR:HG21	1:C:289:TYR:CZ	2.52	0.44
1:H:52:LEU:HD23	1:H:111:ARG:HD2	2.00	0.44
1:F:141:LEU:HD21	1:F:214:LEU:HD13	1.99	0.44
1:A:273:ASP:OD2	1:A:273:ASP:C	2.61	0.44
1:D:273:ASP:OD1	1:D:273:ASP:C	2.60	0.44
1:G:273:ASP:OD2	1:G:273:ASP:C	2.60	0.44
1:A:224:ARG:HD2	3:A:402:EDO:H22	2.00	0.43
1:B:267:ALA:HA	5:B:505:HOH:O	2.19	0.43
1:D:52:LEU:HD23	1:D:111:ARG:HD2	2.02	0.41
1:B:273:ASP:OD2	1:B:273:ASP:C	2.62	0.41
1:H:273:ASP:OD1	1:H:273:ASP:C	2.61	0.41
1:A:249:TYR:CG	1:C:225:LEU:HD12	2.56	0.40
1:C:273:ASP:OD2	1:C:273:ASP:C	2.64	0.40
1:G:207:LYS:HZ3	1:H:224:ARG:CZ	2.35	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	242/304 (80%)	235 (97%)	7 (3%)	0	100	100
1	B	281/304 (92%)	271 (96%)	10 (4%)	0	100	100
1	C	267/304 (88%)	257 (96%)	10 (4%)	0	100	100
1	D	233/304 (77%)	225 (97%)	8 (3%)	0	100	100
1	E	245/304 (81%)	238 (97%)	7 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	244/304 (80%)	238 (98%)	6 (2%)	0	100	100
1	G	258/304 (85%)	252 (98%)	6 (2%)	0	100	100
1	H	243/304 (80%)	236 (97%)	7 (3%)	0	100	100
All	All	2013/2432 (83%)	1952 (97%)	61 (3%)	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	195/248 (79%)	190 (97%)	5 (3%)	40	68
1	B	221/248 (89%)	215 (97%)	6 (3%)	39	67
1	C	210/248 (85%)	204 (97%)	6 (3%)	37	65
1	D	186/248 (75%)	182 (98%)	4 (2%)	45	72
1	E	195/248 (79%)	190 (97%)	5 (3%)	40	68
1	F	199/248 (80%)	195 (98%)	4 (2%)	48	74
1	G	207/248 (84%)	202 (98%)	5 (2%)	43	70
1	H	195/248 (79%)	190 (97%)	5 (3%)	40	68
All	All	1608/1984 (81%)	1568 (98%)	40 (2%)	42	69

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	LYS
1	A	22	VAL
1	A	68	THR
1	A	209	THR
1	A	274	VAL
1	B	8	LYS
1	B	22	VAL
1	B	62	GLU

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Mol	Chain	Res	Type
1	B	90	LYS
1	B	209	THR
1	B	274	VAL
1	C	8	LYS
1	C	22	VAL
1	C	62	GLU
1	C	68	THR
1	C	69	VAL
1	C	274	VAL
1	D	8	LYS
1	D	22	VAL
1	D	209	THR
1	D	274	VAL
1	E	5	THR
1	E	8	LYS
1	E	22	VAL
1	E	62	GLU
1	E	274	VAL
1	F	8	LYS
1	F	22	VAL
1	F	68	THR
1	F	274	VAL
1	G	8	LYS
1	G	22	VAL
1	G	62	GLU
1	G	85	ASP
1	G	274	VAL
1	H	8	LYS
1	H	22	VAL
1	H	68	THR
1	H	132	GLN
1	H	274	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	132	GLN
1	D	132	GLN
1	H	132	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](#)

Of 10 ligands modelled in this entry, 9 are monoatomic - leaving 1 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	EDO	A	402	-	3,3,3	0.41	0	2,2,2	0.42	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	EDO	A	402	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	402	EDO	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	402	EDO	4	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	248/304 (81%)	-0.06	8 (3%)	50	44	15, 23, 45, 58	0
1	B	285/304 (93%)	-0.19	2 (0%)	84	82	11, 20, 35, 44	0
1	C	271/304 (89%)	0.17	12 (4%)	39	33	16, 29, 56, 72	0
1	D	239/304 (78%)	0.05	6 (2%)	58	52	14, 27, 44, 56	0
1	E	250/304 (82%)	-0.06	6 (2%)	59	54	13, 24, 45, 68	1 (0%)
1	F	250/304 (82%)	-0.12	4 (1%)	70	66	14, 22, 38, 62	0
1	G	264/304 (86%)	0.27	17 (6%)	25	20	15, 31, 64, 79	0
1	H	249/304 (81%)	-0.02	3 (1%)	76	73	14, 27, 43, 58	0
All	All	2056/2432 (84%)	0.01	58 (2%)	55	49	11, 25, 47, 79	1 (0%)

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	91	ALA	4.9
1	D	91	ALA	4.8
1	G	91	ALA	4.5
1	A	69	VAL	4.2
1	G	152	ARG	3.9
1	C	90	LYS	3.8
1	A	174	ASP	3.8
1	G	73	THR	3.6
1	C	89	ASP	3.6
1	C	81	LEU	3.6
1	C	83	PRO	3.5
1	G	74	LEU	3.5
1	D	174	ASP	3.4
1	H	288	SER	3.4
1	G	3	GLN	3.4
1	C	86	SER	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	160	ILE	3.3
1	G	75	ALA	3.3
1	A	288	SER	3.3
1	G	78	ALA	3.3
1	F	295	PRO	3.2
1	D	68	THR	3.1
1	D	284	GLY	3.1
1	F	3	GLN	3.0
1	G	71	HIS	3.0
1	A	292	VAL	2.9
1	G	76	GLU	2.9
1	G	292	VAL	2.8
1	G	81	LEU	2.8
1	C	78	ALA	2.8
1	C	174	ASP	2.7
1	E	288	SER	2.7
1	G	92	ASN	2.7
1	H	91	ALA	2.7
1	E	3	GLN	2.7
1	G	83	PRO	2.6
1	E	92	ASN	2.6
1	G	288	SER	2.6
1	D	92	ASN	2.5
1	C	88	TYR	2.5
1	C	152	ARG	2.4
1	G	72	PRO	2.4
1	A	3	GLN	2.4
1	C	91	ALA	2.4
1	F	256	ASP	2.4
1	C	84	ILE	2.3
1	G	82	LEU	2.3
1	E	287	SER	2.3
1	F	91	ALA	2.2
1	E	69	VAL	2.2
1	A	256	ASP	2.2
1	D	3	GLN	2.1
1	A	291	PRO	2.1
1	B	2	VAL	2.1
1	C	82	LEU	2.1
1	E	289	TYR	2.1
1	G	84	ILE	2.0
1	H	68	THR	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	FE	B	401	1/1	0.92	0.07	54,54,54,54	0
3	EDO	A	402	4/4	0.93	0.09	27,29,29,30	0
2	FE	G	401	1/1	0.96	0.07	56,56,56,56	0
2	FE	C	401	1/1	0.97	0.07	46,46,46,46	0
2	FE	F	401	1/1	0.97	0.06	38,38,38,38	0
2	FE	A	401	1/1	0.98	0.04	33,33,33,33	0
2	FE	E	401	1/1	0.99	0.02	41,41,41,41	0
2	FE	H	401	1/1	0.99	0.03	43,43,43,43	0
2	FE	D	401	1/1	0.99	0.07	35,35,35,35	0
4	MG	F	402	1/1	0.99	0.07	3,3,3,3	0

6.5 Other polymers [i](#)

There are no such residues in this entry.