



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 8, 2016 – 12:27 PM EDT

PDB ID : 5H80
Title : Biotin Carboxylase domain of single-chain bacterial carboxylase
Authors : Hagmann, A.; Hunkeler, M.; Stuttfeld, E.; Maier, T.
Deposited on : 2015-12-23
Resolution : 1.70 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027939
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20027939

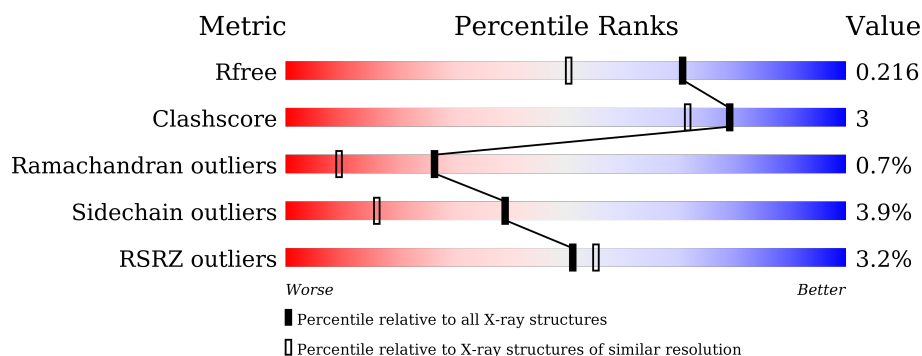
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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}	91344	3190 (1.70-1.70)
Clashscore	102246	3585 (1.70-1.70)
Ramachandran outliers	100387	3527 (1.70-1.70)
Sidechain outliers	100360	3527 (1.70-1.70)
RSRZ outliers	91569	3200 (1.70-1.70)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	494	<div> <div>4%</div> <div>75%</div> <div>10%</div> <div>••</div> <div>13%</div> </div>
1	B	494	<div> <div>2%</div> <div>76%</div> <div>9%</div> <div>•</div> <div>14%</div> </div>

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	EDO	A	501	-	-	-	X
2	EDO	A	502	-	-	-	X
2	EDO	A	503	-	-	-	X
2	EDO	A	505	-	-	-	X
2	EDO	B	501	-	-	-	X
2	EDO	B	502	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7305 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 Carboxylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	430	Total	C	N	O	S	0	2	0
			3256	2049	591	609	7			
1	B	426	Total	C	N	O	S	0	2	0
			3235	2040	585	603	7			

There are 60 discrepancies between the modelled and reference sequences:

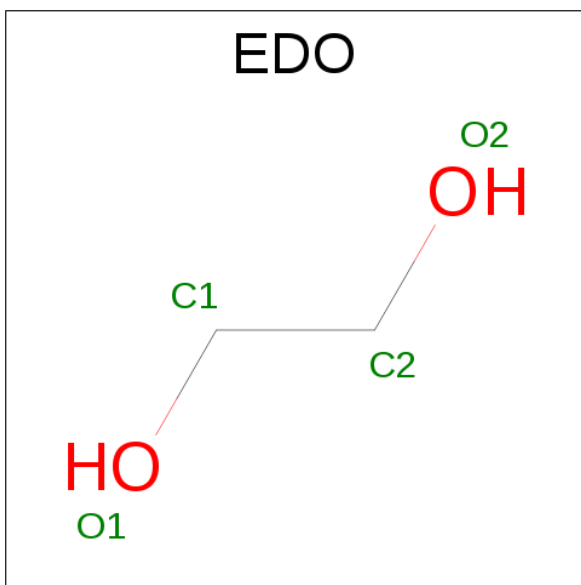
Chain	Residue	Modelled	Actual	Comment	Reference
A	-29	MET	-	initiating methionine	UNP Q9RYK2
A	-28	ALA	-	expression tag	UNP Q9RYK2
A	-27	HIS	-	expression tag	UNP Q9RYK2
A	-26	HIS	-	expression tag	UNP Q9RYK2
A	-25	HIS	-	expression tag	UNP Q9RYK2
A	-24	HIS	-	expression tag	UNP Q9RYK2
A	-23	HIS	-	expression tag	UNP Q9RYK2
A	-22	HIS	-	expression tag	UNP Q9RYK2
A	-21	HIS	-	expression tag	UNP Q9RYK2
A	-20	HIS	-	expression tag	UNP Q9RYK2
A	-19	HIS	-	expression tag	UNP Q9RYK2
A	-18	HIS	-	expression tag	UNP Q9RYK2
A	-17	LYS	-	expression tag	UNP Q9RYK2
A	-16	LEU	-	expression tag	UNP Q9RYK2
A	-15	THR	-	expression tag	UNP Q9RYK2
A	-14	SER	-	expression tag	UNP Q9RYK2
A	-13	LEU	-	expression tag	UNP Q9RYK2
A	-12	TYR	-	expression tag	UNP Q9RYK2
A	-11	LYS	-	expression tag	UNP Q9RYK2
A	-10	LYS	-	expression tag	UNP Q9RYK2
A	-9	ALA	-	expression tag	UNP Q9RYK2
A	-8	GLY	-	expression tag	UNP Q9RYK2
A	-7	LEU	-	expression tag	UNP Q9RYK2
A	-6	GLU	-	expression tag	UNP Q9RYK2
A	-5	ASN	-	expression tag	UNP Q9RYK2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	LEU	-	expression tag	UNP Q9RYK2
A	-3	TYR	-	expression tag	UNP Q9RYK2
A	-2	PHE	-	expression tag	UNP Q9RYK2
A	-1	GLN	-	expression tag	UNP Q9RYK2
A	0	GLY	-	expression tag	UNP Q9RYK2
B	-29	MET	-	initiating methionine	UNP Q9RYK2
B	-28	ALA	-	expression tag	UNP Q9RYK2
B	-27	HIS	-	expression tag	UNP Q9RYK2
B	-26	HIS	-	expression tag	UNP Q9RYK2
B	-25	HIS	-	expression tag	UNP Q9RYK2
B	-24	HIS	-	expression tag	UNP Q9RYK2
B	-23	HIS	-	expression tag	UNP Q9RYK2
B	-22	HIS	-	expression tag	UNP Q9RYK2
B	-21	HIS	-	expression tag	UNP Q9RYK2
B	-20	HIS	-	expression tag	UNP Q9RYK2
B	-19	HIS	-	expression tag	UNP Q9RYK2
B	-18	HIS	-	expression tag	UNP Q9RYK2
B	-17	LYS	-	expression tag	UNP Q9RYK2
B	-16	LEU	-	expression tag	UNP Q9RYK2
B	-15	THR	-	expression tag	UNP Q9RYK2
B	-14	SER	-	expression tag	UNP Q9RYK2
B	-13	LEU	-	expression tag	UNP Q9RYK2
B	-12	TYR	-	expression tag	UNP Q9RYK2
B	-11	LYS	-	expression tag	UNP Q9RYK2
B	-10	LYS	-	expression tag	UNP Q9RYK2
B	-9	ALA	-	expression tag	UNP Q9RYK2
B	-8	GLY	-	expression tag	UNP Q9RYK2
B	-7	LEU	-	expression tag	UNP Q9RYK2
B	-6	GLU	-	expression tag	UNP Q9RYK2
B	-5	ASN	-	expression tag	UNP Q9RYK2
B	-4	LEU	-	expression tag	UNP Q9RYK2
B	-3	TYR	-	expression tag	UNP Q9RYK2
B	-2	PHE	-	expression tag	UNP Q9RYK2
B	-1	GLN	-	expression tag	UNP Q9RYK2
B	0	GLY	-	expression tag	UNP Q9RYK2

- Molecule 2 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



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

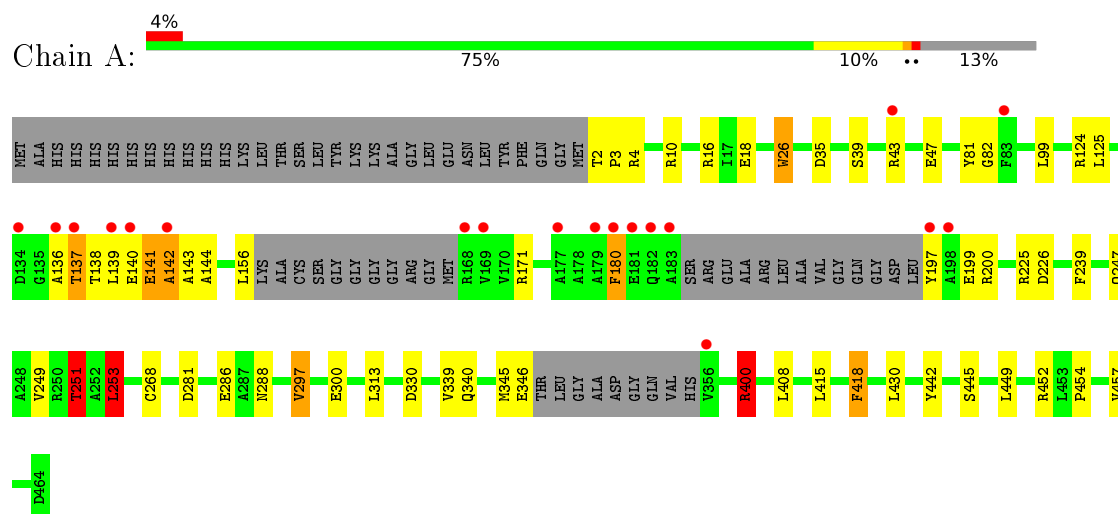
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	412	Total	O	0	0
			412	412		
3	B	374	Total	O	0	0
			374	374		

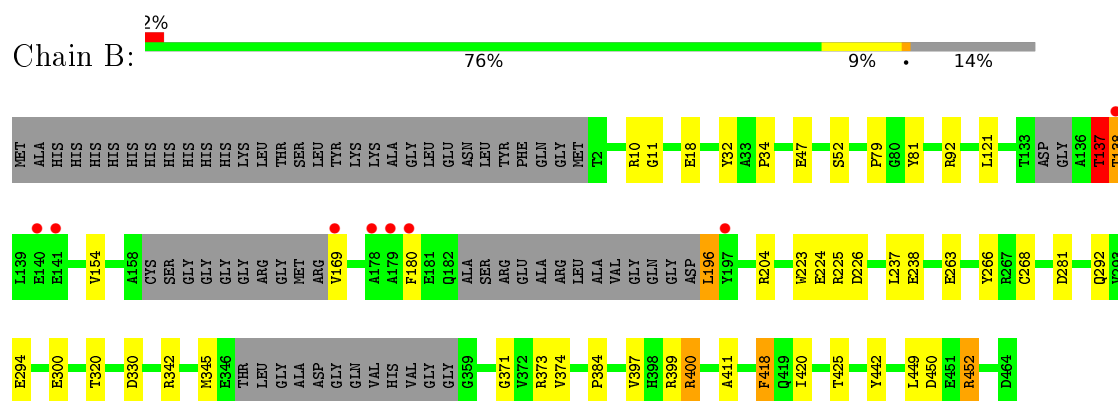
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: Carboxylase



• Molecule 1: Carboxylase



4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	138.10 Å 138.10 Å 97.87 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	75.74 – 1.70 75.74 – 1.70	Depositor EDS
% Data completeness (in resolution range)	99.9 (75.74-1.70) 99.7 (75.74-1.70)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.53 (at 1.70 Å)	Xtriage
Refinement program	REFMAC 5.8.0131	Depositor
R, R_{free}	0.149 , 0.184 0.178 , 0.216	Depositor DCC
R_{free} test set	4741 reflections (4.25%)	DCC
Wilson B-factor (Å ²)	26.0	Xtriage
Anisotropy	0.073	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 43.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	0.000 for h,-h-k,-l	Xtriage
Reported twinning fraction	0.694 for H, K, L 0.306 for K, H, -L	Depositor
Outliers	0 of 116229 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7305	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.87% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical 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: EDO

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	1.27	8/3323 (0.2%)	1.26	23/4534 (0.5%)
1	B	1.25	10/3301 (0.3%)	1.25	23/4503 (0.5%)
All	All	1.26	18/6624 (0.3%)	1.25	46/9037 (0.5%)

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 (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	294	GLU	CG-CD	7.50	1.63	1.51
1	B	266	TYR	CZ-OH	-7.04	1.25	1.37
1	A	251	THR	CB-CG2	-6.72	1.30	1.52
1	A	445	SER	CA-CB	6.51	1.62	1.52
1	A	300	GLU	CG-CD	-6.38	1.42	1.51
1	B	18	GLU	CD-OE2	6.08	1.32	1.25
1	A	239	PHE	CB-CG	-6.04	1.41	1.51
1	B	300	GLU	CD-OE1	5.81	1.32	1.25
1	A	26	TRP	CB-CG	5.79	1.60	1.50
1	B	11	GLY	C-O	5.55	1.32	1.23
1	A	39	SER	CB-OG	5.54	1.49	1.42
1	B	223	TRP	CE2-CZ2	5.43	1.49	1.39
1	B	320	THR	C-O	5.43	1.33	1.23
1	A	199	GLU	CD-OE2	5.40	1.31	1.25
1	B	32	TYR	CG-CD2	5.32	1.46	1.39
1	B	292	GLN	C-O	5.25	1.33	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	82	GLY	C-O	5.07	1.31	1.23
1	B	224	GLU	CD-OE2	5.07	1.31	1.25

All (46) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	225	ARG	NE-CZ-NH1	10.95	125.78	120.30
1	A	330	ASP	CB-CG-OD1	9.83	127.14	118.30
1	A	452	ARG	NE-CZ-NH2	8.33	124.47	120.30
1	A	400	ARG	NE-CZ-NH2	8.06	124.33	120.30
1	B	373	ARG	NE-CZ-NH2	-7.40	116.60	120.30
1	A	330	ASP	CB-CG-OD2	-6.99	112.01	118.30
1	B	224	GLU	OE1-CD-OE2	6.88	131.55	123.30
1	B	204	ARG	NE-CZ-NH1	-6.84	116.88	120.30
1	A	408	LEU	CB-CG-CD1	-6.60	99.78	111.00
1	B	374	VAL	CG1-CB-CG2	-6.49	100.51	110.90
1	B	400	ARG	NE-CZ-NH2	6.37	123.48	120.30
1	A	339	VAL	CG1-CB-CG2	-6.33	100.78	110.90
1	A	200	ARG	NE-CZ-NH2	-6.29	117.15	120.30
1	A	171	ARG	NE-CZ-NH1	6.27	123.44	120.30
1	A	418	PHE	CB-CG-CD2	6.18	125.12	120.80
1	A	253	LEU	CA-CB-CG	6.16	129.47	115.30
1	A	313	LEU	CB-CG-CD2	6.14	121.44	111.00
1	B	226	ASP	CB-CG-OD2	-6.07	112.84	118.30
1	B	330	ASP	CB-CG-OD1	6.07	123.76	118.30
1	B	330	ASP	CB-CG-OD2	-6.06	112.85	118.30
1	B	400	ARG	NE-CZ-NH1	-6.04	117.28	120.30
1	A	400	ARG	NE-CZ-NH1	-6.02	117.29	120.30
1	A	226	ASP	CB-CG-OD2	-5.91	112.98	118.30
1	A	253	LEU	CB-CG-CD1	5.90	121.03	111.00
1	A	418	PHE	CB-CG-CD1	-5.82	116.73	120.80
1	B	281	ASP	CB-CG-OD1	5.76	123.48	118.30
1	A	99	LEU	CB-CG-CD1	5.67	120.64	111.00
1	B	452	ARG	NE-CZ-NH1	5.50	123.05	120.30
1	B	196	LEU	CA-CB-CG	5.47	127.88	115.30
1	B	204	ARG	NE-CZ-NH2	5.46	123.03	120.30
1	B	342	ARG	NE-CZ-NH1	5.43	123.02	120.30
1	A	16	ARG	NE-CZ-NH1	5.40	123.00	120.30
1	B	418	PHE	CB-CG-CD2	5.37	124.56	120.80
1	A	281	ASP	CB-CG-OD1	5.32	123.08	118.30
1	B	92	ARG	NE-CZ-NH1	5.31	122.96	120.30
1	A	430	LEU	CB-CG-CD2	5.30	120.01	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	452	ARG	NE-CZ-NH2	-5.24	117.68	120.30
1	B	399	ARG	NE-CZ-NH1	5.23	122.92	120.30
1	B	345	MET	CG-SD-CE	-5.23	91.83	100.20
1	A	171	ARG	NE-CZ-NH2	-5.21	117.69	120.30
1	B	92	ARG	NE-CZ-NH2	-5.11	117.75	120.30
1	A	225	ARG	NE-CZ-NH1	5.10	122.85	120.30
1	B	300	GLU	OE1-CD-OE2	-5.08	117.21	123.30
1	B	450	ASP	CB-CG-OD2	5.06	122.85	118.30
1	A	415	LEU	CB-CG-CD1	-5.03	102.45	111.00
1	A	35	ASP	CB-CG-OD1	5.02	122.82	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	18	GLU	Mainchain

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	3256	0	3209	32	0
1	B	3235	0	3197	11	0
2	A	20	0	30	2	0
2	B	8	0	12	1	0
3	A	412	0	0	3	1
3	B	374	0	0	3	0
All	All	7305	0	6448	45	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All (45) 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:A:139:LEU:HD22	1:A:180:PHE:CE1	1.96	0.99

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:141:GLU:O	1:A:144:ALA:N	2.01	0.93
1:A:138:THR:OG1	1:A:141:GLU:HG3	1.76	0.85
1:A:138:THR:OG1	1:A:141:GLU:CG	2.33	0.76
2:B:501:EDO:H11	3:B:685:HOH:O	1.85	0.76
1:B:137:THR:OG1	1:B:138:THR:N	2.22	0.70
1:A:141:GLU:C	1:A:141:GLU:OE1	2.37	0.62
1:A:141:GLU:O	1:A:142:ALA:C	2.39	0.62
1:A:141:GLU:OE1	1:A:142:ALA:N	2.34	0.60
1:B:397[A]:VAL:HG21	1:B:411:ALA:HA	1.84	0.60
1:A:137:THR:HG23	1:A:139:LEU:HG	1.85	0.59
1:A:180:PHE:HD2	1:A:180:PHE:C	2.06	0.59
1:A:180:PHE:CD2	1:A:180:PHE:C	2.77	0.58
1:A:400:ARG:HH21	1:A:400:ARG:HG2	1.68	0.56
1:A:124:ARG:NH1	2:A:504:EDO:O1	2.41	0.54
1:A:156:LEU:O	1:A:197:TYR:N	2.41	0.53
1:A:4:ARG:NH2	1:A:47:GLU:OE1	2.37	0.52
1:A:141:GLU:O	1:A:143:ALA:N	2.43	0.51
1:A:286:GLU:HG2	1:A:288:ASN:HD21	1.76	0.51
1:B:180:PHE:O	1:B:196:LEU:HD11	2.11	0.50
1:A:139:LEU:HD22	1:A:180:PHE:CZ	2.46	0.49
1:A:124:ARG:NH2	3:A:608:HOH:O	2.45	0.49
1:B:452:ARG:NH1	3:B:611:HOH:O	2.45	0.48
1:A:136:ALA:O	1:A:137:THR:HB	2.14	0.48
1:A:249:VAL:HG12	1:A:253:LEU:HD22	1.96	0.48
1:B:137:THR:HG22	1:B:196:LEU:O	2.14	0.47
1:A:247:GLN:O	1:A:251:THR:HG23	2.14	0.47
1:B:371:GLY:O	1:B:397[B]:VAL:HA	2.14	0.47
1:B:420:ILE:HD12	1:B:425:THR:HG21	1.97	0.47
1:A:297[A]:VAL:HG13	1:A:340:GLN:HB2	1.97	0.46
1:A:141:GLU:C	1:A:143:ALA:N	2.69	0.46
1:A:3:PRO:HG2	1:A:26:TRP:CD2	2.51	0.45
1:A:454:PRO:HA	1:A:457:VAL:HG22	1.98	0.45
1:A:400:ARG:CG	1:A:400:ARG:HH21	2.31	0.44
2:A:503:EDO:O2	3:A:602:HOH:O	2.21	0.43
1:B:449:LEU:C	1:B:449:LEU:HD23	2.39	0.43
1:B:121:LEU:HD11	1:B:263:GLU:HG2	2.00	0.43
1:A:449:LEU:C	1:A:449:LEU:HD23	2.40	0.43
1:A:449:LEU:O	1:A:449:LEU:HD23	2.18	0.42
1:B:52:SER:HB3	3:B:941:HOH:O	2.18	0.42
1:A:345:MET:O	1:A:346:GLU:HB3	2.20	0.42
1:A:286:GLU:HG2	1:A:288:ASN:ND2	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:237:LEU:C	1:B:238:GLU:HG3	2.39	0.40
1:A:43:ARG:NE	3:A:612:HOH:O	2.53	0.40
1:A:136:ALA:O	1:A:137:THR:CB	2.69	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:731:HOH:O	3:A:932:HOH:O[6_554]	2.00	0.20

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	424/494 (86%)	409 (96%)	12 (3%)	3 (1%)	26	9
1	B	418/494 (85%)	406 (97%)	9 (2%)	3 (1%)	26	9
All	All	842/988 (85%)	815 (97%)	21 (2%)	6 (1%)	26	9

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	137	THR
1	A	142	ALA
1	A	268	CYS
1	B	268	CYS
1	B	137	THR
1	B	138	THR

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	326/371 (88%)	312 (96%)	14 (4%)	35	14
1	B	325/371 (88%)	313 (96%)	12 (4%)	41	18
All	All	651/742 (88%)	625 (96%)	26 (4%)	39	16

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	THR
1	A	10	ARG
1	A	81	TYR
1	A	125	LEU
1	A	140	GLU
1	A	141	GLU
1	A	180	PHE
1	A	251	THR
1	A	253	LEU
1	A	297[A]	VAL
1	A	297[B]	VAL
1	A	400	ARG
1	A	418	PHE
1	A	442	TYR
1	B	10	ARG
1	B	34	PRO
1	B	47	GLU
1	B	79	PRO
1	B	81	TYR
1	B	137	THR
1	B	154	VAL
1	B	169	VAL
1	B	384	PRO
1	B	400	ARG
1	B	418	PHE
1	B	442	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

7 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	EDO	A	501	-	3,3,3	0.35	0	2,2,2	0.48	0
2	EDO	A	502	-	3,3,3	0.44	0	2,2,2	0.58	0
2	EDO	A	503	-	3,3,3	0.60	0	2,2,2	0.67	0
2	EDO	A	504	-	3,3,3	0.28	0	2,2,2	0.42	0
2	EDO	A	505	-	3,3,3	0.42	0	2,2,2	1.07	0
2	EDO	B	501	-	3,3,3	0.45	0	2,2,2	0.64	0
2	EDO	B	502	-	3,3,3	0.85	0	2,2,2	0.53	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	EDO	A	501	-	-	0/1/1/1	0/0/0/0
2	EDO	A	502	-	-	0/1/1/1	0/0/0/0
2	EDO	A	503	-	-	0/1/1/1	0/0/0/0
2	EDO	A	504	-	-	0/1/1/1	0/0/0/0
2	EDO	A	505	-	-	0/1/1/1	0/0/0/0
2	EDO	B	501	-	-	0/1/1/1	0/0/0/0
2	EDO	B	502	-	-	0/1/1/1	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	503	EDO	1	0
2	A	504	EDO	1	0
2	B	501	EDO	1	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	430/494 (87%)	-0.03	19 (4%) 38 42	19, 29, 66, 93	0
1	B	426/494 (86%)	-0.08	8 (1%) 70 74	19, 29, 63, 106	0
All	All	856/988 (86%)	-0.06	27 (3%) 51 55	19, 29, 64, 106	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	142	ALA	5.1
1	A	134	ASP	4.3
1	A	169	VAL	4.0
1	B	169	VAL	3.9
1	A	198	ALA	3.8
1	A	197	TYR	3.7
1	B	179	ALA	3.3
1	A	180	PHE	3.3
1	A	139	LEU	3.3
1	A	182	GLN	3.2
1	B	197	TYR	3.0
1	A	136	ALA	2.9
1	B	180	PHE	2.8
1	A	168	ARG	2.7
1	B	138	THR	2.6
1	A	179	ALA	2.5
1	B	141	GLU	2.5
1	A	137	THR	2.4
1	A	140	GLU	2.4
1	A	356	VAL	2.4
1	B	140	GLU	2.3
1	A	177	ALA	2.2
1	A	181	GLU	2.2
1	A	43	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	83	PHE	2.1
1	A	183	ALA	2.1
1	B	178	ALA	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	EDO	A	502	4/4	0.84	0.19	8.59	38,45,49,55	0
2	EDO	A	503	4/4	0.76	0.25	6.95	45,53,57,67	0
2	EDO	A	505	4/4	0.77	0.16	6.55	38,41,43,50	0
2	EDO	B	502	4/4	0.88	0.15	3.50	27,28,30,36	0
2	EDO	A	501	4/4	0.92	0.11	3.03	39,40,43,46	0
2	EDO	B	501	4/4	0.91	0.13	2.82	37,37,46,47	0
2	EDO	A	504	4/4	0.83	0.18	-	53,56,57,58	0

6.5 Other polymers [i](#)

There are no such residues in this entry.