



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 12:42 AM UTC

PDB ID : 1PN2 / pdb_00001pn2
Title : Crystal structure analysis of the selenomethionine labelled 2-enoyl-CoA hydratase 2 domain of Candida tropicalis multifunctional enzyme type 2
Authors : Koski, M.K.; Haapalainen, A.M.; Hiltunen, J.K.; Glumoff, T.
Deposited on : 2003-06-12
Resolution : 1.95 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

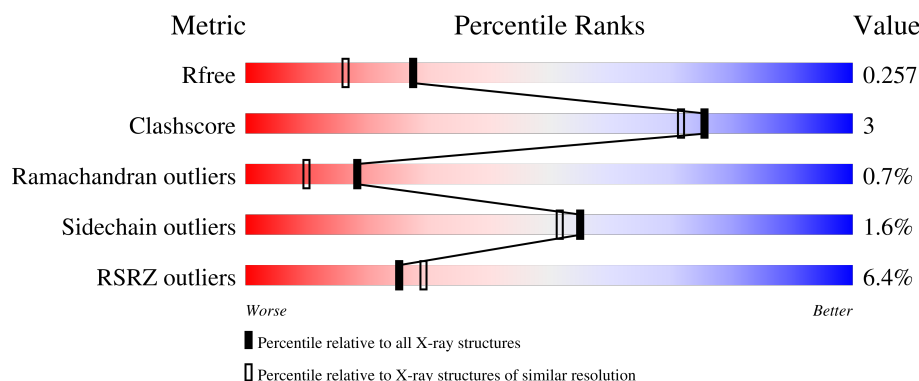
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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	280	 8% 84% 12% . .
1	B	280	 2% 85% 10% . .
1	C	280	 7% 82% 11% . 5%
1	D	280	 6% 89% 8% .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9374 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 Peroxisomal hydratase-dehydrogenase-epimerase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	272	Total	C	N	O	S	Se	98	0	0
			2153	1386	367	395	2	3			
1	B	270	Total	C	N	O	S	Se	51	0	0
			2141	1380	365	391	2	3			
1	C	267	Total	C	N	O	S	Se	36	0	0
			2117	1366	360	386	2	3			
1	D	272	Total	C	N	O	S	Se	32	0	0
			2153	1386	367	395	2	3			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	GLU	engineered mutation	UNP P22414
A	69	MSE	MET	modified residue	UNP P22414
A	206	MSE	MET	modified residue	UNP P22414
A	222	MSE	MET	modified residue	UNP P22414
B	1	MSE	GLU	engineered mutation	UNP P22414
B	69	MSE	MET	modified residue	UNP P22414
B	206	MSE	MET	modified residue	UNP P22414
B	222	MSE	MET	modified residue	UNP P22414
C	1	MSE	GLU	engineered mutation	UNP P22414
C	69	MSE	MET	modified residue	UNP P22414
C	206	MSE	MET	modified residue	UNP P22414
C	222	MSE	MET	modified residue	UNP P22414
D	1	MSE	GLU	engineered mutation	UNP P22414
D	69	MSE	MET	modified residue	UNP P22414
D	206	MSE	MET	modified residue	UNP P22414
D	222	MSE	MET	modified residue	UNP P22414

- Molecule 2 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



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

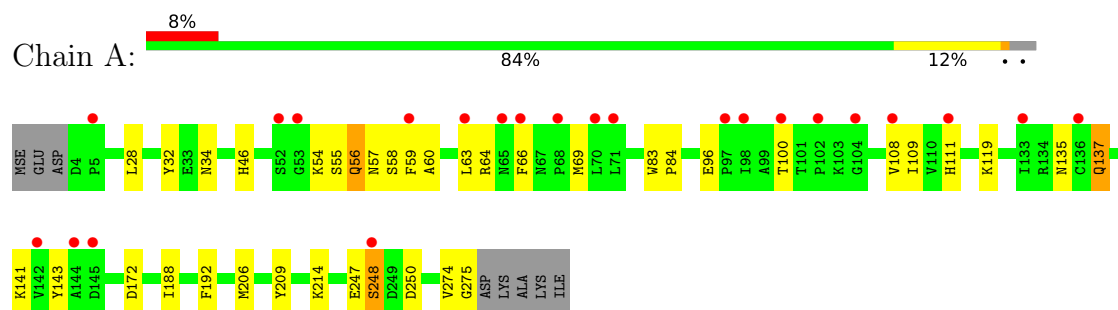
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	153	Total	O	0	0
			153	153		
3	B	147	Total	O	0	0
			147	147		
3	C	236	Total	O	0	0
			236	236		
3	D	254	Total	O	0	0
			254	254		

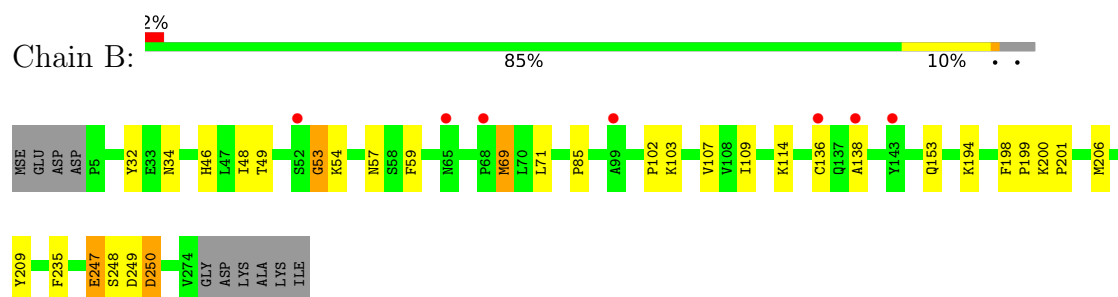
3 Residue-property plots [i](#)

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.

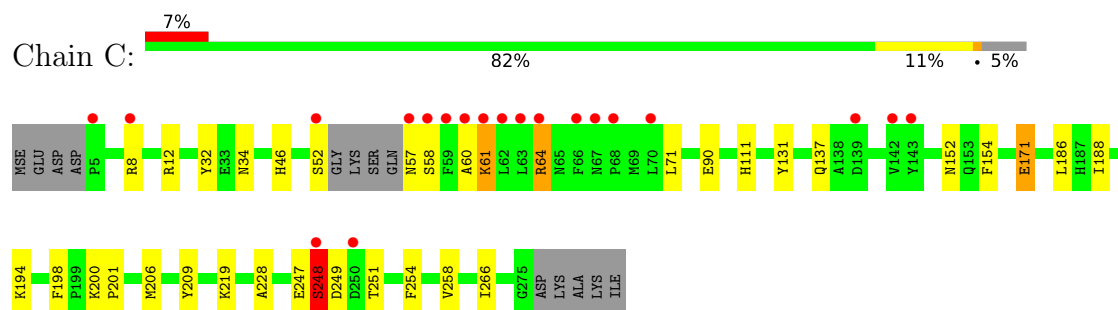
- Molecule 1: Peroxisomal hydratase-dehydrogenase-epimerase



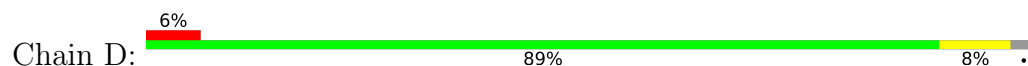
- Molecule 1: Peroxisomal hydratase-dehydrogenase-epimerase

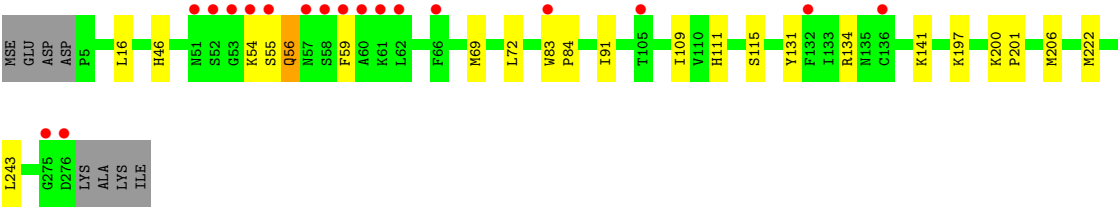


- Molecule 1: Peroxisomal hydratase-dehydrogenase-epimerase



- Molecule 1: Peroxisomal hydratase-dehydrogenase-epimerase





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	178.80Å 60.65Å 131.12Å 90.00° 94.57° 90.00°	Depositor
Resolution (Å)	30.00 – 1.95 30.00 – 1.95	Depositor EDS
% Data completeness (in resolution range)	95.2 (30.00-1.95) 95.2 (30.00-1.95)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.28 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.1	Depositor
R, R_{free}	0.179 , 0.214 0.236 , 0.257	Depositor DCC
R_{free} test set	9761 reflections (10.00%)	wwPDB-VP
Wilson B-factor (Å ²)	26.4	Xtriage
Anisotropy	0.619	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 44.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	9374	wwPDB-VP
Average B, all atoms (Å ²)	11.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.11% 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: 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	1.12	8/2207 (0.4%)	1.29	19/2996 (0.6%)
1	B	0.83	3/2195 (0.1%)	1.00	5/2979 (0.2%)
1	C	1.16	8/2170 (0.4%)	1.07	9/2945 (0.3%)
1	D	1.02	5/2207 (0.2%)	1.04	6/2995 (0.2%)
All	All	1.04	24/8779 (0.3%)	1.11	39/11915 (0.3%)

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
1	B	0	1
1	C	0	1
All	All	0	3

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	61	LYS	CA-CB	-18.76	1.21	1.53
1	A	56	GLN	C-N	-16.46	1.09	1.33
1	A	60	ALA	C-N	-14.62	1.12	1.33
1	A	135	ASN	CA-CB	-13.71	1.32	1.53
1	A	96	GLU	CD-OE1	-12.61	1.01	1.25
1	C	60	ALA	CA-CB	-12.12	1.31	1.53
1	D	134	ARG	NE-CZ	-10.69	1.21	1.33
1	D	222	MSE	SE-CE	-10.07	1.65	1.95
1	C	171	GLU	CD-OE2	-8.12	1.09	1.25
1	B	114	LYS	CE-NZ	-7.72	1.26	1.49
1	A	100	THR	CA-CB	-7.49	1.38	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	249	ASP	CA-CB	-7.22	1.40	1.53
1	A	109	ILE	CB-CG1	6.96	1.67	1.53
1	A	137	GLN	CA-CB	6.96	1.65	1.53
1	C	249	ASP	CG-OD1	6.62	1.38	1.25
1	D	54	LYS	CA-CB	6.41	1.64	1.53
1	A	209	TYR	CA-CB	-6.32	1.43	1.53
1	C	171	GLU	CD-OE1	6.20	1.37	1.25
1	C	209	TYR	CA-CB	-5.84	1.44	1.53
1	B	103	LYS	CB-CG	-5.82	1.34	1.52
1	C	52	SER	CA-CB	-5.70	1.42	1.53
1	D	141	LYS	CD-CE	-5.68	1.35	1.52
1	D	69	MSE	SE-CE	5.15	2.10	1.95
1	C	258	VAL	CA-C	5.02	1.59	1.52

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	60	ALA	CA-C-N	21.77	157.70	120.68
1	A	60	ALA	C-N-CA	21.77	157.70	120.68
1	A	56	GLN	O-C-N	-14.55	100.96	122.20
1	A	135	ASN	CB-CA-C	12.06	130.56	111.51
1	A	60	ALA	O-C-N	-11.47	107.33	122.59
1	D	56	GLN	OE1-CD-NE2	-10.64	111.95	122.60
1	C	57	ASN	CB-CA-C	-10.25	90.63	110.10
1	C	249	ASP	N-CA-C	-9.04	99.25	110.65
1	C	60	ALA	N-CA-CB	-8.50	95.52	110.39
1	A	56	GLN	CA-C-N	8.01	135.16	121.14
1	A	56	GLN	C-N-CA	8.01	135.16	121.14
1	A	135	ASN	CA-CB-CG	7.90	120.50	112.60
1	C	61	LYS	CB-CA-C	7.87	123.23	110.09
1	B	53	GLY	N-CA-C	7.53	131.02	113.18
1	D	54	LYS	CB-CA-C	-7.50	95.50	110.42
1	A	96	GLU	OE1-CD-OE2	7.20	140.18	122.90
1	A	96	GLU	CG-CD-OE2	-7.09	102.10	118.40
1	D	56	GLN	CG-CD-NE2	6.81	126.61	116.40
1	A	143	TYR	CB-CA-C	-6.80	97.65	109.07
1	B	85	PRO	O-C-N	6.70	124.30	121.15
1	A	119	LYS	CA-CB-CG	6.49	127.08	114.10
1	A	108	VAL	N-CA-CB	6.40	119.41	111.41
1	A	108	VAL	CB-CA-C	-6.26	102.54	111.31
1	C	60	ALA	CB-CA-C	6.08	122.58	110.17
1	A	141	LYS	N-CA-C	6.01	118.31	109.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	56	GLN	CB-CA-C	-5.94	99.76	109.56
1	C	209	TYR	CA-CB-CG	5.88	124.49	113.90
1	B	103	LYS	CA-CB-CG	5.88	125.85	114.10
1	C	61	LYS	N-CA-CB	5.77	119.04	110.49
1	A	57	ASN	CB-CA-C	-5.73	98.01	109.99
1	A	109	ILE	CA-CB-CG2	5.67	120.14	110.50
1	C	249	ASP	CB-CG-OD2	5.63	131.35	118.40
1	B	209	TYR	CA-CB-CG	5.58	123.95	113.90
1	B	57	ASN	CA-CB-CG	5.36	117.96	112.60
1	D	134	ARG	CD-NE-CZ	5.33	131.86	124.40
1	C	254	PHE	N-CA-C	5.27	116.81	108.96
1	D	141	LYS	CG-CD-CE	5.26	123.39	111.30
1	D	56	GLN	CB-CG-CD	5.25	121.52	112.60
1	A	55	SER	CB-CA-C	-5.07	99.49	109.67

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	56	GLN	Mainchain
1	B	250	ASP	Peptide
1	C	248	SER	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	2153	0	2131	12	0
1	B	2141	0	2127	14	0
1	C	2117	0	2100	15	0
1	D	2153	0	2134	9	0
2	A	4	0	6	1	0
2	C	8	0	12	0	0
2	D	8	0	12	0	0
3	A	153	0	0	1	0
3	B	147	0	0	1	0
3	C	236	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	254	0	0	0	0
All	All	9374	0	8522	49	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 3.

All (49) 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:54:LYS:O	1:A:58:SER:HB3	1.89	0.73
1:B:247:GLU:OE1	1:B:247:GLU:HA	1.91	0.69
1:C:58:SER:HB2	1:C:111:HIS:HE1	1.63	0.63
1:C:64:ARG:HB2	1:C:137:GLN:HB3	1.84	0.60
1:A:247:GLU:O	1:A:248:SER:HB3	2.02	0.59
1:C:12:ARG:HH21	1:D:16:LEU:HD23	1.69	0.58
1:C:32:TYR:CZ	1:C:34:ASN:HB2	2.40	0.56
1:B:69:MSE:HB2	3:B:402:HOH:O	2.05	0.56
1:D:46:HIS:CG	1:D:206:MSE:HE2	2.41	0.55
1:A:214:LYS:NZ	3:A:927:HOH:O	2.31	0.55
1:D:83:TRP:HA	1:D:84:PRO:C	2.32	0.54
1:A:32:TYR:CZ	1:A:34:ASN:HB2	2.42	0.54
1:B:59:PHE:HB2	1:B:109:ILE:CD1	2.39	0.53
1:D:83:TRP:HB2	1:D:84:PRO:HA	1.94	0.50
1:B:48:ILE:HD12	1:B:49:THR:N	2.27	0.50
1:B:107:VAL:HG12	1:B:136:CYS:HB3	1.93	0.50
1:A:69:MSE:HE2	1:A:192:PHE:HB3	1.95	0.49
1:A:172:ASP:OD2	2:A:805:EDO:O2	2.30	0.49
1:B:200:LYS:HB2	1:B:201:PRO:CD	2.43	0.48
1:A:46:HIS:CG	1:A:206:MSE:HE2	2.48	0.48
1:B:59:PHE:HB2	1:B:109:ILE:HD11	1.95	0.48
1:D:59:PHE:HB3	1:D:111:HIS:CD2	2.48	0.48
1:B:194:LYS:HA	1:B:198:PHE:O	2.13	0.47
1:C:46:HIS:CG	1:C:206:MSE:HE2	2.50	0.47
1:A:64:ARG:HB2	1:A:137:GLN:O	2.16	0.46
1:B:46:HIS:CG	1:B:206:MSE:HE2	2.50	0.46
1:C:152:ASN:HB3	1:C:154:PHE:CE2	2.51	0.46
1:D:109:ILE:HG23	1:D:131:TYR:HB2	1.98	0.45
1:C:58:SER:HB2	1:C:111:HIS:CE1	2.47	0.45
1:D:200:LYS:HB2	1:D:201:PRO:CD	2.48	0.44
1:C:219:LYS:NZ	3:C:979:HOH:O	2.48	0.44
1:A:63:LEU:HD23	1:A:66:PHE:CG	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:248:SER:HA	1:C:251:THR:H	1.83	0.43
1:B:71:LEU:C	1:B:71:LEU:HD23	2.43	0.43
1:C:8:ARG:NH1	1:C:90:GLU:OE2	2.52	0.43
1:C:71:LEU:HD11	1:C:131:TYR:HB3	2.00	0.42
1:B:102:PRO:HG3	1:B:138:ALA:O	2.19	0.42
1:D:91:ILE:CG2	1:D:115:SER:HB3	2.49	0.42
1:A:59:PHE:HB3	1:A:111:HIS:CD2	2.55	0.42
1:C:228:ALA:HB1	1:C:266:ILE:HG23	2.01	0.42
1:A:83:TRP:CG	1:A:84:PRO:HA	2.55	0.42
1:A:274:VAL:HG12	1:A:275:GLY:N	2.35	0.42
1:B:200:LYS:HB2	1:B:201:PRO:HD2	2.02	0.42
1:C:247:GLU:O	1:C:248:SER:C	2.62	0.42
1:B:32:TYR:CZ	1:B:34:ASN:HB2	2.55	0.42
1:B:199:PRO:HG2	1:B:235:PHE:CZ	2.55	0.41
1:D:91:ILE:HG21	1:D:115:SER:HB3	2.03	0.41
1:C:194:LYS:HA	1:C:198:PHE:O	2.21	0.40
1:C:200:LYS:HB2	1:C:201:PRO:CD	2.51	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	270/280 (96%)	263 (97%)	5 (2%)	2 (1%)	18	10
1	B	268/280 (96%)	259 (97%)	7 (3%)	2 (1%)	18	10
1	C	263/280 (94%)	257 (98%)	4 (2%)	2 (1%)	16	8
1	D	270/280 (96%)	265 (98%)	4 (2%)	1 (0%)	30	21
All	All	1071/1120 (96%)	1044 (98%)	20 (2%)	7 (1%)	18	10

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	248	SER
1	B	53	GLY
1	B	54	LYS
1	C	248	SER
1	D	55	SER
1	C	188	ILE
1	A	188	ILE

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	233/236 (99%)	231 (99%)	2 (1%)	70	70
1	B	232/236 (98%)	227 (98%)	5 (2%)	45	40
1	C	229/236 (97%)	225 (98%)	4 (2%)	53	49
1	D	233/236 (99%)	229 (98%)	4 (2%)	53	49
All	All	927/944 (98%)	912 (98%)	15 (2%)	55	52

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	LEU
1	A	250	ASP
1	B	69	MSE
1	B	153	GLN
1	B	247	GLU
1	B	248	SER
1	B	250	ASP
1	C	61	LYS
1	C	64	ARG
1	C	171	GLU
1	C	186	LEU
1	D	56	GLN
1	D	72	LEU
1	D	197	LYS
1	D	243	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	73	HIS
1	B	111	HIS
1	C	111	HIS
1	C	140	ASN
1	C	191	ASN
1	D	111	HIS
1	D	135	ASN
1	D	140	ASN

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](#)

5 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	C	803	-	3,3,3	0.18	0	2,2,2	0.50	0
2	EDO	D	804	-	3,3,3	0.29	0	2,2,2	0.73	0
2	EDO	A	805	-	3,3,3	0.37	0	2,2,2	0.03	0
2	EDO	C	801	-	3,3,3	0.24	0	2,2,2	1.11	0
2	EDO	D	802	-	3,3,3	0.44	0	2,2,2	0.23	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	C	803	-	-	0/1/1/1	-
2	EDO	D	804	-	-	1/1/1/1	-
2	EDO	A	805	-	-	1/1/1/1	-
2	EDO	C	801	-	-	0/1/1/1	-
2	EDO	D	802	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	805	EDO	O1-C1-C2-O2
2	D	804	EDO	O1-C1-C2-O2
2	D	802	EDO	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	805	EDO	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	2

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	60:ALA	C	61:LYS	N	1.12
1	A	56:GLN	C	57:ASN	N	1.09

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	267/280 (95%)	0.44	23 (8%) 16 19	2, 6, 21, 25	27 (10%)
1	B	267/280 (95%)	0.16	7 (2%) 57 64	2, 5, 20, 37	16 (5%)
1	C	264/280 (94%)	0.24	20 (7%) 20 22	2, 6, 24, 39	13 (4%)
1	D	269/280 (96%)	0.24	18 (6%) 24 27	2, 6, 21, 43	10 (3%)
All	All	1067/1120 (95%)	0.27	68 (6%) 25 29	2, 6, 21, 43	66 (6%)

All (68) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	54	LYS	5.5
1	A	65	ASN	4.7
1	C	59	PHE	4.6
1	D	53	GLY	4.6
1	C	66	PHE	4.4
1	C	58	SER	4.2
1	D	51	ASN	3.8
1	A	136	CYS	3.7
1	D	132	PHE	3.6
1	C	62	LEU	3.5
1	C	248	SER	3.4
1	D	105	THR	3.3
1	A	71	LEU	3.1
1	A	100	THR	3.1
1	C	250	ASP	3.0
1	A	66	PHE	3.0
1	A	102	PRO	3.0
1	A	97	PRO	2.9
1	A	108	VAL	2.9
1	C	68	PRO	2.8
1	D	57	ASN	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	275	GLY	2.8
1	C	60	ALA	2.7
1	A	133	ILE	2.7
1	A	248	SER	2.7
1	C	67	ASN	2.7
1	A	145	ASP	2.6
1	D	276	ASP	2.5
1	C	57	ASN	2.5
1	C	139	ASP	2.5
1	A	63	LEU	2.5
1	C	63	LEU	2.5
1	B	52	SER	2.5
1	D	66	PHE	2.5
1	D	52	SER	2.4
1	D	55	SER	2.4
1	A	52	SER	2.4
1	A	70	LEU	2.4
1	B	143	TYR	2.4
1	A	144	ALA	2.4
1	A	53	GLY	2.4
1	D	59	PHE	2.3
1	B	68	PRO	2.3
1	C	143	TYR	2.3
1	D	60	ALA	2.3
1	C	52	SER	2.3
1	A	111	HIS	2.3
1	A	59	PHE	2.3
1	A	5	PRO	2.3
1	A	98	ILE	2.3
1	C	8	ARG	2.2
1	C	61	LYS	2.2
1	A	68	PRO	2.2
1	A	104	GLY	2.2
1	C	5	PRO	2.2
1	B	136	CYS	2.2
1	C	70	LEU	2.2
1	B	138	ALA	2.2
1	D	61	LYS	2.1
1	B	65	ASN	2.1
1	D	136	CYS	2.1
1	D	83	TRP	2.1
1	C	64	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	58	SER	2.0
1	A	142	VAL	2.0
1	D	62	LEU	2.0
1	B	99	ALA	2.0
1	C	142	VAL	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	EDO	A	805	4/4	0.70	0.24	44,47,49,49	0
2	EDO	D	804	4/4	0.90	0.16	34,37,40,44	0
2	EDO	D	802	4/4	0.91	0.18	21,22,26,26	0
2	EDO	C	801	4/4	0.94	0.14	29,33,34,35	0
2	EDO	C	803	4/4	0.94	0.18	34,39,40,44	0

6.5 Other polymers [i](#)

There are no such residues in this entry.