



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 05:00 PM UTC

PDB ID : 4IP2 / pdb_00004ip2
Title : Putative Aromatic Acid Decarboxylase
Authors : Schneider, G.; Brunner, K.; Izumi, A.; Jacewicz, A.
Deposited on : 2013-01-09
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

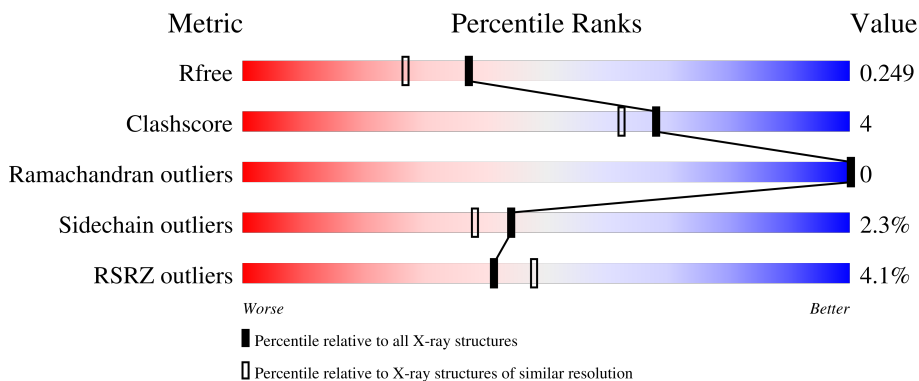
1 Overall quality at a glance

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	518	 87% 8% .
1	B	518	 89% 6% .
1	C	518	 10% 84% 12% ..

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 12473 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 Aromatic Acid Decarboxylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	499	Total 3900	C 2480	N 707	O 695	S 18	0	4	0
1	B	495	Total 3844	C 2439	N 701	O 688	S 16	0	1	0
1	C	502	Total 3913	C 2485	N 710	O 701	S 17	7	2	0

There are 66 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-21	MET	-	expression tag	UNP Q9I6N5
A	-20	HIS	-	expression tag	UNP Q9I6N5
A	-19	HIS	-	expression tag	UNP Q9I6N5
A	-18	HIS	-	expression tag	UNP Q9I6N5
A	-17	HIS	-	expression tag	UNP Q9I6N5
A	-16	HIS	-	expression tag	UNP Q9I6N5
A	-15	HIS	-	expression tag	UNP Q9I6N5
A	-14	SER	-	expression tag	UNP Q9I6N5
A	-13	SER	-	expression tag	UNP Q9I6N5
A	-12	GLY	-	expression tag	UNP Q9I6N5
A	-11	VAL	-	expression tag	UNP Q9I6N5
A	-10	ASP	-	expression tag	UNP Q9I6N5
A	-9	LEU	-	expression tag	UNP Q9I6N5
A	-8	GLY	-	expression tag	UNP Q9I6N5
A	-7	THR	-	expression tag	UNP Q9I6N5
A	-6	GLU	-	expression tag	UNP Q9I6N5
A	-5	ASN	-	expression tag	UNP Q9I6N5
A	-4	LEU	-	expression tag	UNP Q9I6N5
A	-3	TYR	-	expression tag	UNP Q9I6N5
A	-2	PHE	-	expression tag	UNP Q9I6N5
A	-1	GLN	-	expression tag	UNP Q9I6N5
A	0	SER	-	expression tag	UNP Q9I6N5
B	-21	MET	-	expression tag	UNP Q9I6N5

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-20	HIS	-	expression tag	UNP Q9I6N5
B	-19	HIS	-	expression tag	UNP Q9I6N5
B	-18	HIS	-	expression tag	UNP Q9I6N5
B	-17	HIS	-	expression tag	UNP Q9I6N5
B	-16	HIS	-	expression tag	UNP Q9I6N5
B	-15	HIS	-	expression tag	UNP Q9I6N5
B	-14	SER	-	expression tag	UNP Q9I6N5
B	-13	SER	-	expression tag	UNP Q9I6N5
B	-12	GLY	-	expression tag	UNP Q9I6N5
B	-11	VAL	-	expression tag	UNP Q9I6N5
B	-10	ASP	-	expression tag	UNP Q9I6N5
B	-9	LEU	-	expression tag	UNP Q9I6N5
B	-8	GLY	-	expression tag	UNP Q9I6N5
B	-7	THR	-	expression tag	UNP Q9I6N5
B	-6	GLU	-	expression tag	UNP Q9I6N5
B	-5	ASN	-	expression tag	UNP Q9I6N5
B	-4	LEU	-	expression tag	UNP Q9I6N5
B	-3	TYR	-	expression tag	UNP Q9I6N5
B	-2	PHE	-	expression tag	UNP Q9I6N5
B	-1	GLN	-	expression tag	UNP Q9I6N5
B	0	SER	-	expression tag	UNP Q9I6N5
C	-21	MET	-	expression tag	UNP Q9I6N5
C	-20	HIS	-	expression tag	UNP Q9I6N5
C	-19	HIS	-	expression tag	UNP Q9I6N5
C	-18	HIS	-	expression tag	UNP Q9I6N5
C	-17	HIS	-	expression tag	UNP Q9I6N5
C	-16	HIS	-	expression tag	UNP Q9I6N5
C	-15	HIS	-	expression tag	UNP Q9I6N5
C	-14	SER	-	expression tag	UNP Q9I6N5
C	-13	SER	-	expression tag	UNP Q9I6N5
C	-12	GLY	-	expression tag	UNP Q9I6N5
C	-11	VAL	-	expression tag	UNP Q9I6N5
C	-10	ASP	-	expression tag	UNP Q9I6N5
C	-9	LEU	-	expression tag	UNP Q9I6N5
C	-8	GLY	-	expression tag	UNP Q9I6N5
C	-7	THR	-	expression tag	UNP Q9I6N5
C	-6	GLU	-	expression tag	UNP Q9I6N5
C	-5	ASN	-	expression tag	UNP Q9I6N5
C	-4	LEU	-	expression tag	UNP Q9I6N5
C	-3	TYR	-	expression tag	UNP Q9I6N5
C	-2	PHE	-	expression tag	UNP Q9I6N5
C	-1	GLN	-	expression tag	UNP Q9I6N5

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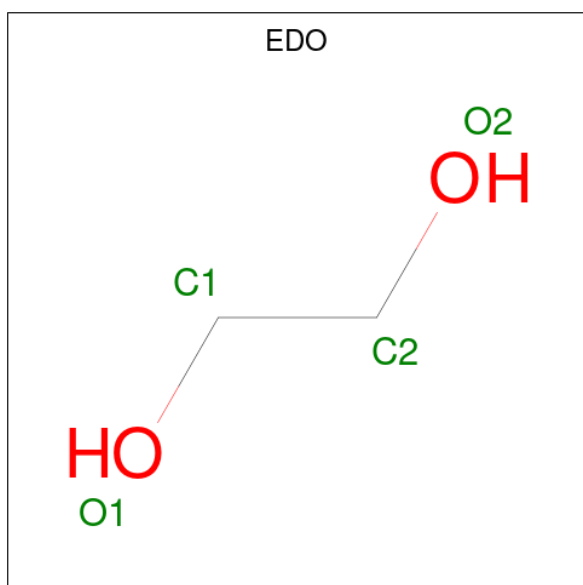
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Chain	Residue	Modelled	Actual	Comment	Reference
C	0	SER	-	expression tag	UNP Q9I6N5

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

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

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



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

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	C O	0	0
			6	3 3		

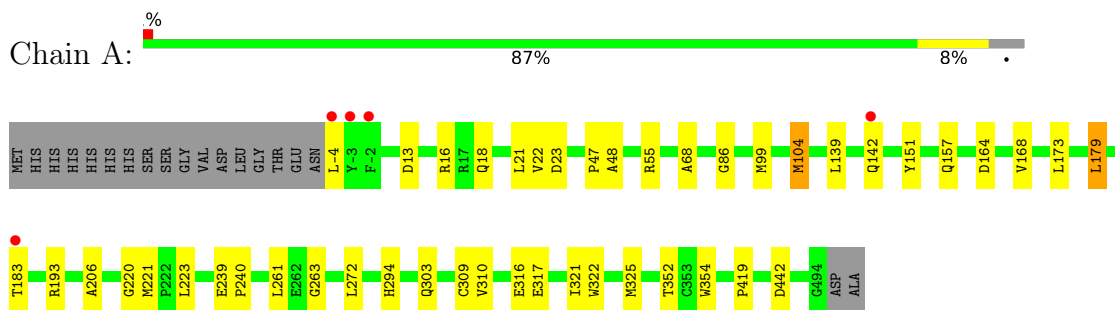
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	336	Total	O	0	0
			336	336		
5	B	290	Total	O	0	0
			290	290		
5	C	161	Total	O	0	0
			161	161		

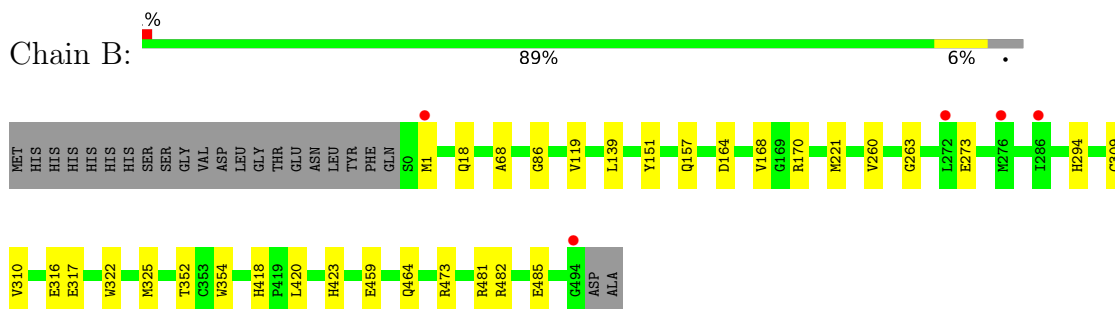
3 Residue-property plots [i](#)

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.

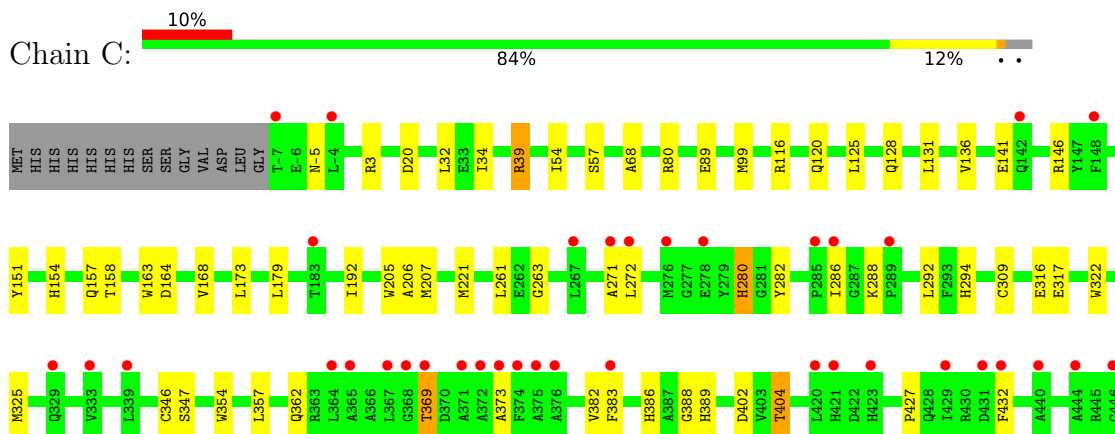
- Molecule 1: Aromatic Acid Decarboxylase

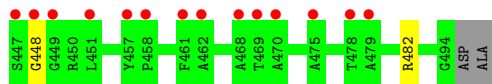


- Molecule 1: Aromatic Acid Decarboxylase



- Molecule 1: Aromatic Acid Decarboxylase





4 Data and refinement statistics i

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	96.85Å 96.85Å 301.48Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.30 – 1.95 50.30 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.6 (50.30-1.95) 99.6 (50.30-1.95)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.77 (at 1.95Å)	Xtrriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.199 , 0.248 0.205 , 0.249	Depositor DCC
R_{free} test set	6027 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	31.7	Xtrriage
Anisotropy	0.470	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 30.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.024 for -h,-k,l	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	12473	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.07% 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, MG, 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	0.60	0/4022	0.79	0/5488
1	B	0.59	0/3955	0.79	0/5399
1	C	0.62	1/4029 (0.0%)	0.82	2/5499 (0.0%)
All	All	0.61	1/12006 (0.0%)	0.80	2/16386 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	383	PHE	CA-CB	-19.56	1.22	1.53

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	383	PHE	N-CA-CB	11.17	126.54	110.12
1	C	89	GLU	N-CA-C	5.58	118.89	111.75

There are no chirality outliers.

There are no planarity outliers.

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	3900	0	3834	35	0
1	B	3844	0	3771	18	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	3913	0	3836	32	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
3	A	8	0	12	0	0
3	B	8	0	12	0	0
3	C	4	0	6	0	0
4	B	6	0	8	0	0
5	A	336	0	0	4	0
5	B	290	0	0	2	0
5	C	161	0	0	1	0
All	All	12473	0	11479	83	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 (83) 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:104[A]:MET:HA	1:A:104[A]:MET:HE2	1.32	1.11
1:A:104[A]:MET:HA	1:A:104[A]:MET:CE	1.91	1.00
1:C:68:ALA:HB3	1:C:317:GLU:HG3	1.62	0.80
1:A:104[A]:MET:HE2	1:A:104[A]:MET:CA	2.12	0.76
1:C:157:GLN:NE2	1:C:164:ASP:OD1	2.21	0.73
1:A:442:ASP:OD2	1:B:418:HIS:HE1	1.77	0.67
1:A:104[A]:MET:HE1	1:A:223:LEU:HD21	1.78	0.65
1:C:32:LEU:HD21	1:C:141:GLU:HA	1.78	0.64
1:C:151:TYR:HB2	1:C:309[B]:CYS:SG	2.37	0.64
1:A:104[A]:MET:HE3	1:A:221:MET:HE2	1.81	0.63
1:C:163:TRP:CZ3	1:C:192:ILE:HD11	2.35	0.62
1:B:418:HIS:HD2	1:B:420:LEU:H	1.48	0.61
1:A:151:TYR:CD1	1:A:309[B]:CYS:SG	2.93	0.61
1:A:13:ASP:OD1	1:A:16:ARG:NH2	2.34	0.60
1:C:163:TRP:CE3	1:C:192:ILE:HD11	2.37	0.59
1:C:382:VAL:O	1:C:388:GLY:HA3	2.03	0.58
1:C:131:LEU:HD13	1:C:179:LEU:HD21	1.88	0.56
1:A:151:TYR:HD1	1:A:309[B]:CYS:HG	1.53	0.55
1:A:22:VAL:HG23	1:A:47:PRO:HB2	1.88	0.55
1:C:125:LEU:H	1:C:128:GLN:HE21	1.54	0.55
1:C:158:THR:HG22	1:C:192:ILE:HD12	1.89	0.54
1:A:68:ALA:HB3	1:A:317:GLU:HG3	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:482:ARG:NH1	1:B:485:GLU:OE1	2.42	0.53
1:A:86:GLY:HA2	5:A:757:HOH:O	2.09	0.53
1:C:263:GLY:HA3	1:C:294:HIS:O	2.09	0.52
1:B:151:TYR:HB2	1:B:309[B]:CYS:SG	2.48	0.52
1:C:20:ASP:OD2	1:C:80:ARG:CG	2.58	0.52
1:A:173:LEU:HA	1:A:179:LEU:HD23	1.92	0.52
1:A:168:VAL:HG21	1:A:322:TRP:CZ2	2.45	0.51
1:C:280:HIS:HD2	1:C:282:TYR:H	1.60	0.50
1:A:183:THR:HG22	1:A:183:THR:O	2.12	0.50
1:A:104[A]:MET:HE1	1:A:223:LEU:CD2	2.42	0.50
1:A:151:TYR:HD1	1:A:309[B]:CYS:SG	2.35	0.49
1:A:104[A]:MET:HE3	1:A:221:MET:CE	2.42	0.49
1:C:154:HIS:CE1	1:C:207:MET:HE3	2.47	0.48
1:C:168:VAL:HG21	1:C:322:TRP:CZ2	2.49	0.48
1:C:34:ILE:HD11	1:C:54:ILE:HG12	1.96	0.48
1:A:263:GLY:HA3	1:A:294:HIS:O	2.14	0.48
1:B:473:ARG:HA	1:B:481:ARG:HD3	1.96	0.47
1:A:157:GLN:NE2	1:A:164:ASP:OD1	2.47	0.46
1:A:220:GLY:HA3	1:A:322:TRP:CZ3	2.50	0.46
1:B:68:ALA:HB3	1:B:317:GLU:HG3	1.98	0.46
1:C:173:LEU:HA	1:C:179:LEU:HD23	1.96	0.46
1:A:419:PRO:O	1:B:423:HIS:HE1	1.98	0.46
1:A:18:GLN:HG3	5:A:687:HOH:O	2.16	0.46
1:A:139:LEU:HD22	1:A:310:VAL:HG22	1.97	0.46
1:C:357:LEU:HD12	1:C:357:LEU:N	2.31	0.46
1:C:402:ASP:OD1	1:C:404:THR:HB	2.15	0.46
1:B:18:GLN:NE2	5:B:743:HOH:O	2.49	0.45
1:C:346:CYS:O	1:C:347:SER:C	2.59	0.45
1:C:39:ARG:HA	1:C:39:ARG:HD2	1.73	0.45
1:C:99[A]:MET:HE3	1:C:99[A]:MET:HB3	1.54	0.45
1:B:170:ARG:NH2	1:B:273:GLU:OE2	2.50	0.45
1:A:99[B]:MET:HE3	1:A:99[B]:MET:HB3	1.69	0.44
1:A:183:THR:HG22	5:A:710:HOH:O	2.18	0.44
1:B:139:LEU:HD22	1:B:310:VAL:HG22	1.99	0.44
1:C:427:PRO:HA	1:C:448:GLY:HA2	1.99	0.44
1:B:221:MET:HA	1:B:325:MET:HE1	1.99	0.44
1:A:16:ARG:HD3	1:A:23:ASP:OD2	2.18	0.44
1:C:271:ALA:HB2	1:C:292:LEU:HD21	2.00	0.44
1:C:369:THR:HG22	1:C:373:ALA:HB3	2.00	0.44
1:A:206:ALA:HA	1:A:261:LEU:O	2.18	0.43
1:B:481:ARG:NH2	5:B:788:HOH:O	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:157:GLN:NE2	1:B:164:ASP:OD1	2.52	0.43
1:B:119:VAL:HG21	1:B:260:VAL:HG23	2.01	0.43
1:C:116:ARG:HA	1:C:120:GLN:OE1	2.19	0.43
1:C:389:HIS:O	1:C:432:PHE:HB2	2.19	0.42
1:B:168:VAL:HG21	1:B:322:TRP:CZ2	2.54	0.42
1:A:321:ILE:HG22	1:A:322:TRP:CE3	2.54	0.42
1:B:263:GLY:HA3	1:B:294:HIS:O	2.19	0.42
1:C:20:ASP:OD2	1:C:80:ARG:HG2	2.20	0.42
1:A:322:TRP:CZ3	1:A:325:MET:HE3	2.55	0.41
1:C:136:VAL:HG12	1:C:146:ARG:HB3	2.02	0.41
1:A:183:THR:HG21	1:A:193:ARG:HD2	2.02	0.41
1:C:206:ALA:HA	1:C:261:LEU:O	2.20	0.41
1:B:459:GLU:HB2	1:B:464:GLN:HB2	2.02	0.41
1:C:3:ARG:HG3	5:C:760:HOH:O	2.21	0.41
1:A:239:GLU:HG2	1:A:240:PRO:HD2	2.03	0.40
1:C:221:MET:HG2	1:C:325:MET:HE1	2.03	0.40
1:A:55:ARG:NH2	5:A:844:HOH:O	2.54	0.40
1:A:21:LEU:HD12	1:A:48:ALA:O	2.22	0.40
1:B:1:MET:HE3	1:B:86:GLY:O	2.21	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	501/518 (97%)	491 (98%)	10 (2%)	0	100	100
1	B	494/518 (95%)	481 (97%)	13 (3%)	0	100	100
1	C	502/518 (97%)	485 (97%)	17 (3%)	0	100	100
All	All	1497/1554 (96%)	1457 (97%)	40 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	401/413 (97%)	390 (97%)	11 (3%)	39	32
1	B	394/413 (95%)	391 (99%)	3 (1%)	73	74
1	C	402/413 (97%)	387 (96%)	15 (4%)	30	20
All	All	1197/1239 (97%)	1168 (98%)	29 (2%)	44	36

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-4	LEU
1	A	104[A]	MET
1	A	104[B]	MET
1	A	142[A]	GLN
1	A	142[B]	GLN
1	A	179	LEU
1	A	272	LEU
1	A	303	GLN
1	A	316	GLU
1	A	352	THR
1	A	354	TRP
1	B	316	GLU
1	B	352	THR
1	B	354	TRP
1	C	-5	ASN
1	C	39	ARG
1	C	57	SER
1	C	205	TRP
1	C	272	LEU
1	C	280	HIS
1	C	286	ILE
1	C	288	LYS
1	C	316	GLU
1	C	354	TRP
1	C	362	GLN
1	C	369	THR

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Mol	Chain	Res	Type
1	C	386	HIS
1	C	404	THR
1	C	482	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	90	HIS
1	A	303	GLN
1	B	18	GLN
1	B	157	GLN
1	B	389	HIS
1	B	418	HIS
1	B	423	HIS
1	C	25	HIS
1	C	128	GLN
1	C	336	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](#)

Of 9 ligands modelled in this entry, 3 are monoatomic - leaving 6 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	502	-	3,3,3	0.47	0	2,2,2	0.28	0
3	EDO	C	502	-	3,3,3	0.44	0	2,2,2	0.35	0
3	EDO	B	504	-	3,3,3	0.43	0	2,2,2	0.32	0
4	GOL	B	502	-	5,5,5	0.17	0	5,5,5	0.47	0
3	EDO	A	503	-	3,3,3	0.45	0	2,2,2	0.28	0
3	EDO	B	503	-	3,3,3	0.44	0	2,2,2	0.35	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	502	-	-	1/1/1/1	-
3	EDO	C	502	-	-	0/1/1/1	-
3	EDO	B	504	-	-	1/1/1/1	-
4	GOL	B	502	-	-	0/4/4/4	-
3	EDO	A	503	-	-	0/1/1/1	-
3	EDO	B	503	-	-	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
3	A	502	EDO	O1-C1-C2-O2
3	B	504	EDO	O1-C1-C2-O2
3	B	503	EDO	O1-C1-C2-O2

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues

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	499/518 (96%)	0.09	5 (1%) 79 84	20, 33, 47, 79	4 (0%)
1	B	495/518 (95%)	0.07	5 (1%) 79 84	22, 34, 49, 64	1 (0%)
1	C	502/518 (96%)	0.74	51 (10%) 12 13	22, 43, 69, 85	3 (0%)
All	All	1496/1554 (96%)	0.30	61 (4%) 41 48	20, 36, 60, 85	8 (0%)

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	-4	LEU	6.6
1	C	286	ILE	4.5
1	B	286	ILE	4.1
1	C	371	ALA	3.6
1	C	429	ILE	3.6
1	C	272	LEU	3.5
1	C	458	PRO	3.4
1	C	271	ALA	3.3
1	C	468	ALA	3.3
1	C	-7	THR	3.3
1	C	448	GLY	3.3
1	C	440	ALA	3.2
1	C	420	LEU	3.2
1	C	375	ALA	3.1
1	C	449	GLY	3.0
1	C	373	ALA	2.9
1	C	446	GLY	2.9
1	C	374	PHE	2.9
1	C	383	PHE	2.9
1	A	-3	TYR	2.9
1	C	367	LEU	2.8
1	A	183	THR	2.8
1	C	333	VAL	2.8

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Mol	Chain	Res	Type	RSRZ
1	C	-4	LEU	2.8
1	C	461	PHE	2.7
1	C	369	THR	2.7
1	A	-2	PHE	2.7
1	C	376	ALA	2.6
1	C	462	ALA	2.6
1	C	447	SER	2.6
1	C	148	PHE	2.5
1	C	479	ALA	2.5
1	C	432	PHE	2.4
1	C	372	ALA	2.4
1	C	285	PRO	2.4
1	C	478	THR	2.4
1	C	142	GLN	2.4
1	C	289	PRO	2.4
1	C	451	LEU	2.3
1	C	470	ALA	2.3
1	C	278	GLU	2.3
1	C	364	LEU	2.3
1	C	423	HIS	2.3
1	C	475	ALA	2.3
1	B	494	GLY	2.3
1	C	267	LEU	2.3
1	A	142[A]	GLN	2.2
1	B	1	MET	2.2
1	C	421	HIS	2.2
1	C	444	ALA	2.1
1	C	431	ASP	2.1
1	B	272	LEU	2.1
1	C	183	THR	2.1
1	C	368	GLY	2.1
1	C	276	MET	2.1
1	C	339	LEU	2.1
1	C	365	ALA	2.1
1	B	276	MET	2.0
1	C	329	GLN	2.0
1	C	469	THR	2.0
1	C	457	TYR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	GOL	B	502	6/6	0.60	0.18	58,61,61,66	0
3	EDO	B	504	4/4	0.83	0.14	48,49,50,51	0
3	EDO	A	503	4/4	0.84	0.16	49,49,49,51	0
3	EDO	A	502	4/4	0.86	0.16	53,56,57,58	0
3	EDO	B	503	4/4	0.86	0.10	59,59,59,61	0
3	EDO	C	502	4/4	0.89	0.10	57,59,60,60	0
2	MG	B	501	1/1	0.96	0.04	31,31,31,31	0
2	MG	C	501	1/1	0.98	0.05	38,38,38,38	0
2	MG	A	501	1/1	0.99	0.03	30,30,30,30	0

6.5 Other polymers [i](#)

There are no such residues in this entry.