



Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 07:34 pm BST

PDB ID : 2NV1
Title : Structure of the synthase subunit Pdx1 (YaaD) of PLP synthase from *Bacillus subtilis*
Authors : Strohmeier, M.; Tews, I.; Sinning, I.
Deposited on : 2006-11-10
Resolution : 2.08 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

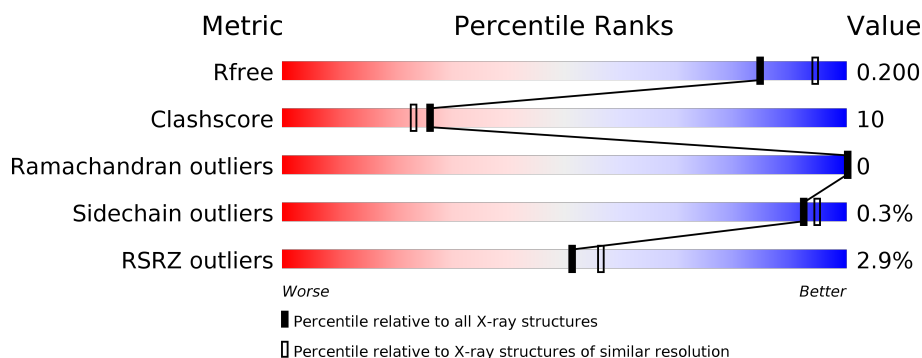
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.08 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}	130704	6189 (2.10-2.06)
Clashscore	141614	6738 (2.10-2.06)
Ramachandran outliers	138981	6663 (2.10-2.06)
Sidechain outliers	138945	6664 (2.10-2.06)
RSRZ outliers	127900	6057 (2.10-2.06)

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 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	305	<div> <div>4%</div> <div> <div></div> <div>74%</div> <div>11%</div> <div>15%</div> </div> </div>
1	B	305	<div> <div>2%</div> <div> <div></div> <div>72%</div> <div>10%</div> <div>17%</div> </div> </div>
1	C	305	<div> <div>3%</div> <div> <div></div> <div>69%</div> <div>14%</div> <div>16%</div> </div> </div>
1	D	305	<div> <div>3%</div> <div> <div></div> <div>70%</div> <div>13%</div> <div>16%</div> </div> </div>
1	E	305	<div> <div>2%</div> <div> <div></div> <div>73%</div> <div>11%</div> <div>16%</div> </div> </div>
1	F	305	<div> <div>2%</div> <div> <div></div> <div>71%</div> <div>10%</div> <div>18%</div> </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	CL	A	3001	-	-	X	-
4	EDO	B	3011	-	-	-	X
4	EDO	C	3017	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 13822 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 Pyridoxal biosynthesis lyase pdxS.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	259	Total	C	N	O	S	0	2	0
			1930	1209	330	375	16			
1	B	252	Total	C	N	O	S	0	2	0
			1875	1177	322	362	14			
1	C	256	Total	C	N	O	S	0	2	0
			1915	1199	329	371	16			
1	D	256	Total	C	N	O	S	0	1	0
			1907	1196	327	368	16			
1	E	256	Total	C	N	O	S	0	2	0
			1910	1196	326	371	17			
1	F	250	Total	C	N	O	S	0	1	0
			1850	1161	317	358	14			

There are 66 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	MET	-	CLONING ARTIFACT	UNP P37527
A	-1	ALA	-	CLONING ARTIFACT	UNP P37527
A	0	SER	-	CLONING ARTIFACT	UNP P37527
A	295	LEU	-	EXPRESSION TAG	UNP P37527
A	296	GLU	-	EXPRESSION TAG	UNP P37527
A	297	HIS	-	EXPRESSION TAG	UNP P37527
A	298	HIS	-	EXPRESSION TAG	UNP P37527
A	299	HIS	-	EXPRESSION TAG	UNP P37527
A	300	HIS	-	EXPRESSION TAG	UNP P37527
A	301	HIS	-	EXPRESSION TAG	UNP P37527
A	302	HIS	-	EXPRESSION TAG	UNP P37527
B	-2	MET	-	CLONING ARTIFACT	UNP P37527
B	-1	ALA	-	CLONING ARTIFACT	UNP P37527
B	0	SER	-	CLONING ARTIFACT	UNP P37527
B	295	LEU	-	EXPRESSION TAG	UNP P37527
B	296	GLU	-	EXPRESSION TAG	UNP P37527
B	297	HIS	-	EXPRESSION TAG	UNP P37527

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Chain	Residue	Modelled	Actual	Comment	Reference
B	298	HIS	-	EXPRESSION TAG	UNP P37527
B	299	HIS	-	EXPRESSION TAG	UNP P37527
B	300	HIS	-	EXPRESSION TAG	UNP P37527
B	301	HIS	-	EXPRESSION TAG	UNP P37527
B	302	HIS	-	EXPRESSION TAG	UNP P37527
C	-2	MET	-	CLONING ARTIFACT	UNP P37527
C	-1	ALA	-	CLONING ARTIFACT	UNP P37527
C	0	SER	-	CLONING ARTIFACT	UNP P37527
C	295	LEU	-	EXPRESSION TAG	UNP P37527
C	296	GLU	-	EXPRESSION TAG	UNP P37527
C	297	HIS	-	EXPRESSION TAG	UNP P37527
C	298	HIS	-	EXPRESSION TAG	UNP P37527
C	299	HIS	-	EXPRESSION TAG	UNP P37527
C	300	HIS	-	EXPRESSION TAG	UNP P37527
C	301	HIS	-	EXPRESSION TAG	UNP P37527
C	302	HIS	-	EXPRESSION TAG	UNP P37527
D	-2	MET	-	CLONING ARTIFACT	UNP P37527
D	-1	ALA	-	CLONING ARTIFACT	UNP P37527
D	0	SER	-	CLONING ARTIFACT	UNP P37527
D	295	LEU	-	EXPRESSION TAG	UNP P37527
D	296	GLU	-	EXPRESSION TAG	UNP P37527
D	297	HIS	-	EXPRESSION TAG	UNP P37527
D	298	HIS	-	EXPRESSION TAG	UNP P37527
D	299	HIS	-	EXPRESSION TAG	UNP P37527
D	300	HIS	-	EXPRESSION TAG	UNP P37527
D	301	HIS	-	EXPRESSION TAG	UNP P37527
D	302	HIS	-	EXPRESSION TAG	UNP P37527
E	-2	MET	-	CLONING ARTIFACT	UNP P37527
E	-1	ALA	-	CLONING ARTIFACT	UNP P37527
E	0	SER	-	CLONING ARTIFACT	UNP P37527
E	295	LEU	-	EXPRESSION TAG	UNP P37527
E	296	GLU	-	EXPRESSION TAG	UNP P37527
E	297	HIS	-	EXPRESSION TAG	UNP P37527
E	298	HIS	-	EXPRESSION TAG	UNP P37527
E	299	HIS	-	EXPRESSION TAG	UNP P37527
E	300	HIS	-	EXPRESSION TAG	UNP P37527
E	301	HIS	-	EXPRESSION TAG	UNP P37527
E	302	HIS	-	EXPRESSION TAG	UNP P37527
F	-2	MET	-	CLONING ARTIFACT	UNP P37527
F	-1	ALA	-	CLONING ARTIFACT	UNP P37527
F	0	SER	-	CLONING ARTIFACT	UNP P37527
F	295	LEU	-	EXPRESSION TAG	UNP P37527

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Chain	Residue	Modelled	Actual	Comment	Reference
F	296	GLU	-	EXPRESSION TAG	UNP P37527
F	297	HIS	-	EXPRESSION TAG	UNP P37527
F	298	HIS	-	EXPRESSION TAG	UNP P37527
F	299	HIS	-	EXPRESSION TAG	UNP P37527
F	300	HIS	-	EXPRESSION TAG	UNP P37527
F	301	HIS	-	EXPRESSION TAG	UNP P37527
F	302	HIS	-	EXPRESSION TAG	UNP P37527

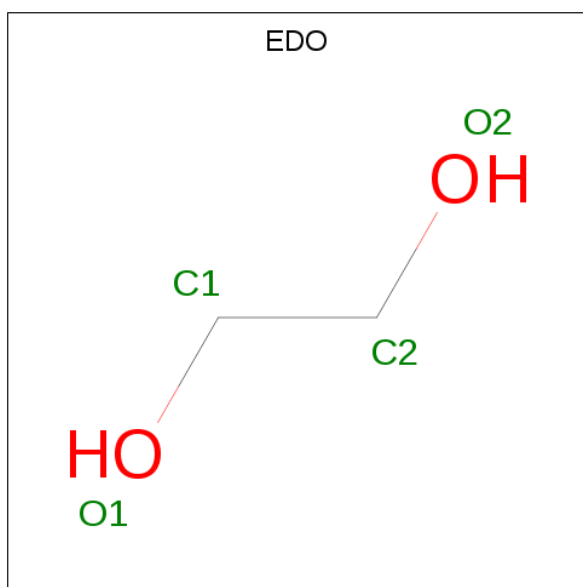
- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	D	2	Total Cl 2 2	0	0
2	E	2	Total Cl 2 2	0	0
2	B	2	Total Cl 2 2	0	0
2	C	2	Total Cl 2 2	0	0
2	A	2	Total Cl 2 2	0	0
2	F	2	Total Cl 2 2	0	0

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	D	2	Total Mg 2 2	0	0
3	E	2	Total Mg 2 2	0	0
3	B	1	Total Mg 1 1	0	0
3	C	3	Total Mg 3 3	0	0
3	A	3	Total Mg 3 3	0	0
3	F	1	Total Mg 1 1	0	0

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	381	Total	O	0	0
			381	381		

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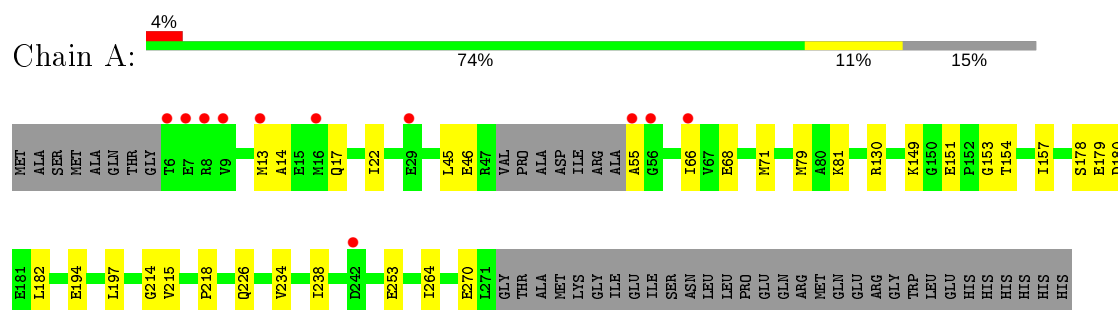
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	425	Total 425	O 425	0	0
5	C	373	Total 373	O 373	0	0
5	D	388	Total 388	O 388	0	0
5	E	398	Total 398	O 398	0	0
5	F	402	Total 402	O 402	0	0

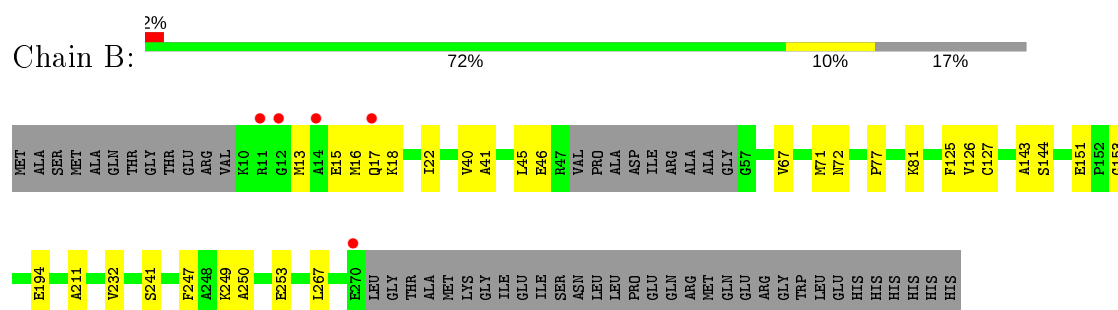
3 Residue-property plots [i](#)

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

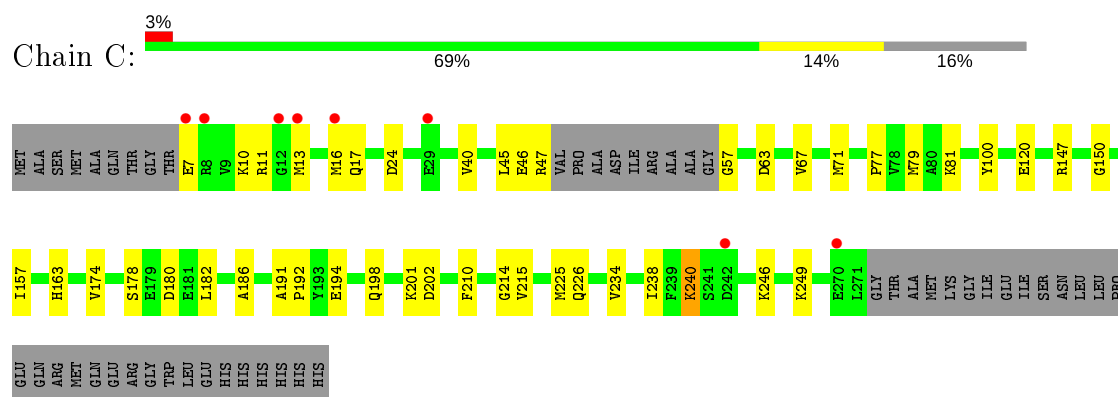
• Molecule 1: Pyridoxal biosynthesis lyase pdxS



• Molecule 1: Pyridoxal biosynthesis lyase pdxS



• Molecule 1: Pyridoxal biosynthesis lyase pdxS



• Molecule 1: Pyridoxal biosynthesis lyase pdxS



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	101.06Å 106.20Å 182.33Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.68 – 2.08 46.76 – 2.08	Depositor EDS
% Data completeness (in resolution range)	98.6 (48.68-2.08) 98.6 (46.76-2.08)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.81 (at 2.08Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.142 , 0.191 0.151 , 0.200	Depositor DCC
R_{free} test set	5782 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	25.1	Xtriage
Anisotropy	0.043	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 69.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.017 for k,h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	13822	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.31% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, EDO, CL

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	0.76	0/1960	0.76	1/2643 (0.0%)
1	B	0.76	0/1905	0.76	0/2570
1	C	0.75	0/1945	0.73	1/2620 (0.0%)
1	D	0.75	0/1932	0.74	1/2604 (0.0%)
1	E	0.80	0/1940	0.74	1/2616 (0.0%)
1	F	0.82	0/1876	0.77	1/2532 (0.0%)
All	All	0.77	0/11558	0.75	5/15585 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	172	ARG	NE-CZ-NH1	5.79	123.20	120.30
1	E	172	ARG	NE-CZ-NH1	5.29	122.94	120.30
1	D	24	ASP	CB-CG-OD2	-5.24	113.58	118.30
1	C	63	ASP	CB-CG-OD1	5.13	122.92	118.30
1	A	182	LEU	CB-CG-CD1	-5.12	102.29	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	1930	0	1942	37	0
1	B	1875	0	1882	25	0
1	C	1915	0	1933	47	0
1	D	1907	0	1931	46	0
1	E	1910	0	1920	39	0
1	F	1850	0	1862	35	1
2	A	2	0	0	2	0
2	B	2	0	0	1	0
2	C	2	0	0	1	0
2	D	2	0	0	1	0
2	E	2	0	0	2	0
2	F	2	0	0	1	0
3	A	3	0	0	0	0
3	B	1	0	0	0	0
3	C	3	0	0	0	0
3	D	2	0	0	0	0
3	E	2	0	0	0	0
3	F	1	0	0	0	0
4	A	8	0	12	0	0
4	B	8	0	12	0	1
4	C	12	0	16	9	0
4	D	4	0	6	0	0
4	E	4	0	5	0	0
4	F	8	0	12	3	0
5	A	381	0	0	25	0
5	B	425	0	0	18	1
5	C	373	0	0	21	1
5	D	388	0	0	28	0
5	E	398	0	0	25	1
5	F	402	0	0	21	1
All	All	13822	0	11533	238	3

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 10.

All (238) 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:211:ALA:HB3	5:B:4411:HOH:O	1.29	1.32
1:D:14:ALA:HB2	5:D:4401:HOH:O	1.40	1.21
1:A:264:ILE:HG23	5:A:4372:HOH:O	1.41	1.19
1:F:120:GLU:HG3	5:F:4169:HOH:O	1.48	1.12

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:13:MET:HE1	5:A:4170:HOH:O	1.52	1.06
1:A:226:GLN:HB2	5:A:4379:HOH:O	1.61	0.99
1:C:17[B]:GLN:OE1	5:C:4325:HOH:O	1.80	0.99
1:C:194:GLU:HB2	4:C:3017:EDO:H12	1.45	0.95
1:C:16:MET:HE3	5:C:4301:HOH:O	1.68	0.94
1:C:192:PRO:HB3	4:C:3017:EDO:H22	1.50	0.94
1:F:14:ALA:HB2	5:F:4401:HOH:O	1.69	0.93
1:B:143:ALA:HA	5:B:4437:HOH:O	1.70	0.92
1:E:13:MET:HE1	5:E:4378:HOH:O	1.71	0.90
1:D:226:GLN:NE2	5:D:4094:HOH:O	2.04	0.88
1:A:226:GLN:OE1	5:A:4169:HOH:O	1.93	0.87
1:F:253[B]:GLU:OE1	5:F:4314:HOH:O	1.92	0.86
1:C:17[B]:GLN:HE22	1:C:79:MET:CE	1.89	0.86
1:C:17[B]:GLN:HE22	1:C:79:MET:HE1	1.43	0.84
1:B:126:VAL:HG22	5:B:4377:HOH:O	1.79	0.83
1:B:232:VAL:HG22	5:B:4411:HOH:O	1.78	0.82
1:B:125:PHE:HB2	5:B:4437:HOH:O	1.79	0.82
2:B:3007:CL:CL	5:B:4100:HOH:O	2.35	0.81
1:C:16:MET:HE1	5:C:4336:HOH:O	1.83	0.78
1:E:10:LYS:HE2	5:E:4437:HOH:O	1.83	0.78
1:D:45:LEU:HD21	1:D:81:LYS:HE3	1.65	0.77
1:B:72:ASN:HB3	5:B:4433:HOH:O	1.85	0.77
1:A:179:GLU:HG2	5:A:4238:HOH:O	1.85	0.76
1:E:66:ILE:HB	5:E:4397:HOH:O	1.84	0.76
1:F:194:GLU:H	4:F:3034:EDO:H22	1.50	0.76
1:F:14:ALA:HA	5:F:4392:HOH:O	1.85	0.76
1:E:66:ILE:HD12	5:E:4397:HOH:O	1.85	0.75
1:A:79:MET:HE3	5:A:4359:HOH:O	1.85	0.74
1:A:197:LEU:HD12	5:A:4349:HOH:O	1.89	0.73
1:C:7:GLU:N	5:C:4280:HOH:O	2.22	0.72
1:E:14:ALA:HB2	5:E:4363:HOH:O	1.88	0.72
1:F:46:GLU:C	5:F:4347:HOH:O	2.27	0.72
1:D:8:ARG:N	5:D:4397:HOH:O	2.22	0.72
2:E:3025:CL:CL	5:E:4141:HOH:O	2.44	0.72
1:D:197:LEU:HD12	5:D:4319:HOH:O	1.89	0.71
2:E:3026:CL:CL	5:E:4428:HOH:O	2.44	0.71
1:A:14:ALA:HA	1:A:17[B]:GLN:HE21	1.55	0.71
2:C:3013:CL:CL	5:C:4107:HOH:O	2.46	0.70
1:B:45:LEU:HD21	1:B:81:LYS:HE2	1.73	0.69
1:E:13:MET:HA	1:E:16[A]:MET:CE	2.23	0.69
1:E:197:LEU:HD12	5:E:4392:HOH:O	1.92	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:3001:CL:CL	5:A:4103:HOH:O	2.47	0.69
1:E:36:GLU:HB2	5:E:4274:HOH:O	1.91	0.68
1:C:45:LEU:HD21	1:C:81:LYS:HE3	1.75	0.68
1:C:194:GLU:CB	4:C:3017:EDO:H12	2.22	0.67
1:D:240:LYS:NZ	5:D:4366:HOH:O	2.26	0.67
1:D:188:ASN:ND2	5:D:4212:HOH:O	2.27	0.67
1:A:238:ILE:HG12	5:A:4290:HOH:O	1.95	0.67
1:A:218:PRO:CB	5:A:4372:HOH:O	2.44	0.66
1:D:188:ASN:HB2	5:D:4419:HOH:O	1.96	0.65
1:E:13:MET:HA	1:E:16[A]:MET:HE3	1.79	0.65
2:F:3030:CL:CL	5:F:4096:HOH:O	2.51	0.65
1:A:194:GLU:OE2	5:A:4121:HOH:O	2.15	0.65
1:E:46:GLU:HG3	1:E:66:ILE:HD12	1.78	0.64
1:A:45:LEU:HD21	1:A:81:LYS:HE3	1.80	0.64
1:F:13:MET:CA	5:F:4313:HOH:O	2.46	0.64
1:F:183:MET:HG3	5:F:4379:HOH:O	1.97	0.63
1:C:24:ASP:OD1	4:C:3019:EDO:O1	2.17	0.62
1:D:262:LYS:NZ	5:D:4216:HOH:O	2.27	0.62
1:E:197:LEU:CD1	5:E:4392:HOH:O	2.45	0.62
1:D:234:VAL:HG21	1:D:238:ILE:HD11	1.82	0.61
1:B:17:GLN:HG3	5:B:4410:HOH:O	1.99	0.61
1:F:150:GLY:H	1:F:163:HIS:HD2	1.48	0.61
1:D:16:MET:HE3	5:D:4391:HOH:O	1.99	0.61
1:A:234:VAL:HG21	1:A:238:ILE:HD11	1.83	0.60
1:C:17[B]:GLN:NE2	1:C:79:MET:HE1	2.15	0.60
1:A:81:LYS:NZ	5:A:4087:HOH:O	2.29	0.59
1:D:14:ALA:O	1:D:17:GLN:HG2	2.01	0.59
1:D:130:ARG:HG3	1:D:149:LYS:HD2	1.85	0.59
1:B:18:LYS:HD2	5:B:4417:HOH:O	2.02	0.59
1:C:192:PRO:CB	4:C:3017:EDO:H22	2.30	0.59
1:E:46:GLU:HG3	1:E:66:ILE:CD1	2.34	0.58
1:A:218:PRO:HB2	5:A:4372:HOH:O	2.03	0.58
1:C:120:GLU:HG3	5:C:4292:HOH:O	2.03	0.58
2:D:3020:CL:CL	5:D:4113:HOH:O	2.54	0.58
1:F:192:PRO:HB3	4:F:3034:EDO:H12	1.85	0.58
1:C:240:LYS:HD3	1:C:240:LYS:N	2.19	0.57
1:E:10:LYS:HG2	5:E:4437:HOH:O	2.04	0.57
1:A:214:GLY:HA2	5:A:4301:HOH:O	2.06	0.56
1:E:256:THR:HB	5:E:4441:HOH:O	2.05	0.55
1:F:18:LYS:HB3	5:F:4371:HOH:O	2.05	0.55
1:E:124:PRO:HA	5:E:4437:HOH:O	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:46:GLU:HG3	1:A:66:ILE:HD12	1.89	0.55
1:A:46:GLU:HG3	1:A:66:ILE:CD1	2.37	0.55
1:D:13:MET:HA	1:D:16:MET:HE2	1.89	0.55
1:B:144:SER:N	5:B:4437:HOH:O	2.38	0.54
1:D:271:LEU:HD13	5:D:4361:HOH:O	2.06	0.54
1:B:15:GLU:O	1:B:18:LYS:HB2	2.07	0.54
1:E:13:MET:O	1:E:16[B]:MET:HG2	2.07	0.54
1:C:191:ALA:O	4:C:3018:EDO:C1	2.55	0.54
1:E:10:LYS:CD	5:E:4437:HOH:O	2.55	0.54
1:E:6:THR:HG23	1:E:9:VAL:H	1.71	0.54
1:B:46:GLU:HB3	5:B:4416:HOH:O	2.08	0.54
1:E:16[B]:MET:CE	5:E:4353:HOH:O	2.56	0.54
1:E:22:ILE:HG12	1:E:41:ALA:HB3	1.90	0.54
1:E:234:VAL:HG21	1:E:238:ILE:HD11	1.91	0.53
1:B:13:MET:O	1:B:16[B]:MET:CG	2.56	0.53
1:F:45:LEU:HD21	1:F:81:LYS:HE3	1.91	0.53
1:E:13:MET:HA	1:E:16[A]:MET:HE2	1.91	0.53
1:B:194:GLU:HG3	5:B:4423:HOH:O	2.09	0.52
1:E:188:ASN:ND2	5:E:4232:HOH:O	2.42	0.52
1:F:150:GLY:H	1:F:163:HIS:CD2	2.26	0.52
1:F:57:GLY:N	5:F:4382:HOH:O	2.42	0.52
1:C:226:GLN:NE2	5:C:4059:HOH:O	2.37	0.52
1:C:198:GLN:NE2	1:C:202:ASP:OD1	2.43	0.52
1:A:226:GLN:NE2	5:A:4129:HOH:O	2.42	0.51
1:D:17:GLN:NE2	1:D:145:MET:SD	2.83	0.51
1:D:201:LYS:HG3	5:D:4319:HOH:O	2.10	0.51
1:F:17:GLN:CD	5:F:4392:HOH:O	2.49	0.51
1:C:246:LYS:NZ	5:C:4395:HOH:O	2.43	0.51
1:C:81:LYS:NZ	4:C:3019:EDO:H12	2.25	0.51
1:F:130:ARG:HG2	1:F:163:HIS:CE1	2.45	0.51
2:A:3001:CL:CL	5:A:4328:HOH:O	2.56	0.51
1:C:13:MET:HG2	5:C:4379:HOH:O	2.09	0.51
1:D:47:ARG:CB	5:D:4244:HOH:O	2.59	0.51
1:C:10:LYS:HA	1:C:13:MET:HE3	1.93	0.50
1:A:68:GLU:HA	1:A:71:MET:HE2	1.93	0.50
1:D:45:LEU:CD2	1:D:81:LYS:HE3	2.37	0.50
1:D:168:ASN:CG	5:D:4356:HOH:O	2.49	0.50
1:B:249:LYS:O	1:B:253[B]:GLU:HG2	2.11	0.50
1:A:55:ALA:N	5:A:4302:HOH:O	2.44	0.50
1:B:127:CYS:CA	5:B:4377:HOH:O	2.60	0.50
1:C:100:TYR:HD2	5:C:4320:HOH:O	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:240:LYS:NZ	5:C:4262:HOH:O	2.24	0.49
1:E:10:LYS:CE	5:E:4437:HOH:O	2.50	0.49
1:D:100:TYR:HD2	5:D:4420:HOH:O	1.94	0.49
1:C:46:GLU:O	1:C:47:ARG:CB	2.61	0.49
1:D:57:GLY:HA3	5:D:4272:HOH:O	2.12	0.49
1:C:11:ARG:HB3	5:C:4305:HOH:O	2.11	0.49
1:A:218:PRO:HB3	5:A:4372:HOH:O	2.11	0.49
1:C:47:ARG:O	5:C:4289:HOH:O	2.20	0.49
1:E:29:GLU:HG3	5:E:4446:HOH:O	2.13	0.49
1:B:232:VAL:HA	5:B:4411:HOH:O	2.12	0.48
1:F:17:GLN:OE1	5:F:4392:HOH:O	2.20	0.48
1:C:186:ALA:HB1	4:C:3018:EDO:H12	1.96	0.48
1:D:197:LEU:CD1	5:D:4319:HOH:O	2.54	0.48
1:F:14:ALA:HB3	5:F:4386:HOH:O	2.14	0.48
1:D:241:SER:HB2	5:D:4412:HOH:O	2.13	0.48
1:D:234:VAL:CG2	1:D:238:ILE:HD11	2.43	0.48
1:A:157:ILE:HD13	1:A:215:VAL:HA	1.96	0.47
1:D:57:GLY:CA	5:D:4272:HOH:O	2.61	0.47
1:F:163:HIS:HE1	5:F:4136:HOH:O	1.97	0.47
1:B:153:GLY:C	5:B:4345:HOH:O	2.51	0.47
1:F:234:VAL:HG21	1:F:238:ILE:HD11	1.96	0.47
1:F:68:GLU:HA	1:F:71:MET:CE	2.45	0.47
1:D:120:GLU:HG3	5:D:4308:HOH:O	2.15	0.47
1:E:45:LEU:HD21	1:E:81:LYS:HE2	1.97	0.47
1:F:188:ASN:HB3	5:F:4404:HOH:O	2.14	0.47
1:D:100:TYR:HE2	5:D:4401:HOH:O	1.97	0.47
1:D:14:ALA:O	1:D:17:GLN:CG	2.62	0.47
1:C:214:GLY:HA2	5:C:4216:HOH:O	2.15	0.47
1:D:182:LEU:HD22	1:D:196:LEU:HD13	1.96	0.47
1:D:13:MET:HG3	5:D:4388:HOH:O	2.14	0.46
1:F:174:VAL:HG13	1:F:182:LEU:HD21	1.96	0.46
1:A:270:GLU:HB2	5:A:4331:HOH:O	2.16	0.46
1:A:14:ALA:HA	1:A:17[B]:GLN:NE2	2.27	0.46
1:A:194:GLU:HB2	5:A:4175:HOH:O	2.16	0.46
1:B:151:GLU:O	5:B:4270:HOH:O	2.21	0.46
1:E:101:ILE:HD12	1:E:123:VAL:HG11	1.97	0.46
1:E:67:VAL:HG12	1:E:71:MET:CE	2.45	0.46
1:F:214:GLY:HA2	5:F:4140:HOH:O	2.15	0.46
1:C:17[B]:GLN:CD	5:C:4325:HOH:O	2.42	0.45
1:C:191:ALA:O	4:C:3018:EDO:H11	2.16	0.45
1:F:68:GLU:HA	1:F:71:MET:HE2	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:18:LYS:HE2	5:B:4279:HOH:O	2.15	0.45
1:A:151:GLU:HG2	1:A:154:THR:HG21	1.98	0.45
1:D:175:VAL:O	1:D:200:LYS:NZ	2.49	0.45
1:D:18:LYS:N	5:D:4289:HOH:O	2.48	0.45
1:A:130:ARG:HG3	1:A:149:LYS:HD2	1.98	0.45
1:F:253[B]:GLU:CD	5:F:4314:HOH:O	2.50	0.45
1:D:68:GLU:HA	1:D:71:MET:HE2	1.99	0.45
1:A:55:ALA:N	5:A:4368:HOH:O	2.50	0.45
1:C:150:GLY:H	1:C:163:HIS:HD2	1.63	0.45
1:D:270:GLU:HG2	5:E:4221:HOH:O	2.16	0.45
1:B:40:VAL:O	1:B:77:PRO:HD2	2.16	0.44
1:C:234:VAL:HG21	1:C:238:ILE:HD11	1.99	0.44
1:C:40:VAL:O	1:C:77:PRO:HD2	2.17	0.44
1:D:145:MET:HG3	1:D:208:VAL:HG23	2.00	0.44
1:C:201:LYS:NZ	5:C:4318:HOH:O	2.50	0.44
1:F:194:GLU:HB2	4:F:3034:EDO:O1	2.17	0.44
1:A:178:SER:OG	1:A:180:ASP:OD1	2.31	0.44
1:A:13:MET:CE	5:A:4170:HOH:O	2.33	0.44
1:C:17[B]:GLN:NE2	1:C:79:MET:CE	2.70	0.44
1:D:71:MET:HA	1:D:78:VAL:HG21	1.99	0.44
1:E:79:MET:HG2	1:E:100:TYR:HB2	2.00	0.44
1:F:197:LEU:HD12	5:F:4403:HOH:O	2.18	0.44
1:A:55:ALA:HB2	5:A:4368:HOH:O	2.18	0.43
1:C:67:VAL:O	1:C:71:MET:HG3	2.18	0.43
1:E:36:GLU:HB3	5:E:4431:HOH:O	2.18	0.43
1:F:57:GLY:HA2	5:F:4382:HOH:O	2.18	0.43
1:C:178:SER:OG	1:C:180:ASP:OD1	2.36	0.43
1:D:16:MET:CE	5:D:4391:HOH:O	2.62	0.43
1:F:194:GLU:OE2	5:F:4080:HOH:O	2.21	0.43
1:C:198:GLN:HG2	5:C:4125:HOH:O	2.19	0.43
1:D:157:ILE:HD13	1:D:215:VAL:HA	2.01	0.43
1:D:240:LYS:CD	1:D:240:LYS:N	2.82	0.43
1:E:16[B]:MET:SD	5:E:4353:HOH:O	2.62	0.43
1:D:122:THR:HG23	5:D:4176:HOH:O	2.19	0.43
1:D:201:LYS:HE3	5:D:4319:HOH:O	2.18	0.43
1:E:101:ILE:CD1	1:E:123:VAL:HG11	2.48	0.43
1:E:151:GLU:HG2	1:E:154:THR:HG21	2.01	0.43
1:E:67:VAL:HG12	1:E:71:MET:HE2	2.02	0.42
1:C:174:VAL:HG13	1:C:182:LEU:HD21	2.00	0.42
1:C:57:GLY:N	5:C:4281:HOH:O	2.51	0.42
1:B:67:VAL:HG12	1:B:71:MET:CE	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:225:MET:HE2	5:C:4155:HOH:O	2.20	0.42
1:C:57:GLY:HA3	5:C:4374:HOH:O	2.19	0.42
1:D:9:VAL:O	1:D:13:MET:HG2	2.19	0.42
1:B:241:SER:HB2	5:B:4395:HOH:O	2.19	0.42
1:B:250:ALA:HA	1:B:267:LEU:HD22	2.01	0.42
1:B:22:ILE:HG12	1:B:41:ALA:HB3	2.01	0.42
1:E:63:ASP:HB3	5:E:4397:HOH:O	2.19	0.42
1:A:17[B]:GLN:OE1	1:A:22:ILE:HD11	2.20	0.42
1:D:240:LYS:N	1:D:240:LYS:HD3	2.35	0.42
1:F:157:ILE:HD13	1:F:215:VAL:HA	2.00	0.42
1:A:253[B]:GLU:CD	5:A:4374:HOH:O	2.59	0.41
1:C:147:ARG:HA	1:C:210:PHE:O	2.20	0.41
1:E:10:LYS:CG	5:E:4437:HOH:O	2.66	0.41
1:D:10:LYS:HE2	5:D:4095:HOH:O	2.20	0.41
1:A:45:LEU:HD21	1:A:81:LYS:CE	2.50	0.41
1:E:270:GLU:HB2	5:E:4429:HOH:O	2.21	0.41
1:C:249:LYS:HG2	5:C:4349:HOH:O	2.20	0.41
1:F:17:GLN:HE22	1:F:79:MET:CE	2.34	0.41
1:E:130:ARG:HG2	1:E:163:HIS:CE1	2.56	0.41
1:F:57:GLY:CA	5:F:4382:HOH:O	2.68	0.41
1:E:17:GLN:NE2	1:E:100:TYR:OH	2.54	0.40
1:C:79:MET:HG2	1:C:100:TYR:HB2	2.04	0.40
1:F:40:VAL:O	1:F:77:PRO:HD2	2.22	0.40
1:D:200:LYS:NZ	5:D:4359:HOH:O	2.53	0.40
1:A:153:GLY:C	5:A:4319:HOH:O	2.59	0.40
1:A:157:ILE:HD11	1:A:215:VAL:HG22	2.02	0.40
1:C:157:ILE:HD13	1:C:215:VAL:HA	2.04	0.40

All (3) 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)
5:C:4168:HOH:O	5:E:4131:HOH:O[2_665]	2.10	0.10
1:F:188:ASN:O	4:B:3011:EDO:O1[2_665]	2.14	0.06
5:B:4100:HOH:O	5:F:4379:HOH:O[2_665]	2.18	0.02

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	257/305 (84%)	254 (99%)	3 (1%)	0	100	100
1	B	250/305 (82%)	249 (100%)	1 (0%)	0	100	100
1	C	254/305 (83%)	250 (98%)	4 (2%)	0	100	100
1	D	253/305 (83%)	250 (99%)	3 (1%)	0	100	100
1	E	254/305 (83%)	252 (99%)	2 (1%)	0	100	100
1	F	247/305 (81%)	245 (99%)	2 (1%)	0	100	100
All	All	1515/1830 (83%)	1500 (99%)	15 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	198/237 (84%)	198 (100%)	0	100	100
1	B	190/237 (80%)	189 (100%)	1 (0%)	88	92
1	C	198/237 (84%)	197 (100%)	1 (0%)	88	92
1	D	196/237 (83%)	196 (100%)	0	100	100
1	E	197/237 (83%)	196 (100%)	1 (0%)	88	92
1	F	189/237 (80%)	188 (100%)	1 (0%)	88	92
All	All	1168/1422 (82%)	1164 (100%)	4 (0%)	92	95

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	247	PHE
1	C	240	LYS
1	E	270	GLU
1	F	196	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	209	ASN
1	B	226	GLN
1	C	72	ASN
1	C	163	HIS
1	C	209	ASN
1	D	72	ASN
1	D	188	ASN
1	E	17	GLN
1	E	188	ASN
1	E	226	GLN
1	F	17	GLN
1	F	72	ASN
1	F	163	HIS
1	F	188	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 35 ligands modelled in this entry, 24 are monoatomic - leaving 11 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$
4	EDO	C	3018	-	3,3,3	0.63	0	2,2,2	0.89	0
4	EDO	F	3034	-	3,3,3	0.51	0	2,2,2	0.64	0
4	EDO	B	3012	-	3,3,3	0.28	0	2,2,2	0.57	0
4	EDO	D	3024	-	3,3,3	0.33	0	2,2,2	0.69	0
4	EDO	A	3006	-	3,3,3	0.40	0	2,2,2	0.39	0
4	EDO	C	3017	-	3,3,3	0.80	0	2,2,2	0.29	0
4	EDO	E	3029	-	3,3,3	0.74	0	2,2,2	0.89	0
4	EDO	B	3011	-	3,3,3	0.41	0	2,2,2	0.60	0
4	EDO	C	3019	-	3,3,3	0.27	0	2,2,2	0.77	0
4	EDO	A	3005	-	3,3,3	0.17	0	2,2,2	0.24	0
4	EDO	F	3035	-	3,3,3	0.39	0	2,2,2	0.85	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
4	EDO	C	3018	-	-	0/1/1/1	-
4	EDO	F	3034	-	-	1/1/1/1	-
4	EDO	B	3012	-	-	0/1/1/1	-
4	EDO	D	3024	-	-	0/1/1/1	-
4	EDO	A	3006	-	-	0/1/1/1	-
4	EDO	C	3017	-	-	1/1/1/1	-
4	EDO	E	3029	-	-	1/1/1/1	-
4	EDO	B	3011	-	-	1/1/1/1	-
4	EDO	C	3019	-	-	0/1/1/1	-
4	EDO	A	3005	-	-	1/1/1/1	-
4	EDO	F	3035	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	3011	EDO	O1-C1-C2-O2
4	F	3034	EDO	O1-C1-C2-O2
4	A	3005	EDO	O1-C1-C2-O2
4	E	3029	EDO	O1-C1-C2-O2
4	F	3035	EDO	O1-C1-C2-O2
4	C	3017	EDO	O1-C1-C2-O2

There are no ring outliers.

5 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	3018	EDO	3	0
4	F	3034	EDO	3	0
4	C	3017	EDO	4	0
4	B	3011	EDO	0	1
4	C	3019	EDO	2	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	259/305 (84%)	-0.25	11 (4%) 36 41	28, 34, 63, 78	0
1	B	252/305 (82%)	-0.42	5 (1%) 65 68	27, 34, 58, 82	0
1	C	256/305 (83%)	-0.32	8 (3%) 49 54	28, 34, 59, 70	0
1	D	256/305 (83%)	-0.30	8 (3%) 49 54	27, 34, 62, 77	0
1	E	256/305 (83%)	-0.35	6 (2%) 60 64	27, 34, 61, 74	0
1	F	250/305 (81%)	-0.32	7 (2%) 53 58	28, 34, 54, 85	0
All	All	1529/1830 (83%)	-0.33	45 (2%) 51 56	27, 34, 60, 85	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	12	GLY	6.0
1	A	56	GLY	5.2
1	A	6	THR	4.9
1	A	55	ALA	4.3
1	E	13	MET	3.9
1	E	270	GLU	3.8
1	C	12	GLY	3.6
1	E	6	THR	3.5
1	F	271	LEU	3.5
1	A	9	VAL	3.3
1	E	16[A]	MET	3.2
1	F	14	ALA	3.2
1	D	9	VAL	3.2
1	C	13	MET	3.1
1	F	16	MET	3.1
1	B	11	ARG	3.1
1	D	242	ASP	3.1
1	F	12	GLY	2.9
1	D	272	GLY	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	17	GLN	2.7
1	C	8	ARG	2.7
1	F	153	GLY	2.6
1	C	16	MET	2.6
1	D	271	LEU	2.5
1	E	242	ASP	2.5
1	A	29	GLU	2.5
1	B	270	GLU	2.5
1	D	13	MET	2.5
1	A	242	ASP	2.5
1	D	16	MET	2.4
1	A	8	ARG	2.4
1	C	270	GLU	2.3
1	D	240	LYS	2.3
1	D	17	GLN	2.3
1	E	36	GLU	2.3
1	F	13	MET	2.3
1	A	66	ILE	2.2
1	F	242	ASP	2.2
1	A	16	MET	2.1
1	C	29	GLU	2.1
1	C	7	GLU	2.1
1	A	7	GLU	2.1
1	C	242	ASP	2.0
1	B	14	ALA	2.0
1	A	13	MET	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. 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(\AA^2)	Q<0.9
3	MG	C	3016	1/1	0.62	0.16	65,65,65,65	0
4	EDO	B	3011	4/4	0.64	0.44	84,85,86,87	0
3	MG	A	3009	1/1	0.69	0.29	61,61,61,61	0
3	MG	E	3032	1/1	0.70	0.18	63,63,63,63	0
3	MG	D	3027	1/1	0.73	0.32	66,66,66,66	0
3	MG	A	3003	1/1	0.75	0.25	61,61,61,61	0
4	EDO	A	3006	4/4	0.75	0.28	73,74,74,75	0
3	MG	A	3004	1/1	0.76	0.23	67,67,67,67	0
3	MG	C	3022	1/1	0.80	0.25	61,61,61,61	0
3	MG	B	3010	1/1	0.80	0.26	58,58,58,58	0
4	EDO	F	3034	4/4	0.85	0.20	32,39,46,47	0
4	EDO	C	3017	4/4	0.86	0.29	37,39,48,50	0
3	MG	D	3023	1/1	0.87	0.25	64,64,64,64	0
4	EDO	C	3019	4/4	0.89	0.19	57,60,61,63	0
4	EDO	D	3024	4/4	0.90	0.23	38,43,46,53	0
4	EDO	A	3005	4/4	0.91	0.18	39,40,44,46	0
3	MG	E	3028	1/1	0.91	0.11	56,56,56,56	0
4	EDO	C	3018	4/4	0.91	0.16	44,45,47,50	0
4	EDO	B	3012	4/4	0.92	0.21	37,44,47,56	0
3	MG	F	3033	1/1	0.93	0.21	60,60,60,60	0
2	CL	D	3021	1/1	0.94	0.08	53,53,53,53	0
4	EDO	F	3035	4/4	0.94	0.12	41,43,45,48	0
4	EDO	E	3029	4/4	0.94	0.18	39,41,44,49	0
3	MG	C	3015	1/1	0.94	0.19	65,65,65,65	0
2	CL	B	3008	1/1	0.95	0.06	54,54,54,54	0
2	CL	E	3025	1/1	0.96	0.05	51,51,51,51	0
2	CL	C	3014	1/1	0.97	0.06	54,54,54,54	0
2	CL	E	3026	1/1	0.97	0.06	51,51,51,51	0
2	CL	F	3031	1/1	0.97	0.09	51,51,51,51	0
2	CL	A	3001	1/1	0.97	0.08	56,56,56,56	0
2	CL	A	3002	1/1	0.98	0.07	58,58,58,58	0
2	CL	C	3013	1/1	0.98	0.09	46,46,46,46	0
2	CL	D	3020	1/1	0.99	0.07	51,51,51,51	0
2	CL	B	3007	1/1	0.99	0.04	50,50,50,50	0
2	CL	F	3030	1/1	0.99	0.07	47,47,47,47	0

6.5 Other polymers ⓘ

There are no such residues in this entry.