



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 25, 2026 – 04:39 AM UTC

PDB ID : 9JVL / pdb_00009jvl
Title : Co-crystal structure of PHBDD-M1-L163G-T234A with N-cyclopropyl-4-methylbenzamide
Authors : Yang, J.; Gao, L.; Qiu, X.
Deposited on : 2024-10-09
Resolution : 2.10 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

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

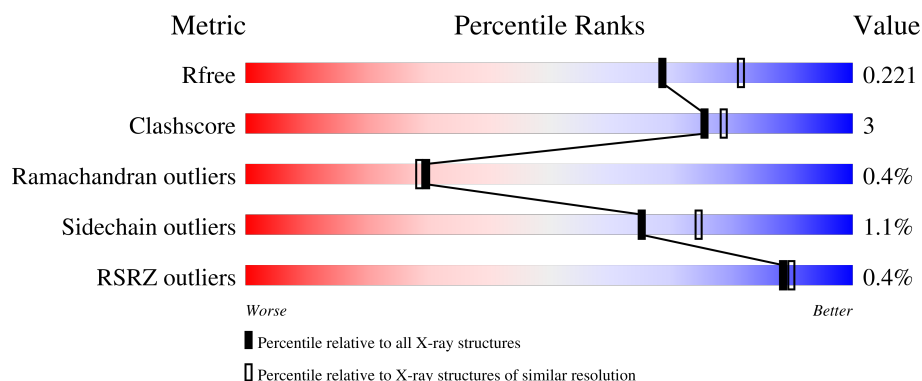
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	493	92% 6% .
1	B	493	93% 6% .
1	C	493	% 92% 6% .
1	D	493	90% 9% .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 15838 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 4-hydroxybenzaldehyde dehydrogenase (NADP(+)).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	486	Total	C	N	O	S	0	0	0
			3691	2325	657	699	10			
1	B	488	Total	C	N	O	S	0	0	0
			3710	2337	662	701	10			
1	C	486	Total	C	N	O	S	0	0	0
			3687	2323	657	697	10			
1	D	487	Total	C	N	O	S	0	0	0
			3699	2331	658	700	10			

There are 24 discrepancies between the modelled and reference sequences:

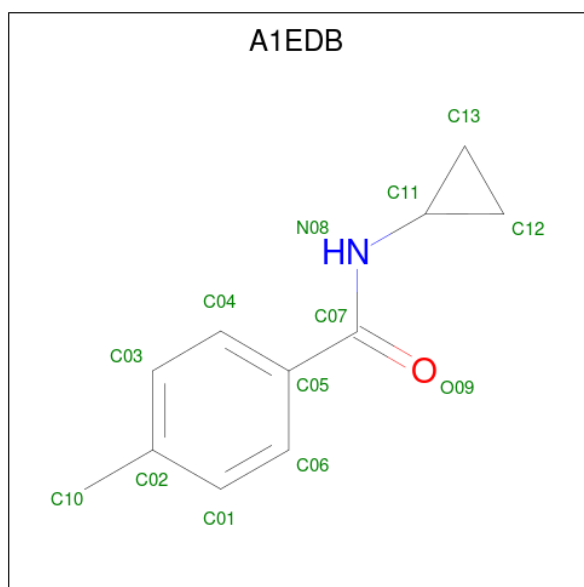
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP Q59702
A	0	SER	-	expression tag	UNP Q59702
A	163	GLY	LEU	engineered mutation	UNP Q59702
A	166	MET	ARG	conflict	UNP Q59702
A	234	ALA	THR	engineered mutation	UNP Q59702
A	259	LEU	GLU	conflict	UNP Q59702
B	-1	GLY	-	expression tag	UNP Q59702
B	0	SER	-	expression tag	UNP Q59702
B	163	GLY	LEU	engineered mutation	UNP Q59702
B	166	MET	ARG	conflict	UNP Q59702
B	234	ALA	THR	engineered mutation	UNP Q59702
B	259	LEU	GLU	conflict	UNP Q59702
C	-1	GLY	-	expression tag	UNP Q59702
C	0	SER	-	expression tag	UNP Q59702
C	163	GLY	LEU	engineered mutation	UNP Q59702
C	166	MET	ARG	conflict	UNP Q59702
C	234	ALA	THR	engineered mutation	UNP Q59702
C	259	LEU	GLU	conflict	UNP Q59702
D	-1	GLY	-	expression tag	UNP Q59702
D	0	SER	-	expression tag	UNP Q59702
D	163	GLY	LEU	engineered mutation	UNP Q59702

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Chain	Residue	Modelled	Actual	Comment	Reference
D	166	MET	ARG	conflict	UNP Q59702
D	234	ALA	THR	engineered mutation	UNP Q59702
D	259	LEU	GLU	conflict	UNP Q59702

- Molecule 2 is {N}-cyclopropyl-4-methyl-benzamide (CCD ID: A1EDB) (formula: C₁₁H₁₃NO).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			13	11	1	1		
2	B	1	Total	C	N	O	0	0
			13	11	1	1		
2	C	1	Total	C	N	O	0	0
			13	11	1	1		
2	D	1	Total	C	N	O	0	0
			13	11	1	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	304	Total	O	0	0
			304	304		
3	B	214	Total	O	0	0
			214	214		
3	C	218	Total	O	0	0
			218	218		

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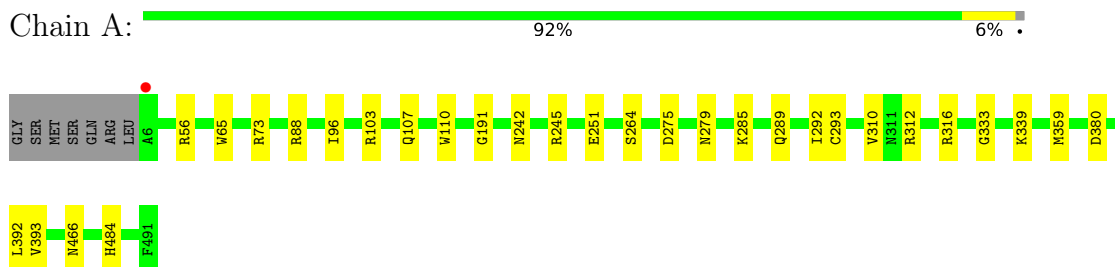
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	263	Total 263	O 263	0	0

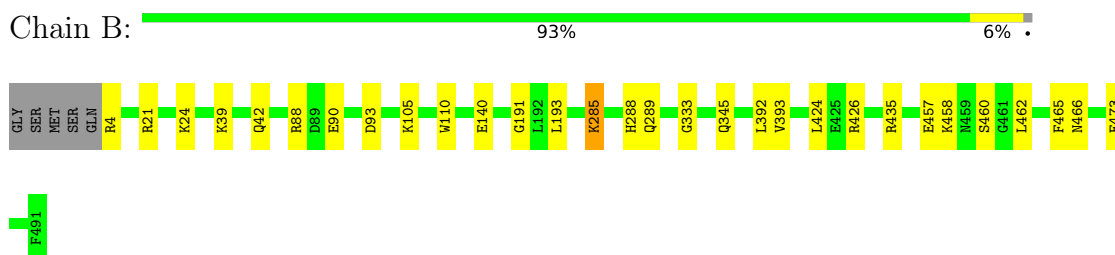
3 Residue-property plots [i](#)

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.

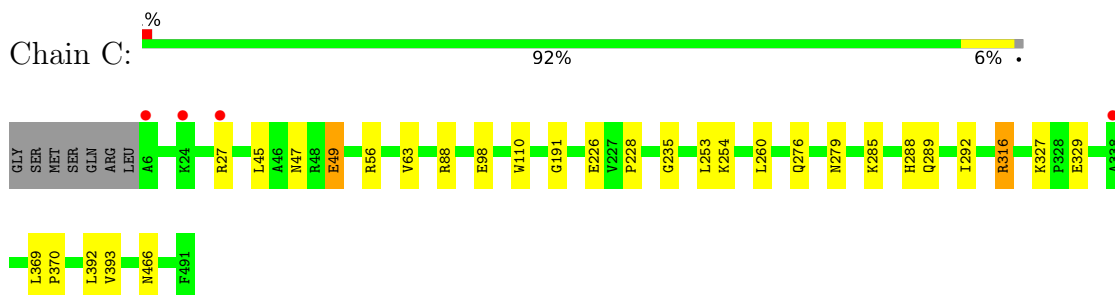
- Molecule 1: 4-hydroxybenzaldehyde dehydrogenase (NADP(+))



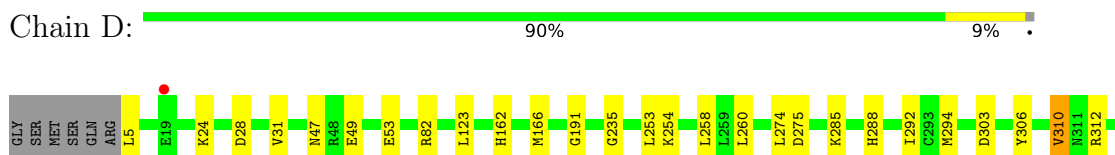
- Molecule 1: 4-hydroxybenzaldehyde dehydrogenase (NADP(+))



- Molecule 1: 4-hydroxybenzaldehyde dehydrogenase (NADP(+))



- Molecule 1: 4-hydroxybenzaldehyde dehydrogenase (NADP(+))





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	90.25Å 124.89Å 90.68Å 90.00° 105.69° 90.00°	Depositor
Resolution (Å)	43.92 – 2.10 43.92 – 2.10	Depositor EDS
% Data completeness (in resolution range)	74.0 (43.92-2.10) 93.5 (43.92-2.10)	Depositor EDS
R_{merge}	0.20	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.90 (at 2.10Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.180 , 0.221 0.180 , 0.221	Depositor DCC
R_{free} test set	1998 reflections (1.78%)	wwPDB-VP
Wilson B-factor (Å ²)	25.1	Xtriage
Anisotropy	0.096	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 35.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.020 for l,-k,h	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	15838	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.41% 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: A1EDB

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	0.45	0/3768	0.55	0/5117
1	B	0.42	0/3787	0.52	0/5142
1	C	0.40	0/3764	0.51	0/5112
1	D	0.43	0/3776	0.55	0/5128
All	All	0.42	0/15095	0.53	0/20499

There are no bond length outliers.

There are no bond angle outliers.

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	3691	0	3644	19	0
1	B	3710	0	3668	26	0
1	C	3687	0	3640	17	0
1	D	3699	0	3655	27	0
2	A	13	0	0	0	0
2	B	13	0	0	0	0
2	C	13	0	0	0	0
2	D	13	0	0	0	0
3	A	304	0	0	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	214	0	0	9	0
3	C	218	0	0	6	0
3	D	263	0	0	9	0
All	All	15838	0	14607	87	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 (87) 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:B:458:LYS:NZ	3:B:602:HOH:O	2.09	0.86
1:A:484:HIS:O	3:A:601:HOH:O	1.92	0.85
1:D:82:ARG:NH2	3:D:602:HOH:O	2.12	0.82
1:D:123:LEU:HD22	1:D:471:ILE:HD11	1.64	0.79
1:B:473:GLU:OE1	3:B:601:HOH:O	2.01	0.76
1:C:98:GLU:O	3:C:601:HOH:O	2.05	0.74
1:B:345:GLN:NE2	3:B:604:HOH:O	2.20	0.72
1:D:435:ARG:NH2	3:D:603:HOH:O	2.13	0.72
1:A:380:ASP:OD2	3:A:602:HOH:O	2.07	0.72
1:D:53:GLU:OE1	3:D:601:HOH:O	2.10	0.69
1:A:285:LYS:HD3	1:A:393:VAL:HB	1.75	0.68
1:B:435:ARG:HD3	1:B:458:LYS:HD2	1.76	0.68
1:B:42:GLN:NE2	3:B:607:HOH:O	2.25	0.68
1:D:275:ASP:OD1	1:D:312:ARG:NH1	2.27	0.67
1:B:285:LYS:HD3	1:B:393:VAL:HB	1.81	0.62
1:C:226:GLU:OE1	3:C:602:HOH:O	2.16	0.62
1:B:4:ARG:NH2	1:B:90:GLU:H	1.99	0.59
1:B:93:ASP:HB3	3:B:623:HOH:O	2.01	0.59
1:D:28:ASP:OD2	3:D:604:HOH:O	2.17	0.59
1:A:251:GLU:OE2	1:A:251:GLU:N	2.31	0.58
1:D:285:LYS:HD3	1:D:393:VAL:HB	1.86	0.57
1:C:279:ASN:OD1	1:C:316:ARG:NH1	2.39	0.55
1:B:289:GLN:NE2	3:B:605:HOH:O	2.24	0.55
1:C:27:ARG:HB3	1:C:45:LEU:HD12	1.89	0.54
1:C:327:LYS:NZ	1:C:329:GLU:OE2	2.33	0.54
1:B:88:ARG:HG3	1:B:110:TRP:CE2	2.43	0.54
1:D:288:HIS:HA	3:D:606:HOH:O	2.07	0.54
1:A:242:ASN:OD1	1:A:245:ARG:NH1	2.41	0.53
1:B:21:ARG:NH1	1:B:24:LYS:HE2	2.23	0.52
1:B:39:LYS:HD3	3:B:607:HOH:O	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:306:TYR:O	1:D:310:VAL:HG13	2.10	0.52
1:D:162:HIS:NE2	1:D:166:MET:HE3	2.25	0.51
1:D:316:ARG:NH2	3:D:609:HOH:O	2.39	0.51
1:B:462:LEU:HG	1:C:253:LEU:HD21	1.93	0.50
1:D:425:GLU:HG3	1:D:426:ARG:N	2.26	0.50
1:C:289:GLN:NE2	3:C:611:HOH:O	2.42	0.49
1:D:457:GLU:OE2	3:D:605:HOH:O	2.19	0.49
1:B:289:GLN:NE2	1:B:333:GLY:H	2.11	0.48
1:A:279:ASN:OD1	1:A:316:ARG:NH1	2.33	0.48
1:C:88:ARG:HG3	1:C:110:TRP:CZ2	2.48	0.48
1:C:285:LYS:HD3	1:C:393:VAL:HB	1.95	0.48
1:D:435:ARG:N	1:D:435:ARG:HD3	2.29	0.48
1:A:96:ILE:HD11	1:A:103:ARG:HG2	1.96	0.47
1:B:457:GLU:O	1:B:460:SER:HB2	2.14	0.47
1:B:426:ARG:NH1	1:B:426:ARG:HG3	2.29	0.47
1:B:288:HIS:HA	3:B:605:HOH:O	2.13	0.47
1:B:140:GLU:OE1	3:B:603:HOH:O	2.20	0.47
1:D:47:ASN:OD1	1:D:49:GLU:HG3	2.15	0.47
1:A:107:GLN:NE2	3:A:611:HOH:O	2.46	0.46
1:D:399:ARG:HG2	1:D:403:HIS:ND1	2.30	0.46
1:A:88:ARG:HG3	1:A:110:TRP:CE2	2.50	0.46
1:D:24:LYS:HE3	1:D:24:LYS:HA	1.97	0.46
1:A:88:ARG:HG3	1:A:110:TRP:CZ2	2.51	0.46
1:A:65:TRP:CH2	1:A:73:ARG:HG2	2.51	0.46
1:B:21:ARG:HH11	1:B:24:LYS:HE2	1.81	0.45
1:C:56:ARG:HD3	3:C:610:HOH:O	2.17	0.45
1:A:310:VAL:HG12	1:A:359:MET:HE2	1.97	0.45
1:A:310:VAL:CG1	1:A:359:MET:HE2	2.46	0.45
1:B:285:LYS:O	1:B:285:LYS:HE3	2.17	0.45
1:A:289:GLN:HG2	1:A:333:GLY:O	2.17	0.44
1:B:424:LEU:HD22	1:D:425:GLU:HB2	1.99	0.44
1:C:63:VAL:HB	3:C:712:HOH:O	2.18	0.44
1:C:316:ARG:HE	1:C:316:ARG:HB3	1.64	0.44
1:A:275:ASP:OD1	1:A:312:ARG:NH1	2.47	0.44
1:D:453:PRO:HB3	1:D:469:TRP:CE2	2.53	0.43
1:D:274:LEU:HA	1:D:274:LEU:HD23	1.80	0.43
1:A:293:CYS:SG	3:A:681:HOH:O	2.49	0.43
1:D:435:ARG:N	1:D:435:ARG:CD	2.82	0.43
1:C:235:GLY:O	1:C:260:LEU:HA	2.19	0.42
1:B:105:LYS:HE3	1:B:288:HIS:CD2	2.54	0.42
1:D:235:GLY:O	1:D:260:LEU:HA	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:258:LEU:HD12	1:D:462:LEU:HD12	2.01	0.42
1:A:56:ARG:NH2	3:A:617:HOH:O	2.52	0.42
1:C:288:HIS:HA	3:C:611:HOH:O	2.20	0.42
1:A:65:TRP:CZ2	1:A:73:ARG:HG2	2.55	0.41
1:A:339:LYS:HE2	3:A:693:HOH:O	2.20	0.41
1:B:39:LYS:O	1:B:39:LYS:HG3	2.20	0.41
1:C:47:ASN:OD1	1:C:49:GLU:HB2	2.21	0.41
1:D:327:LYS:HD2	3:D:843:HOH:O	2.21	0.41
1:D:429:LYS:NZ	3:D:626:HOH:O	2.52	0.41
1:B:465:PHE:O	1:B:466:ASN:HB2	2.21	0.41
1:B:289:GLN:HG2	1:B:333:GLY:O	2.21	0.40
1:C:228:PRO:HG2	1:C:254:LYS:HE2	2.02	0.40
1:D:253:LEU:HD13	1:D:254:LYS:N	2.36	0.40
1:D:292:ILE:HG22	1:D:294:MET:H	1.87	0.40
1:C:369:LEU:HA	1:C:370:PRO:HD3	1.96	0.40
1:B:435:ARG:HH11	1:B:458:LYS:HG3	1.86	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	484/493 (98%)	471 (97%)	11 (2%)	2 (0%)	30	28
1	B	486/493 (99%)	476 (98%)	9 (2%)	1 (0%)	43	44
1	C	484/493 (98%)	469 (97%)	13 (3%)	2 (0%)	30	28
1	D	485/493 (98%)	472 (97%)	11 (2%)	2 (0%)	30	28
All	All	1939/1972 (98%)	1888 (97%)	44 (2%)	7 (0%)	30	28

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	466	ASN
1	C	466	ASN
1	D	466	ASN
1	A	191	GLY
1	C	191	GLY
1	D	191	GLY
1	B	191	GLY

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	381/388 (98%)	378 (99%)	3 (1%)	73	81
1	B	383/388 (99%)	380 (99%)	3 (1%)	73	81
1	C	380/388 (98%)	375 (99%)	5 (1%)	61	69
1	D	382/388 (98%)	376 (98%)	6 (2%)	55	64
All	All	1526/1552 (98%)	1509 (99%)	17 (1%)	65	74

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	264	SER
1	A	292	ILE
1	A	392	LEU
1	B	193	LEU
1	B	285	LYS
1	B	392	LEU
1	C	49	GLU
1	C	276	GLN
1	C	292	ILE
1	C	316	ARG
1	C	392	LEU
1	D	5	LEU
1	D	31	VAL
1	D	303	ASP
1	D	310	VAL

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Mol	Chain	Res	Type
1	D	392	LEU
1	D	476	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	107	GLN
1	A	162	HIS
1	A	291	GLN
1	A	365	GLN
1	B	42	GLN
1	B	60	GLN
1	B	75	GLN
1	B	252	HIS
1	B	291	GLN
1	C	162	HIS
1	C	238	GLN
1	C	345	GLN
1	C	484	HIS
1	D	42	GLN
1	D	311	ASN
1	D	365	GLN
1	D	484	HIS

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](#)

4 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	A1EDB	C	501	-	14,14,14	5.74	8 (57%)	19,19,19	1.68	3 (15%)
2	A1EDB	B	501	-	14,14,14	5.74	8 (57%)	19,19,19	1.58	4 (21%)
2	A1EDB	D	501	-	14,14,14	5.71	8 (57%)	19,19,19	1.60	5 (26%)
2	A1EDB	A	501	-	14,14,14	5.71	7 (50%)	19,19,19	1.66	3 (15%)

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	A1EDB	C	501	-	-	0/8/10/10	0/2/2/2
2	A1EDB	B	501	-	-	0/8/10/10	0/2/2/2
2	A1EDB	D	501	-	-	0/8/10/10	0/2/2/2
2	A1EDB	A	501	-	-	0/8/10/10	0/2/2/2

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	A1EDB	C04-C05	9.83	1.54	1.39
2	A	501	A1EDB	C04-C03	9.45	1.54	1.38
2	C	501	A1EDB	C06-C05	9.42	1.53	1.39
2	B	501	A1EDB	C04-C03	9.34	1.53	1.38
2	D	501	A1EDB	C04-C03	9.29	1.53	1.38
2	D	501	A1EDB	C06-C05	9.28	1.53	1.39
2	B	501	A1EDB	C06-C05	9.26	1.53	1.39
2	C	501	A1EDB	C04-C05	9.18	1.53	1.39
2	C	501	A1EDB	C04-C03	9.15	1.53	1.38
2	D	501	A1EDB	C06-C01	9.13	1.53	1.38
2	A	501	A1EDB	C06-C01	9.09	1.53	1.38
2	D	501	A1EDB	C04-C05	9.09	1.53	1.39
2	A	501	A1EDB	C06-C05	9.06	1.53	1.39
2	C	501	A1EDB	C06-C01	9.04	1.53	1.38
2	A	501	A1EDB	C04-C05	9.01	1.53	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	A1EDB	C06-C01	8.40	1.52	1.38
2	B	501	A1EDB	C03-C02	6.52	1.54	1.38
2	A	501	A1EDB	C01-C02	6.43	1.54	1.38
2	C	501	A1EDB	C01-C02	6.38	1.54	1.38
2	C	501	A1EDB	C03-C02	6.34	1.53	1.38
2	A	501	A1EDB	C03-C02	6.18	1.53	1.38
2	D	501	A1EDB	C01-C02	6.00	1.53	1.38
2	D	501	A1EDB	C03-C02	5.94	1.52	1.38
2	B	501	A1EDB	C01-C02	5.81	1.52	1.38
2	D	501	A1EDB	C07-N08	5.46	1.46	1.34
2	B	501	A1EDB	C07-N08	5.40	1.46	1.34
2	A	501	A1EDB	C07-N08	5.20	1.46	1.34
2	C	501	A1EDB	C07-N08	5.15	1.46	1.34
2	D	501	A1EDB	C13-C11	2.41	1.54	1.48
2	C	501	A1EDB	C13-C11	2.24	1.54	1.48
2	B	501	A1EDB	C13-C11	2.16	1.53	1.48

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	A1EDB	C12-C11-N08	-5.05	111.89	118.52
2	C	501	A1EDB	C12-C11-N08	-4.45	112.69	118.52
2	B	501	A1EDB	C13-C11-N08	-3.67	113.71	118.52
2	C	501	A1EDB	C13-C11-N08	-3.56	113.84	118.52
2	D	501	A1EDB	C11-N08-C07	-3.41	116.09	122.56
2	B	501	A1EDB	C11-N08-C07	-3.39	116.14	122.56
2	A	501	A1EDB	C13-C11-N08	-2.64	115.05	118.52
2	A	501	A1EDB	C11-N08-C07	-2.44	117.95	122.56
2	D	501	A1EDB	C06-C05-C04	2.35	121.55	118.57
2	D	501	A1EDB	C05-C07-N08	2.31	121.33	117.04
2	C	501	A1EDB	C06-C01-C02	-2.30	118.72	121.37
2	B	501	A1EDB	C03-C04-C05	-2.29	118.35	120.80
2	D	501	A1EDB	C12-C11-N08	-2.14	115.71	118.52
2	B	501	A1EDB	C06-C05-C04	2.06	121.18	118.57
2	D	501	A1EDB	C13-C11-N08	-2.03	115.86	118.52

There are no chirality outliers.

There are no torsion outliers.

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 [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	486/493 (98%)	-0.30	1 (0%) 91 92	16, 22, 34, 47	0
1	B	488/493 (98%)	-0.16	0 100 100	17, 26, 40, 52	0
1	C	486/493 (98%)	-0.05	4 (0%) 82 84	18, 28, 44, 65	0
1	D	487/493 (98%)	-0.22	2 (0%) 88 90	17, 24, 38, 48	0
All	All	1947/1972 (98%)	-0.18	7 (0%) 88 90	16, 25, 41, 65	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	24	LYS	3.7
1	D	424	LEU	2.8
1	C	27	ARG	2.6
1	A	6	ALA	2.5
1	C	6	ALA	2.5
1	C	338	ALA	2.4
1	D	19	GLU	2.1

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

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	A1EDB	C	501	13/13	0.93	0.09	27,30,32,36	0
2	A1EDB	D	501	13/13	0.94	0.07	19,24,26,27	0
2	A1EDB	B	501	13/13	0.95	0.07	22,25,28,30	0
2	A1EDB	A	501	13/13	0.96	0.06	20,23,30,34	0

6.5 Other polymers

There are no such residues in this entry.