



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 25, 2020 – 04:39 PM BST

PDB ID : 4HN0
Title : Crystal Structure of ChmJ, a 3'-monoepimerase apoenzyme from *Streptomyces bikiniensis*
Authors : Holden, H.M.; Kubiak, R.L.
Deposited on : 2012-10-18
Resolution : 2.20 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	:	2.13
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)	:	2.13

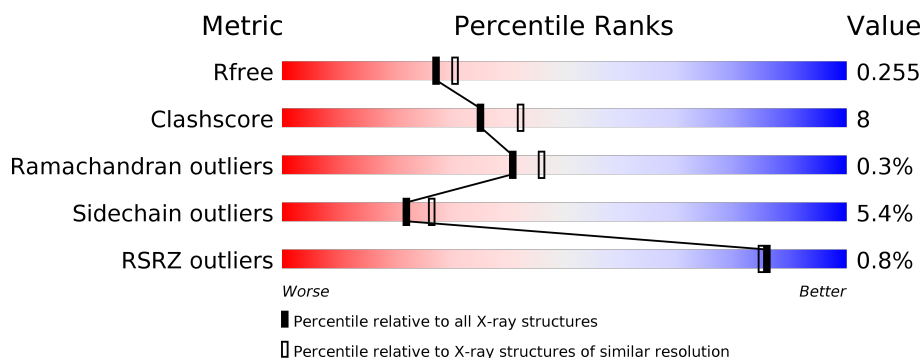
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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	198	<div> <div>%</div> <div> <div></div> <div>86%</div> <div>11%</div> <div>.</div> </div> </div>
1	B	198	<div> <div>85%</div> <div>15%</div> <div>.</div> </div>
1	C	198	<div> <div>%</div> <div> <div></div> <div>72%</div> <div>25%</div> <div>.</div> </div> </div>
1	D	198	<div> <div>2%</div> <div> <div></div> <div>73%</div> <div>25%</div> <div>..</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6456 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 Putative 3-epimerase in D-allose pathway.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	198	Total	C	N	O	S	0	0	0
			1549	974	279	292	4			
1	B	198	Total	C	N	O	S	0	0	0
			1549	974	279	292	4			
1	C	198	Total	C	N	O	S	0	0	0
			1549	974	279	292	4			
1	D	197	Total	C	N	O	S	0	1	0
			1543	971	278	290	4			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	197	LEU	-	EXPRESSION TAG	UNP Q5SFD1
A	198	GLU	-	EXPRESSION TAG	UNP Q5SFD1
B	197	LEU	-	EXPRESSION TAG	UNP Q5SFD1
B	198	GLU	-	EXPRESSION TAG	UNP Q5SFD1
C	197	LEU	-	EXPRESSION TAG	UNP Q5SFD1
C	198	GLU	-	EXPRESSION TAG	UNP Q5SFD1
D	197	LEU	-	EXPRESSION TAG	UNP Q5SFD1
D	198	GLU	-	EXPRESSION TAG	UNP Q5SFD1

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	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 CHLORIDE ION (three-letter code: CL) (formula: Cl).

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

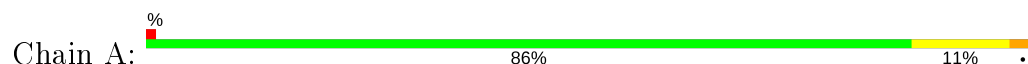
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	65	Total 65	O 65	0	0
4	B	63	Total 63	O 63	0	0
4	C	51	Total 51	O 51	0	0
4	D	35	Total 35	O 35	0	0

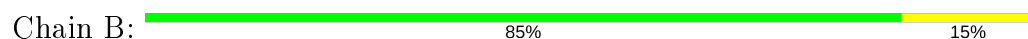
3 Residue-property plots

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.

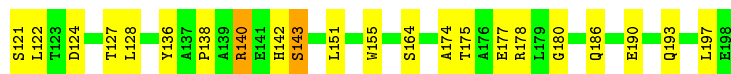
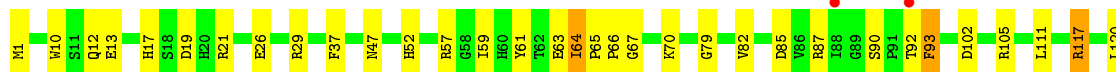
- Molecule 1: Putative 3-epimerase in D-allose pathway



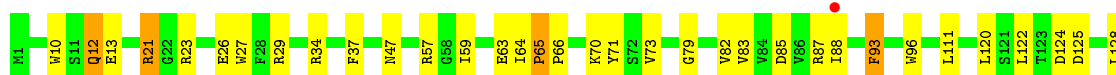
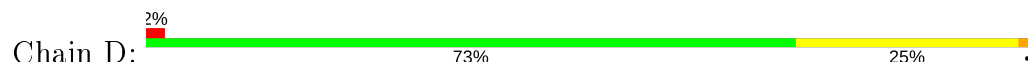
- Molecule 1: Putative 3-epimerase in D-allose pathway



- Molecule 1: Putative 3-epimerase in D-allose pathway



- Molecule 1: Putative 3-epimerase in D-allose pathway



4 Data and refinement statistics

Property	Value	Source
Space group	I 4	Depositor
Cell constants a, b, c, α , β , γ	141.59Å 141.59Å 115.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.20 27.76 – 2.20	Depositor EDS
% Data completeness (in resolution range)	97.8 (30.00-2.20) 97.9 (27.76-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	13.10 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.215 , 0.259 0.212 , 0.255	Depositor DCC
R_{free} test set	2917 reflections (5.16%)	wwPDB-VP
Wilson B-factor (Å ²)	44.9	Xtriage
Anisotropy	0.054	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 24.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.487 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6456	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.85% 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: 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	0.46	0/1592	1.14	5/2174 (0.2%)
1	B	0.47	0/1592	1.16	9/2174 (0.4%)
1	C	0.43	0/1592	1.14	6/2174 (0.3%)
1	D	0.42	0/1589	1.14	9/2170 (0.4%)
All	All	0.45	0/6365	1.15	29/8692 (0.3%)

There are no bond length outliers.

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	105	ARG	NE-CZ-NH1	7.85	124.23	120.30
1	D	23	ARG	NE-CZ-NH2	-7.13	116.73	120.30
1	A	21	ARG	NE-CZ-NH1	-6.95	116.82	120.30
1	A	140	ARG	NE-CZ-NH1	6.67	123.63	120.30
1	A	165	ASP	CB-CG-OD1	6.58	124.22	118.30
1	C	117	ARG	CG-CD-NE	-6.48	98.20	111.80
1	C	85	ASP	CB-CG-OD1	6.47	124.12	118.30
1	A	19	ASP	CB-CG-OD1	6.46	124.12	118.30
1	C	19	ASP	CB-CG-OD2	6.42	124.08	118.30
1	D	21	ARG	NE-CZ-NH1	6.41	123.50	120.30
1	C	57	ARG	NE-CZ-NH1	6.33	123.47	120.30
1	D	158	ASP	CB-CG-OD1	6.24	123.91	118.30
1	A	120	LEU	CB-CG-CD2	6.14	121.45	111.00
1	D	85	ASP	CB-CG-OD1	5.86	123.58	118.30
1	B	105	ARG	NE-CZ-NH2	-5.83	117.38	120.30
1	B	21	ARG	NE-CZ-NH1	-5.74	117.43	120.30
1	C	21	ARG	NE-CZ-NH1	5.70	123.15	120.30
1	B	125	ASP	CB-CG-OD1	5.68	123.41	118.30
1	D	57	ARG	NE-CZ-NH2	-5.60	117.50	120.30
1	B	165	ASP	CB-CG-OD1	5.48	123.23	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	23	ARG	NE-CZ-NH1	5.40	123.00	120.30
1	D	150	ASP	CB-CG-OD1	5.38	123.14	118.30
1	B	124	ASP	CB-CG-OD2	5.28	123.06	118.30
1	C	124	ASP	CB-CG-OD1	5.20	122.98	118.30
1	B	122	LEU	CB-CG-CD2	-5.12	102.30	111.00
1	B	57	ARG	NE-CZ-NH2	-5.09	117.75	120.30
1	D	128	LEU	CA-CB-CG	-5.09	103.59	115.30
1	B	128	LEU	CA-CB-CG	-5.04	103.70	115.30
1	D	57	ARG	NE-CZ-NH1	5.02	122.81	120.30

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	1549	0	1481	21	0
1	B	1549	0	1481	19	0
1	C	1549	0	1478	44	0
1	D	1543	0	1477	29	0
2	A	16	0	24	3	0
2	B	20	0	30	3	0
2	C	4	0	6	0	0
2	D	8	0	12	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	2	0	0	0	0
4	A	65	0	0	1	0
4	B	63	0	0	0	0
4	C	51	0	0	0	0
4	D	35	0	0	0	0
All	All	6456	0	5989	100	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 8.

All (100) 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:D:70:LYS:NZ	1:D:132:CYS:SG	1.97	1.36
1:C:70:LYS:HE3	1:C:117:ARG:HH22	1.04	1.14
1:C:70:LYS:HE3	1:C:117:ARG:NH2	1.75	1.01
1:D:65:PRO:HB3	1:D:66:PRO:HA	1.46	0.96
1:A:17:HIS:HD2	1:C:17:HIS:HD2	1.07	0.93
1:A:17:HIS:HD2	1:C:17:HIS:CD2	1.85	0.93
1:B:10:TRP:CZ3	2:B:205:EDO:H11	2.15	0.81
1:A:95:ARG:HB2	2:A:204:EDO:H11	1.64	0.79
1:A:17:HIS:CD2	1:C:17:HIS:CD2	2.72	0.78
1:C:186:GLN:HG2	1:C:190:GLU:OE2	1.83	0.78
1:D:71:TYR:OH	2:D:202:EDO:H22	1.88	0.74
1:A:17:HIS:CD2	1:C:17:HIS:HD2	1.99	0.73
1:D:65:PRO:CB	1:D:66:PRO:HA	2.19	0.73
1:D:196:ARG:N	1:D:196:ARG:HD3	2.04	0.72
1:A:70:LYS:HE2	1:A:132:CYS:SG	2.31	0.71
1:B:70:LYS:HE2	1:B:132:CYS:SG	2.32	0.70
1:B:70:LYS:NZ	1:B:117:ARG:NH2	2.40	0.69
1:C:1:MET:HE3	1:C:12:GLN:HB3	1.75	0.68
1:A:26:GLU:HA	1:B:47:ASN:HD22	1.60	0.66
1:A:2:HIS:NE2	4:A:301:HOH:O	2.29	0.65
1:D:193:GLN:O	1:D:196:ARG:HG2	1.97	0.65
1:D:140:ARG:HH11	1:D:140:ARG:HG3	1.66	0.61
1:C:186:GLN:O	1:C:190:GLU:HG3	2.00	0.61
1:A:47:ASN:HD22	1:B:26:GLU:HA	1.66	0.60
1:C:70:LYS:HG3	1:C:117:ARG:NH2	2.15	0.60
1:B:47:ASN:HB2	2:B:203:EDO:H12	1.83	0.59
1:C:47:ASN:HD22	1:D:26:GLU:HA	1.67	0.59
1:C:186:GLN:CG	1:C:190:GLU:OE2	2.50	0.59
1:D:88:ILE:HD11	1:D:182:LEU:HB2	1.85	0.59
1:D:191:GLN:O	1:D:195:GLN:HG3	2.04	0.58
1:C:79:GLY:HA2	1:C:122:LEU:HG	1.85	0.57
1:A:76:ARG:HG3	1:B:76:ARG:HD2	1.86	0.57
1:C:26:GLU:HA	1:D:47:ASN:HD22	1.69	0.56
1:C:177:GLU:O	1:C:180:GLY:N	2.31	0.56
1:C:87:ARG:O	1:C:93:PHE:HB2	2.05	0.56
1:D:196:ARG:H	1:D:196:ARG:HD3	1.73	0.54
1:C:70:LYS:CE	1:C:117:ARG:NH2	2.61	0.53
1:C:70:LYS:CE	1:C:117:ARG:HH22	1.97	0.53
1:A:1:MET:HA	1:A:12:GLN:HE21	1.74	0.52
1:A:70:LYS:NZ	1:A:117:ARG:NH1	2.58	0.52
1:B:47:ASN:HB2	2:B:203:EDO:C1	2.39	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:87:ARG:O	1:D:93:PHE:HD1	1.92	0.52
1:C:64:ILE:HA	1:C:65:PRO:C	2.29	0.51
1:A:95:ARG:HE	2:A:204:EDO:H12	1.74	0.51
1:D:10:TRP:CE2	1:D:37:PHE:HE1	2.28	0.51
1:A:45:GLN:HE21	1:A:47:ASN:HD21	1.58	0.50
1:C:175:THR:HG22	1:C:175:THR:O	2.10	0.50
1:D:63:GLU:OE1	1:D:140:ARG:NH2	2.45	0.49
1:D:79:GLY:HA2	1:D:122:LEU:HG	1.95	0.49
1:D:65:PRO:CB	1:D:66:PRO:CA	2.90	0.48
1:A:26:GLU:HA	1:B:47:ASN:ND2	2.27	0.48
1:D:186:GLN:HG2	1:D:190:GLU:OE2	2.14	0.48
1:A:95:ARG:HE	2:A:204:EDO:C1	2.28	0.47
1:C:61:TYR:CZ	1:C:142:HIS:HB2	2.51	0.47
1:D:174:ALA:O	1:D:177:GLU:HB3	2.15	0.46
1:C:90:SER:O	1:C:92:THR:N	2.49	0.46
1:B:1:MET:HA	1:B:12:GLN:HE21	1.82	0.44
1:C:52:HIS:CD2	1:C:52:HIS:N	2.84	0.44
1:B:145:ASN:ND2	1:B:168:GLU:O	2.50	0.44
1:C:174:ALA:O	1:C:178:ARG:HG3	2.17	0.44
1:D:189:GLN:O	1:D:193:GLN:NE2	2.51	0.44
1:A:47:ASN:ND2	1:B:26:GLU:HA	2.32	0.44
1:B:1:MET:HG2	1:B:12:GLN:HE21	1.83	0.43
1:C:164:SER:HB3	1:D:21:ARG:CZ	2.48	0.43
1:B:59:ILE:O	1:B:143:SER:HA	2.18	0.43
1:C:90:SER:C	1:C:92:THR:N	2.71	0.43
1:D:27:TRP:CE2	1:D:73:VAL:HG12	2.53	0.43
1:C:136:TYR:CZ	1:C:138:PRO:HG3	2.54	0.43
1:B:45:GLN:HE21	1:B:47:ASN:HD21	1.67	0.42
1:C:63:GLU:HB2	1:C:140:ARG:HG3	1.99	0.42
1:B:70:LYS:NZ	1:B:117:ARG:CZ	2.82	0.42
1:D:83:VAL:O	1:D:96:TRP:HA	2.20	0.42
1:D:87:ARG:O	1:D:93:PHE:HB2	2.19	0.42
1:C:175:THR:O	1:C:175:THR:CG2	2.66	0.42
1:C:82:VAL:HG21	1:C:155:TRP:CH2	2.55	0.42
1:B:51:SER:HB3	1:B:55:ALA:HB3	2.02	0.41
1:C:26:GLU:HG3	1:D:47:ASN:ND2	2.35	0.41
1:A:160:GLU:HA	1:A:161:PRO:HD3	1.81	0.41
1:C:117:ARG:O	1:C:117:ARG:HG3	2.20	0.41
1:C:66:PRO:HD2	1:C:67:GLY:H	1.84	0.41
1:D:12:GLN:HB2	1:D:12:GLN:HE21	1.53	0.41
1:C:63:GLU:OE1	1:C:140:ARG:NE	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:66:PRO:CD	1:C:67:GLY:N	2.83	0.41
1:B:151:LEU:HD23	1:B:151:LEU:HA	1.86	0.41
1:D:59:ILE:O	1:D:143:SER:HA	2.20	0.41
1:B:148:ASP:HB3	1:B:151:LEU:HB2	2.03	0.41
1:D:82:VAL:HG21	1:D:155:TRP:CH2	2.56	0.41
1:A:194:ALA:O	1:A:198:GLU:HG3	2.21	0.41
1:C:10:TRP:N	1:C:10:TRP:CD1	2.88	0.41
1:C:59:ILE:O	1:C:143:SER:HA	2.21	0.41
1:C:87:ARG:O	1:C:93:PHE:HD1	2.04	0.41
1:A:151:LEU:HD23	1:A:151:LEU:HA	1.75	0.41
1:C:66:PRO:CD	1:C:67:GLY:H	2.34	0.41
1:C:87:ARG:O	1:C:93:PHE:CD1	2.74	0.41
1:C:102:ASP:OD2	1:C:105:ARG:NH1	2.54	0.40
1:D:124:ASP:O	1:D:125:ASP:HB2	2.22	0.40
1:C:127:THR:C	1:C:128:LEU:HD12	2.42	0.40
1:C:66:PRO:HD3	1:C:140:ARG:HG2	2.03	0.40
1:A:59:ILE:O	1:A:143:SER:HA	2.22	0.40
1:C:10:TRP:CE2	1:C:37:PHE:HE1	2.39	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	196/198 (99%)	192 (98%)	4 (2%)	0	100	100
1	B	196/198 (99%)	193 (98%)	3 (2%)	0	100	100
1	C	196/198 (99%)	186 (95%)	9 (5%)	1 (0%)	29	31
1	D	196/198 (99%)	188 (96%)	7 (4%)	1 (0%)	29	31
All	All	784/792 (99%)	759 (97%)	23 (3%)	2 (0%)	41	46

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	93	PHE
1	D	93	PHE

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	162/162 (100%)	154 (95%)	8 (5%)	25	31
1	B	162/162 (100%)	157 (97%)	5 (3%)	40	51
1	C	162/162 (100%)	151 (93%)	11 (7%)	16	17
1	D	162/162 (100%)	151 (93%)	11 (7%)	16	17
All	All	648/648 (100%)	613 (95%)	35 (5%)	22	26

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	76	ARG
1	A	111	LEU
1	A	120	LEU
1	A	140	ARG
1	A	143	SER
1	A	151	LEU
1	A	157	ASP
1	A	178	ARG
1	B	29	ARG
1	B	121	SER
1	B	144	VAL
1	B	151	LEU
1	B	157	ASP
1	C	13	GLU
1	C	29	ARG
1	C	64	ILE
1	C	111	LEU
1	C	120	LEU
1	C	121	SER
1	C	140	ARG

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Mol	Chain	Res	Type
1	C	143	SER
1	C	151	LEU
1	C	193	GLN
1	C	197	LEU
1	D	12	GLN
1	D	13	GLU
1	D	29	ARG
1	D	34	ARG
1	D	64	ILE
1	D	65	PRO
1	D	111	LEU
1	D	120	LEU
1	D	136	TYR
1	D	151	LEU
1	D	166	ARG

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	12	GLN
1	A	17	HIS
1	A	45	GLN
1	A	47	ASN
1	B	12	GLN
1	B	45	GLN
1	B	47	ASN
1	C	17	HIS
1	C	47	ASN
1	C	52	HIS
1	D	12	GLN
1	D	39	HIS
1	D	47	ASN
1	D	186	GLN

5.3.3 RNA ⓘ

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 16 ligands modelled in this entry, 4 are monoatomic - leaving 12 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	$\# Z > 2$
2	EDO	D	201	-	3,3,3	0.46	0	2,2,2	1.04	0
2	EDO	B	201	-	3,3,3	0.42	0	2,2,2	0.55	0
2	EDO	A	203	-	3,3,3	0.51	0	2,2,2	0.15	0
2	EDO	A	204	-	3,3,3	0.49	0	2,2,2	0.26	0
2	EDO	B	204	-	3,3,3	0.43	0	2,2,2	0.28	0
2	EDO	A	201	-	3,3,3	0.46	0	2,2,2	0.35	0
2	EDO	B	203	-	3,3,3	0.44	0	2,2,2	0.38	0
2	EDO	C	201	-	3,3,3	0.50	0	2,2,2	0.82	0
2	EDO	B	202	-	3,3,3	0.42	0	2,2,2	0.17	0
2	EDO	D	202	-	3,3,3	0.50	0	2,2,2	0.98	0
2	EDO	A	202	-	3,3,3	0.40	0	2,2,2	0.47	0
2	EDO	B	205	-	3,3,3	0.42	0	2,2,2	0.46	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	D	201	-	-	0/1/1/1	-
2	EDO	B	201	-	-	1/1/1/1	-
2	EDO	A	203	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	EDO	A	204	-	-	1/1/1/1	-
2	EDO	B	204	-	-	1/1/1/1	-
2	EDO	A	201	-	-	0/1/1/1	-
2	EDO	B	203	-	-	1/1/1/1	-
2	EDO	C	201	-	-	0/1/1/1	-
2	EDO	B	202	-	-	1/1/1/1	-
2	EDO	D	202	-	-	1/1/1/1	-
2	EDO	A	202	-	-	1/1/1/1	-
2	EDO	B	205	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	204	EDO	O1-C1-C2-O2
2	B	203	EDO	O1-C1-C2-O2
2	D	202	EDO	O1-C1-C2-O2
2	A	203	EDO	O1-C1-C2-O2
2	A	204	EDO	O1-C1-C2-O2
2	A	202	EDO	O1-C1-C2-O2
2	B	201	EDO	O1-C1-C2-O2
2	B	202	EDO	O1-C1-C2-O2

There are no ring outliers.

4 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	204	EDO	3	0
2	B	203	EDO	2	0
2	D	202	EDO	1	0
2	B	205	EDO	1	0

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 [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	198/198 (100%)	-0.16	1 (0%) 91 90	25, 41, 62, 68	0
1	B	198/198 (100%)	-0.14	0 100 100	25, 40, 61, 67	0
1	C	198/198 (100%)	0.07	2 (1%) 82 81	28, 49, 76, 85	0
1	D	197/198 (99%)	0.02	3 (1%) 73 72	27, 50, 76, 81	0
All	All	791/792 (99%)	-0.05	6 (0%) 86 85	25, 45, 71, 85	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	88	ILE	3.3
1	D	88	ILE	3.2
1	D	197	LEU	3.2
1	D	136	TYR	2.2
1	C	92	THR	2.0
1	A	197	LEU	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 monosaccharides 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
2	EDO	A	204	4/4	0.75	0.12	69,71,72,73	0
2	EDO	B	204	4/4	0.84	0.17	60,61,62,63	0
3	CL	B	206	1/1	0.89	0.13	54,54,54,54	0
2	EDO	B	205	4/4	0.90	0.14	50,54,57,58	0
2	EDO	A	203	4/4	0.91	0.16	53,60,63,64	0
2	EDO	A	201	4/4	0.92	0.16	40,41,43,52	0
2	EDO	D	202	4/4	0.93	0.12	47,47,48,54	0
2	EDO	A	202	4/4	0.93	0.16	33,42,42,49	0
2	EDO	B	203	4/4	0.94	0.14	44,53,56,62	0
2	EDO	B	201	4/4	0.95	0.14	39,40,43,50	0
3	CL	C	203	1/1	0.96	0.16	47,47,47,47	0
2	EDO	C	201	4/4	0.97	0.12	32,32,34,41	0
2	EDO	B	202	4/4	0.97	0.29	35,43,44,49	0
2	EDO	D	201	4/4	0.97	0.12	30,33,34,42	0
3	CL	C	202	1/1	0.98	0.12	45,45,45,45	0
3	CL	A	205	1/1	0.98	0.09	52,52,52,52	0

6.5 Other polymers ⓘ

There are no such residues in this entry.