



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:01 AM UTC

PDB ID : 3OV2 / pdb_00003ov2
Title : Curcumin synthase 1 from Curcuma longa
Authors : Katsuyama, Y.; Miyazono, K.; Tanokura, M.; Ohnishi, Y.; Horinouchi, S.
Deposited on : 2010-09-15
Resolution : 2.32 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

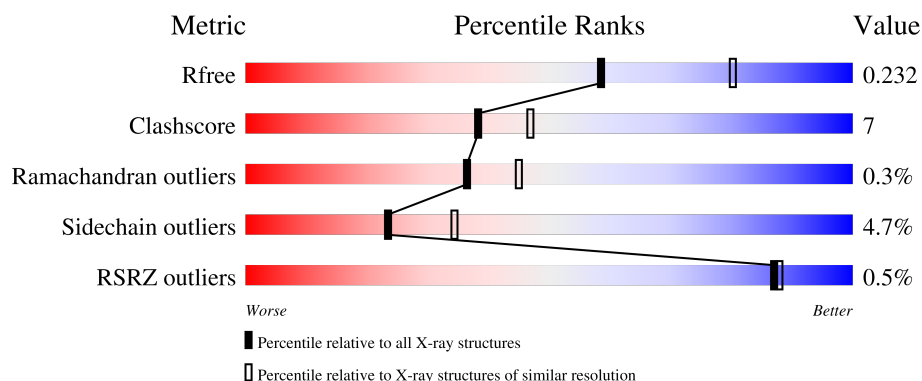
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.32 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	7754 (2.34-2.30)
Clashscore	190562	8383 (2.34-2.30)
Ramachandran outliers	187476	8303 (2.34-2.30)
Sidechain outliers	187428	8303 (2.34-2.30)
RSRZ outliers	180081	7760 (2.34-2.30)

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	393	% 82% 15% ..
1	B	393	86% 12% ..
1	C	393	% 84% 12% .
1	D	393	86% 11% ..

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	MLI	A	396	-	-	X	-
2	MLI	A	398	-	-	X	-
2	MLI	B	397	-	-	X	-
2	MLI	D	394	-	-	X	-
3	EDO	B	412	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 13043 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 Curcumin synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	389	Total	C	N	O	S	0	0	0
			3032	1924	535	557	16			
1	B	389	Total	C	N	O	S	0	0	0
			3032	1924	535	557	16			
1	C	392	Total	C	N	O	S	0	0	0
			3057	1939	542	560	16			
1	D	389	Total	C	N	O	S	0	0	0
			3032	1924	535	557	16			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	390	HIS	-	expression tag	UNP C0SVZ6
A	391	HIS	-	expression tag	UNP C0SVZ6
A	392	HIS	-	expression tag	UNP C0SVZ6
A	393	HIS	-	expression tag	UNP C0SVZ6
B	390	HIS	-	expression tag	UNP C0SVZ6
B	391	HIS	-	expression tag	UNP C0SVZ6
B	392	HIS	-	expression tag	UNP C0SVZ6
B	393	HIS	-	expression tag	UNP C0SVZ6
C	390	HIS	-	expression tag	UNP C0SVZ6
C	391	HIS	-	expression tag	UNP C0SVZ6
C	392	HIS	-	expression tag	UNP C0SVZ6
C	393	HIS	-	expression tag	UNP C0SVZ6
D	390	HIS	-	expression tag	UNP C0SVZ6
D	391	HIS	-	expression tag	UNP C0SVZ6
D	392	HIS	-	expression tag	UNP C0SVZ6
D	393	HIS	-	expression tag	UNP C0SVZ6

- Molecule 2 is MALONATE ION (CCD ID: MLI) (formula: C₃H₂O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			7	3	4		
2	A	1	Total	C	O	0	0
			7	3	4		
2	A	1	Total	C	O	0	0
			7	3	4		
2	A	1	Total	C	O	0	0
			7	3	4		
2	B	1	Total	C	O	0	0
			7	3	4		
2	B	1	Total	C	O	0	0
			7	3	4		
2	B	1	Total	C	O	0	0
			7	3	4		
2	C	1	Total	C	O	0	0
			7	3	4		
2	D	1	Total	C	O	0	0
			7	3	4		
2	D	1	Total	C	O	0	0
			7	3	4		
2	D	1	Total	C	O	0	0
			7	3	4		
2	D	1	Total	C	O	0	0
			7	3	4		

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



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

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

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

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

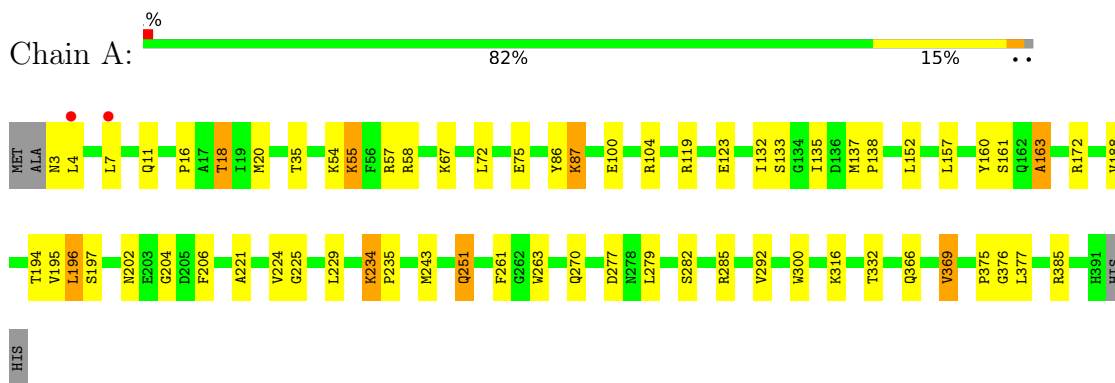
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	141	Total O 141 141	0	0
4	B	124	Total O 124 124	0	0
4	C	149	Total O 149 149	0	0
4	D	126	Total O 126 126	0	0

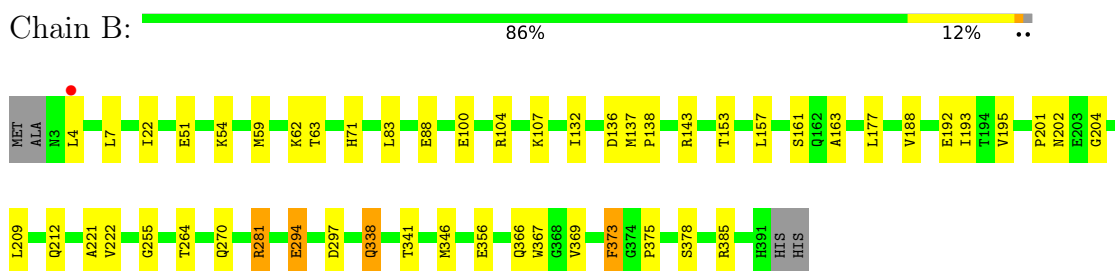
3 Residue-property plots [i](#)

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.

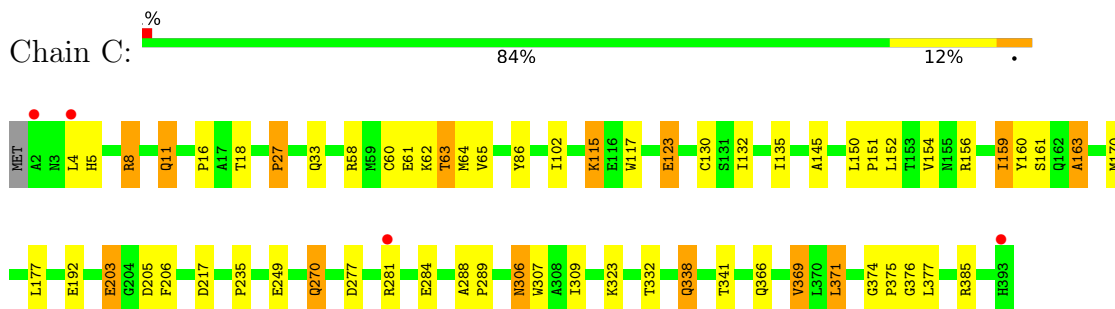
- Molecule 1: Curcumin synthase



- Molecule 1: Curcumin synthase

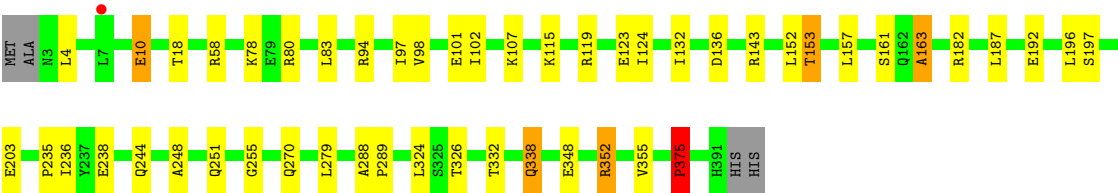


- Molecule 1: Curcumin synthase



- Molecule 1: Curcumin synthase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	77.18Å 115.81Å 221.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.82 – 2.32 19.82 – 2.32	Depositor EDS
% Data completeness (in resolution range)	100.0 (19.82-2.32) 99.8 (19.82-2.32)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.65 (at 2.33Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.183 , 0.235 0.183 , 0.232	Depositor DCC
R_{free} test set	4333 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	27.6	Xtriage
Anisotropy	0.084	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 42.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	13043	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.52% 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, MLI

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	0.69	0/3096	0.87	1/4189 (0.0%)
1	B	0.66	0/3096	0.86	0/4189
1	C	0.70	2/3123 (0.1%)	0.89	2/4226 (0.0%)
1	D	0.67	1/3096 (0.0%)	0.87	1/4189 (0.0%)
All	All	0.68	3/12411 (0.0%)	0.87	4/16793 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	102	ILE	CA-CB	7.66	1.57	1.54
1	C	102	ILE	CA-CB	7.18	1.57	1.54
1	C	27	PRO	CA-C	6.63	1.55	1.51

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	87	LYS	N-CA-C	6.61	120.57	111.71
1	C	374	GLY	CA-C-N	6.34	126.36	119.90
1	C	374	GLY	C-N-CA	6.34	126.36	119.90
1	D	236	ILE	N-CA-C	-5.58	107.01	111.81

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	3032	0	3033	55	0
1	B	3032	0	3033	35	0
1	C	3057	0	3052	54	0
1	D	3032	0	3033	38	0
2	A	35	0	10	8	0
2	B	28	0	8	3	0
2	C	7	0	2	0	0
2	D	28	0	8	5	0
3	A	76	0	114	6	0
3	B	68	0	102	9	0
3	C	52	0	78	8	0
3	D	56	0	84	3	0
4	A	141	0	0	2	0
4	B	124	0	0	2	0
4	C	149	0	0	1	0
4	D	126	0	0	2	0
All	All	13043	0	12557	172	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 7.

All (172) 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:234:LYS:O	1:A:234:LYS:HD2	1.36	1.24
1:A:194:THR:HA	2:A:398:MLI:H11	1.49	0.93
1:C:63:THR:HG23	1:C:65:VAL:H	1.38	0.87
1:A:234:LYS:O	1:A:234:LYS:CD	2.25	0.85
1:A:251:GLN:H	1:A:251:GLN:HE21	1.19	0.85
1:A:172:ARG:HH12	3:B:412:EDO:H22	1.45	0.81
1:C:156:ARG:H	1:D:244:GLN:HE22	1.29	0.79
1:C:156:ARG:H	1:D:244:GLN:NE2	1.79	0.79
1:A:4:LEU:HD22	1:B:385:ARG:HH12	1.49	0.77
1:A:172:ARG:NH1	3:B:412:EDO:H22	1.98	0.76
1:B:22:ILE:CD1	1:B:346:MET:HE3	2.16	0.76
1:D:115:LYS:HD2	1:D:115:LYS:C	2.11	0.75
1:A:119:ARG:NH1	1:A:123:GLU:OE2	2.23	0.71
1:D:152:LEU:HA	2:D:395:MLI:H11	1.71	0.71
1:A:67:LYS:HE2	4:A:554:HOH:O	1.91	0.70
1:B:22:ILE:HG12	1:B:346:MET:CE	2.22	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:61:GLU:CG	3:C:399:EDO:H22	2.23	0.68
1:C:63:THR:CG2	1:C:65:VAL:H	2.07	0.67
1:D:153:THR:HG22	4:D:517:HOH:O	1.95	0.67
1:A:18:THR:HG22	1:A:20:MET:HG3	1.78	0.63
1:A:86:TYR:CZ	1:A:87:LYS:HD2	2.34	0.63
1:C:306:ASN:C	1:C:306:ASN:HD22	2.06	0.63
1:B:212:GLN:HA	3:B:404:EDO:H11	1.79	0.63
1:C:60:CYS:O	1:C:63:THR:HG22	1.99	0.62
1:C:61:GLU:HG2	3:C:399:EDO:H22	1.81	0.62
1:A:202:ASN:ND2	1:A:204:GLY:H	1.98	0.62
1:D:326:THR:HG23	3:D:399:EDO:H21	1.80	0.62
1:A:157:LEU:HA	2:A:396:MLI:O8	2.00	0.62
1:A:163:ALA:CB	1:A:375:PRO:HD2	2.31	0.61
1:C:33:GLN:HE21	1:C:63:THR:HG21	1.64	0.60
1:D:251:GLN:H	1:D:251:GLN:CD	2.09	0.59
1:B:71:HIS:ND1	2:B:395:MLI:H12	2.18	0.58
1:B:157:LEU:HB2	3:B:412:EDO:H21	1.85	0.58
1:C:61:GLU:HG3	3:C:399:EDO:H22	1.86	0.58
1:D:197:SER:HB2	3:D:408:EDO:H21	1.85	0.58
1:C:270:GLN:HE21	1:C:270:GLN:H	1.52	0.57
1:D:196:LEU:HB2	2:D:394:MLI:C2	2.35	0.56
1:A:196:LEU:HB2	2:A:398:MLI:O9	2.06	0.56
1:B:22:ILE:HG12	1:B:346:MET:HE3	1.87	0.56
1:B:22:ILE:HD11	1:B:346:MET:HE3	1.87	0.56
1:A:194:THR:HA	2:A:398:MLI:C1	2.30	0.56
1:D:119:ARG:NH2	1:D:123:GLU:HG3	2.19	0.56
1:A:87:LYS:HG2	1:A:261:PHE:CD2	2.41	0.56
1:B:163:ALA:HB1	4:B:428:HOH:O	2.06	0.55
1:D:192:GLU:HG3	1:D:338:GLN:HB2	1.89	0.55
1:A:55:LYS:HG2	1:A:206:PHE:CE1	2.41	0.55
1:C:369:VAL:CG2	1:C:371:LEU:HD13	2.37	0.54
1:A:100:GLU:HG3	1:A:104:ARG:NH2	2.23	0.54
1:B:202:ASN:ND2	1:B:204:GLY:H	2.05	0.54
1:C:11:GLN:NE2	1:C:11:GLN:HA	2.23	0.53
1:D:196:LEU:HB2	2:D:394:MLI:O7	2.08	0.53
1:A:57:ARG:HE	3:A:404:EDO:C1	2.21	0.53
1:A:72:LEU:HG	1:A:195:VAL:HG22	1.91	0.53
1:D:348:GLU:OE2	1:D:352:ARG:HD2	2.09	0.52
1:D:203:GLU:H	1:D:203:GLU:CD	2.17	0.52
1:A:119:ARG:CZ	1:A:123:GLU:OE2	2.58	0.52
1:B:22:ILE:CG1	1:B:346:MET:HE3	2.39	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:130:CYS:SG	1:C:170:MET:HE2	2.50	0.52
1:D:338:GLN:NE2	2:D:394:MLI:O8	2.42	0.52
1:A:104:ARG:HG2	4:A:476:HOH:O	2.09	0.52
1:A:163:ALA:HB3	1:A:375:PRO:HD2	1.92	0.51
1:D:18:THR:HB	1:D:235:PRO:HB3	1.92	0.51
1:A:197:SER:O	3:A:407:EDO:H11	2.10	0.51
1:C:163:ALA:CB	1:C:375:PRO:HD2	2.41	0.51
1:C:306:ASN:ND2	1:C:309:ILE:H	2.09	0.51
1:A:16:PRO:HG3	1:B:4:LEU:HD23	1.93	0.51
1:C:4:LEU:HD12	1:D:238:GLU:OE1	2.11	0.50
1:D:163:ALA:HB1	1:D:375:PRO:HD2	1.93	0.50
1:A:376:GLY:N	1:A:377:LEU:HA	2.26	0.50
1:B:222:VAL:HG11	1:B:346:MET:HE1	1.92	0.50
1:A:157:LEU:HD12	2:A:396:MLI:O8	2.12	0.50
1:C:117:TRP:HH2	1:C:123:GLU:HB3	1.76	0.50
1:C:307:TRP:HE1	3:C:397:EDO:H21	1.76	0.50
1:A:234:LYS:HD2	1:A:234:LYS:C	2.28	0.50
1:C:159:ILE:HD13	1:D:157:LEU:HD11	1.94	0.50
1:A:18:THR:HB	1:A:225:GLY:O	2.12	0.50
1:A:152:LEU:HA	2:A:397:MLI:H11	1.94	0.49
1:A:132:ILE:HA	1:A:161:SER:HA	1.95	0.49
1:B:177:LEU:CD1	3:B:412:EDO:H11	2.42	0.49
1:C:366:GLN:HE21	1:C:385:ARG:HD2	1.78	0.49
1:A:366:GLN:HE21	1:A:385:ARG:HD2	1.78	0.48
1:D:115:LYS:HD2	1:D:115:LYS:O	2.12	0.48
1:C:8:ARG:HH11	1:C:8:ARG:HG3	1.79	0.48
1:B:373:PHE:HA	1:B:378:SER:O	2.13	0.48
1:A:57:ARG:HE	3:A:404:EDO:H11	1.77	0.48
1:C:369:VAL:HG22	1:C:371:LEU:HD13	1.95	0.48
1:B:264:THR:OG1	3:B:414:EDO:H11	2.12	0.47
1:B:366:GLN:H	3:B:403:EDO:H21	1.78	0.47
1:D:115:LYS:C	1:D:115:LYS:CD	2.86	0.47
1:B:132:ILE:HA	1:B:161:SER:HA	1.97	0.47
1:B:192:GLU:HG3	1:B:338:GLN:HB2	1.96	0.47
1:C:145:ALA:HA	1:C:150:LEU:HD12	1.96	0.47
1:A:277:ASP:HA	1:A:316:LYS:HE2	1.97	0.47
1:C:18:THR:HB	1:C:235:PRO:HB3	1.96	0.47
1:C:163:ALA:HB3	1:C:375:PRO:HD2	1.97	0.47
1:D:101:GLU:CG	3:D:405:EDO:H21	2.45	0.47
1:A:196:LEU:HB2	2:A:398:MLI:C3	2.45	0.47
1:B:136:ASP:OD2	1:B:143:ARG:NH2	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:202:ASN:HD22	1:B:204:GLY:H	1.62	0.47
1:C:192:GLU:HG3	1:C:338:GLN:HB2	1.96	0.47
1:A:18:THR:HG23	1:A:235:PRO:HB3	1.97	0.47
1:C:8:ARG:HG3	1:C:8:ARG:NH1	2.29	0.46
1:B:201:PRO:HB3	1:B:209:LEU:HD13	1.97	0.46
1:A:18:THR:O	1:A:224:VAL:HA	2.15	0.46
1:B:367:TRP:H	3:B:403:EDO:H21	1.80	0.46
1:B:356:GLU:O	3:B:405:EDO:H11	2.16	0.45
1:C:58:ARG:HG3	1:C:62:LYS:HE2	1.98	0.45
1:A:75:GLU:HG2	1:C:277:ASP:CG	2.41	0.45
1:C:270:GLN:HE21	1:C:270:GLN:N	2.14	0.45
1:D:94:ARG:O	1:D:98:VAL:HG22	2.17	0.45
1:C:288:ALA:HB3	1:C:289:PRO:HD3	1.98	0.45
1:D:255:GLY:HA2	1:D:375:PRO:HG3	1.98	0.45
1:A:54:LYS:HA	3:A:404:EDO:H21	1.98	0.44
1:B:255:GLY:HA2	1:B:375:PRO:HG3	1.98	0.44
1:D:132:ILE:HA	1:D:161:SER:HA	2.00	0.44
1:C:135:ILE:HG13	3:C:405:EDO:H12	2.00	0.44
1:D:163:ALA:HB2	1:D:375:PRO:HG2	2.00	0.44
1:C:306:ASN:C	1:C:306:ASN:ND2	2.76	0.44
1:A:58:ARG:HA	1:A:58:ARG:HD2	1.84	0.43
1:A:35:THR:OG1	3:A:411:EDO:H11	2.18	0.43
1:D:136:ASP:OD2	1:D:143:ARG:NH2	2.51	0.43
1:C:163:ALA:HB2	1:C:375:PRO:HG2	2.00	0.43
1:A:133:SER:HB3	1:A:196:LEU:HD12	2.00	0.43
1:C:135:ILE:HG12	1:C:160:TYR:CG	2.54	0.43
1:C:151:PRO:HG2	1:C:154:VAL:HG23	2.00	0.43
1:C:152:LEU:HD11	1:D:248:ALA:HB2	2.00	0.43
1:A:243:MET:HE2	1:A:243:MET:HB3	1.90	0.43
1:C:11:GLN:HA	1:C:11:GLN:HE21	1.81	0.43
1:C:135:ILE:HG12	1:C:160:TYR:CD2	2.54	0.43
1:C:16:PRO:HG3	1:D:4:LEU:HD23	2.01	0.42
1:D:163:ALA:CB	1:D:375:PRO:HD2	2.50	0.42
1:A:7:LEU:O	1:A:11:GLN:HG2	2.20	0.42
1:A:202:ASN:HD22	1:A:204:GLY:H	1.63	0.42
1:A:243:MET:HB2	2:B:397:MLI:H11	2.00	0.42
1:C:177:LEU:HD11	3:C:403:EDO:H12	2.02	0.42
1:B:137:MET:HA	1:B:138:PRO:C	2.44	0.42
1:B:294:GLU:O	1:B:297:ASP:HB2	2.19	0.42
1:C:86:TYR:OH	1:C:205:ASP:OD1	2.26	0.42
1:A:263:TRP:CH2	1:B:137:MET:HE3	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:323:LYS:HA	1:C:323:LYS:HD3	1.81	0.42
1:D:197:SER:N	2:D:394:MLI:O6	2.53	0.42
1:A:194:THR:CA	2:A:398:MLI:H11	2.36	0.42
1:C:27:PRO:HG2	1:C:217:ASP:HB2	2.01	0.42
1:A:188:VAL:O	1:A:221:ALA:HA	2.19	0.42
1:A:3:ASN:HD22	1:A:4:LEU:N	2.18	0.42
1:B:188:VAL:O	1:B:221:ALA:HA	2.20	0.42
1:C:307:TRP:HE1	3:C:397:EDO:C2	2.33	0.42
1:B:100:GLU:O	1:B:104:ARG:HG3	2.20	0.41
1:B:193:ILE:HG22	1:B:195:VAL:HG23	2.01	0.41
1:C:63:THR:O	1:C:64:MET:HB2	2.20	0.41
1:A:4:LEU:H	1:A:4:LEU:HD23	1.85	0.41
3:A:410:EDO:H21	1:B:157:LEU:HD12	2.02	0.41
1:C:132:ILE:HA	1:C:161:SER:HA	2.03	0.41
1:C:203:GLU:H	1:C:203:GLU:HG2	1.50	0.41
1:A:137:MET:HA	1:A:138:PRO:C	2.45	0.41
1:C:115:LYS:NZ	4:C:492:HOH:O	2.54	0.41
1:D:78:LYS:HD3	1:D:78:LYS:HA	1.67	0.41
1:D:78:LYS:NZ	4:D:503:HOH:O	2.54	0.41
1:A:300:TRP:CZ3	1:A:369:VAL:HG21	2.56	0.41
1:B:59:MET:O	1:B:63:THR:HG23	2.21	0.41
1:C:249:GLU:O	3:C:406:EDO:H12	2.21	0.41
1:D:10:GLU:H	1:D:10:GLU:HG2	1.67	0.41
1:D:288:ALA:N	1:D:289:PRO:HD2	2.36	0.41
1:B:281:ARG:NH1	4:B:470:HOH:O	2.54	0.41
1:B:153:THR:O	2:B:397:MLI:O9	2.39	0.40
1:C:60:CYS:O	1:C:63:THR:CG2	2.68	0.40
1:C:376:GLY:HA3	1:C:377:LEU:C	2.45	0.40
1:D:94:ARG:HD3	1:D:196:LEU:HA	2.02	0.40
1:D:80:ARG:HD2	1:D:97:ILE:HD11	2.03	0.40
1:D:124:ILE:HD12	1:D:187:LEU:HB2	2.03	0.40
1:A:135:ILE:HA	1:A:160:TYR:CZ	2.57	0.40
1:C:117:TRP:CH2	1:C:123:GLU:HB3	2.56	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	387/393 (98%)	374 (97%)	12 (3%)	1 (0%)	36	45
1	B	387/393 (98%)	373 (96%)	14 (4%)	0	100	100
1	C	390/393 (99%)	379 (97%)	10 (3%)	1 (0%)	36	45
1	D	387/393 (98%)	374 (97%)	11 (3%)	2 (0%)	24	30
All	All	1551/1572 (99%)	1500 (97%)	47 (3%)	4 (0%)	36	45

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	163	ALA
1	D	163	ALA
1	A	163	ALA
1	D	375	PRO

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	315/318 (99%)	302 (96%)	13 (4%)	27	40
1	B	315/318 (99%)	301 (96%)	14 (4%)	25	37
1	C	317/318 (100%)	299 (94%)	18 (6%)	18	26
1	D	315/318 (99%)	301 (96%)	14 (4%)	25	37
All	All	1262/1272 (99%)	1203 (95%)	59 (5%)	23	34

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	THR
1	A	55	LYS
1	A	196	LEU
1	A	229	LEU
1	A	234	LYS
1	A	251	GLN
1	A	270	GLN
1	A	279	LEU
1	A	282	SER
1	A	285	ARG
1	A	292	VAL
1	A	332	THR
1	A	369	VAL
1	B	7	LEU
1	B	51	GLU
1	B	54	LYS
1	B	62	LYS
1	B	83	LEU
1	B	88	GLU
1	B	107	LYS
1	B	270	GLN
1	B	281	ARG
1	B	294	GLU
1	B	338	GLN
1	B	341	THR
1	B	369	VAL
1	B	373	PHE
1	C	5	HIS
1	C	8	ARG
1	C	11	GLN
1	C	63	THR
1	C	115	LYS
1	C	123	GLU
1	C	159	ILE
1	C	203	GLU
1	C	206	PHE
1	C	270	GLN
1	C	281	ARG
1	C	284	GLU
1	C	306	ASN
1	C	332	THR
1	C	338	GLN

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Mol	Chain	Res	Type
1	C	341	THR
1	C	369	VAL
1	C	371	LEU
1	D	10	GLU
1	D	58	ARG
1	D	83	LEU
1	D	107	LYS
1	D	153	THR
1	D	182	ARG
1	D	270	GLN
1	D	279	LEU
1	D	324	LEU
1	D	332	THR
1	D	338	GLN
1	D	352	ARG
1	D	355	VAL
1	D	375	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	3	ASN
1	A	202	ASN
1	A	251	GLN
1	A	366	GLN
1	A	390	HIS
1	B	3	ASN
1	B	5	HIS
1	B	202	ASN
1	B	391	HIS
1	C	11	GLN
1	C	45	ASN
1	C	202	ASN
1	C	270	GLN
1	C	303	HIS
1	C	306	ASN
1	C	329	HIS
1	C	366	GLN
1	D	5	HIS
1	D	202	ASN
1	D	244	GLN
1	D	303	HIS

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Mol	Chain	Res	Type
1	D	338	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

77 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	$\# Z > 2$
3	EDO	A	415	-	3,3,3	0.37	0	2,2,2	0.56	0
2	MLI	A	397	-	6,6,6	1.17	0	7,7,7	1.46	1 (14%)
3	EDO	B	406	-	3,3,3	0.38	0	2,2,2	0.55	0
3	EDO	D	410	-	3,3,3	0.58	0	2,2,2	0.16	0
2	MLI	B	395	-	6,6,6	1.13	0	7,7,7	1.44	2 (28%)
3	EDO	D	406	-	3,3,3	0.42	0	2,2,2	0.55	0
3	EDO	D	398	-	3,3,3	0.43	0	2,2,2	0.53	0
3	EDO	A	412	-	3,3,3	0.44	0	2,2,2	0.42	0
3	EDO	A	407	-	3,3,3	0.41	0	2,2,2	0.27	0
2	MLI	A	396	-	6,6,6	1.27	0	7,7,7	1.69	2 (28%)
2	MLI	A	398	-	6,6,6	1.62	1 (16%)	7,7,7	2.02	3 (42%)
3	EDO	C	396	-	3,3,3	0.43	0	2,2,2	0.48	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	EDO	C	404	-	3,3,3	0.60	0	2,2,2	0.18	0
3	EDO	C	402	-	3,3,3	0.58	0	2,2,2	0.13	0
3	EDO	B	404	-	3,3,3	0.50	0	2,2,2	0.47	0
3	EDO	C	398	-	3,3,3	0.48	0	2,2,2	0.34	0
3	EDO	B	413	-	3,3,3	0.48	0	2,2,2	0.18	0
3	EDO	C	400	-	3,3,3	0.43	0	2,2,2	0.44	0
2	MLI	C	394	-	6,6,6	1.15	0	7,7,7	1.04	0
2	MLI	D	396	-	6,6,6	1.18	0	7,7,7	1.36	1 (14%)
3	EDO	C	397	-	3,3,3	0.57	0	2,2,2	0.06	0
3	EDO	C	403	-	3,3,3	0.50	0	2,2,2	0.05	0
3	EDO	B	412	-	3,3,3	0.42	0	2,2,2	0.20	0
3	EDO	B	403	-	3,3,3	0.50	0	2,2,2	0.43	0
3	EDO	B	402	-	3,3,3	0.43	0	2,2,2	0.38	0
3	EDO	B	405	-	3,3,3	0.51	0	2,2,2	0.25	0
3	EDO	B	410	-	3,3,3	0.44	0	2,2,2	0.37	0
3	EDO	D	399	-	3,3,3	0.44	0	2,2,2	0.36	0
3	EDO	D	404	-	3,3,3	0.55	0	2,2,2	0.09	0
2	MLI	D	394	-	6,6,6	1.13	0	7,7,7	1.82	2 (28%)
3	EDO	C	399	-	3,3,3	0.42	0	2,2,2	0.34	0
2	MLI	D	395	-	6,6,6	1.16	0	7,7,7	1.53	1 (14%)
3	EDO	A	403	-	3,3,3	0.45	0	2,2,2	0.43	0
3	EDO	A	410	-	3,3,3	0.55	0	2,2,2	0.24	0
3	EDO	A	406	-	3,3,3	0.45	0	2,2,2	0.19	0
3	EDO	D	400	-	3,3,3	0.44	0	2,2,2	0.45	0
3	EDO	C	395	-	3,3,3	0.46	0	2,2,2	0.17	0
3	EDO	B	409	-	3,3,3	0.53	0	2,2,2	0.06	0
3	EDO	B	400	-	3,3,3	0.59	0	2,2,2	0.08	0
3	EDO	A	413	-	3,3,3	0.48	0	2,2,2	0.51	0
3	EDO	D	407	-	3,3,3	0.36	0	2,2,2	0.67	0
2	MLI	B	396	-	6,6,6	1.31	0	7,7,7	1.12	0
3	EDO	B	399	-	3,3,3	0.50	0	2,2,2	0.49	0
3	EDO	A	404	-	3,3,3	0.31	0	2,2,2	0.58	0
3	EDO	D	403	-	3,3,3	0.45	0	2,2,2	0.33	0
3	EDO	A	409	-	3,3,3	0.49	0	2,2,2	0.30	0
3	EDO	D	408	-	3,3,3	0.41	0	2,2,2	0.73	0
3	EDO	A	405	-	3,3,3	0.36	0	2,2,2	0.54	0
3	EDO	A	408	-	3,3,3	0.57	0	2,2,2	0.16	0
3	EDO	A	399	-	3,3,3	0.36	0	2,2,2	0.60	0
3	EDO	C	401	-	3,3,3	0.52	0	2,2,2	0.23	0
3	EDO	B	408	-	3,3,3	0.49	0	2,2,2	0.19	0
3	EDO	B	414	-	3,3,3	0.39	0	2,2,2	0.72	0
3	EDO	B	398	-	3,3,3	0.46	0	2,2,2	0.32	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	EDO	A	402	-	3,3,3	0.47	0	2,2,2	0.35	0
3	EDO	D	411	-	3,3,3	0.53	0	2,2,2	0.33	0
3	EDO	D	401	-	3,3,3	0.42	0	2,2,2	0.41	0
3	EDO	A	411	-	3,3,3	0.48	0	2,2,2	0.22	0
3	EDO	A	401	-	3,3,3	0.71	0	2,2,2	0.33	0
2	MLI	B	397	-	6,6,6	1.21	0	7,7,7	1.13	1 (14%)
2	MLI	B	394	-	6,6,6	1.12	0	7,7,7	1.39	0
3	EDO	B	407	-	3,3,3	0.43	0	2,2,2	0.33	0
3	EDO	B	401	-	3,3,3	0.41	0	2,2,2	0.61	0
3	EDO	C	406	-	3,3,3	0.56	0	2,2,2	0.14	0
2	MLI	A	395	-	6,6,6	1.02	0	7,7,7	1.41	2 (28%)
3	EDO	D	402	-	3,3,3	0.45	0	2,2,2	0.31	0
3	EDO	A	400	-	3,3,3	0.46	0	2,2,2	0.39	0
3	EDO	A	416	-	3,3,3	0.41	0	2,2,2	0.48	0
3	EDO	B	411	-	3,3,3	0.53	0	2,2,2	0.27	0
2	MLI	D	397	-	6,6,6	1.21	0	7,7,7	1.11	0
3	EDO	A	414	-	3,3,3	0.49	0	2,2,2	0.22	0
3	EDO	D	409	-	3,3,3	0.39	0	2,2,2	0.52	0
3	EDO	C	407	-	3,3,3	0.43	0	2,2,2	0.42	0
3	EDO	D	405	-	3,3,3	0.49	0	2,2,2	0.39	0
3	EDO	A	417	-	3,3,3	0.39	0	2,2,2	0.53	0
2	MLI	A	394	-	6,6,6	1.10	0	7,7,7	1.18	0
3	EDO	C	405	-	3,3,3	0.58	0	2,2,2	0.24	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	415	-	-	1/1/1/1	-
2	MLI	A	397	-	-	3/4/4/4	-
3	EDO	B	406	-	-	1/1/1/1	-
3	EDO	D	410	-	-	1/1/1/1	-
2	MLI	B	395	-	-	0/4/4/4	-
3	EDO	D	406	-	-	0/1/1/1	-
3	EDO	D	398	-	-	0/1/1/1	-
3	EDO	A	412	-	-	1/1/1/1	-
3	EDO	A	407	-	-	1/1/1/1	-
2	MLI	A	396	-	-	0/4/4/4	-
2	MLI	A	398	-	-	2/4/4/4	-
3	EDO	C	396	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EDO	C	404	-	-	0/1/1/1	-
3	EDO	C	402	-	-	1/1/1/1	-
3	EDO	B	404	-	-	1/1/1/1	-
3	EDO	C	398	-	-	1/1/1/1	-
3	EDO	B	413	-	-	1/1/1/1	-
3	EDO	C	400	-	-	1/1/1/1	-
2	MLI	C	394	-	-	0/4/4/4	-
2	MLI	D	396	-	-	2/4/4/4	-
3	EDO	C	397	-	-	1/1/1/1	-
3	EDO	C	403	-	-	1/1/1/1	-
3	EDO	B	412	-	-	0/1/1/1	-
3	EDO	B	403	-	-	1/1/1/1	-
3	EDO	B	402	-	-	1/1/1/1	-
3	EDO	B	405	-	-	1/1/1/1	-
3	EDO	B	410	-	-	0/1/1/1	-
3	EDO	D	399	-	-	0/1/1/1	-
3	EDO	D	404	-	-	0/1/1/1	-
2	MLI	D	394	-	-	2/4/4/4	-
3	EDO	C	399	-	-	1/1/1/1	-
2	MLI	D	395	-	-	2/4/4/4	-
3	EDO	A	403	-	-	1/1/1/1	-
3	EDO	A	410	-	-	1/1/1/1	-
3	EDO	A	406	-	-	1/1/1/1	-
3	EDO	D	400	-	-	0/1/1/1	-
3	EDO	C	395	-	-	0/1/1/1	-
3	EDO	B	409	-	-	1/1/1/1	-
3	EDO	B	400	-	-	0/1/1/1	-
3	EDO	A	413	-	-	0/1/1/1	-
3	EDO	D	407	-	-	1/1/1/1	-
2	MLI	B	396	-	-	4/4/4/4	-
3	EDO	B	399	-	-	1/1/1/1	-
3	EDO	A	404	-	-	1/1/1/1	-
3	EDO	D	403	-	-	0/1/1/1	-
3	EDO	A	409	-	-	1/1/1/1	-
3	EDO	D	408	-	-	1/1/1/1	-
3	EDO	A	405	-	-	1/1/1/1	-
3	EDO	A	408	-	-	1/1/1/1	-
3	EDO	A	399	-	-	1/1/1/1	-
3	EDO	C	401	-	-	1/1/1/1	-
3	EDO	B	408	-	-	1/1/1/1	-
3	EDO	B	414	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EDO	B	398	-	-	0/1/1/1	-
3	EDO	A	402	-	-	1/1/1/1	-
3	EDO	D	411	-	-	0/1/1/1	-
3	EDO	D	401	-	-	1/1/1/1	-
3	EDO	A	411	-	-	0/1/1/1	-
3	EDO	A	401	-	-	0/1/1/1	-
2	MLI	B	397	-	-	0/4/4/4	-
2	MLI	B	394	-	-	2/4/4/4	-
3	EDO	B	407	-	-	1/1/1/1	-
3	EDO	B	401	-	-	1/1/1/1	-
3	EDO	C	406	-	-	1/1/1/1	-
2	MLI	A	395	-	-	4/4/4/4	-
3	EDO	D	402	-	-	1/1/1/1	-
3	EDO	A	400	-	-	1/1/1/1	-
3	EDO	A	416	-	-	1/1/1/1	-
3	EDO	B	411	-	-	0/1/1/1	-
2	MLI	D	397	-	-	2/4/4/4	-
3	EDO	A	414	-	-	0/1/1/1	-
3	EDO	D	409	-	-	0/1/1/1	-
3	EDO	C	407	-	-	1/1/1/1	-
3	EDO	D	405	-	-	1/1/1/1	-
3	EDO	A	417	-	-	0/1/1/1	-
2	MLI	A	394	-	-	0/4/4/4	-
3	EDO	C	405	-	-	0/1/1/1	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	398	MLI	O9-C3	-2.10	1.23	1.30

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	398	MLI	C3-C1-C2	-3.74	99.73	112.95
2	D	394	MLI	C3-C1-C2	-3.32	101.23	112.95
2	A	396	MLI	C3-C1-C2	-2.71	103.39	112.95
2	B	395	MLI	O7-C2-C1	2.54	122.39	114.51
2	D	395	MLI	O9-C3-C1	2.37	121.86	114.51
2	A	398	MLI	O6-C2-C1	-2.27	115.64	122.11
2	D	394	MLI	O6-C2-C1	-2.17	115.93	122.11
2	A	395	MLI	C3-C1-C2	-2.15	105.37	112.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	398	MLI	O8-C3-C1	-2.13	116.06	122.11
2	B	395	MLI	O6-C2-C1	-2.12	116.08	122.11
2	A	397	MLI	O9-C3-C1	2.11	121.06	114.51
2	D	396	MLI	O7-C2-C1	2.05	120.85	114.51
2	A	396	MLI	O6-C2-C1	-2.04	116.30	122.11
2	A	395	MLI	O8-C3-C1	-2.04	116.31	122.11
2	B	397	MLI	O7-C2-C1	2.01	120.73	114.51

There are no chirality outliers.

All (64) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	403	EDO	O1-C1-C2-O2
3	A	404	EDO	O1-C1-C2-O2
3	A	405	EDO	O1-C1-C2-O2
3	A	409	EDO	O1-C1-C2-O2
3	A	412	EDO	O1-C1-C2-O2
3	B	399	EDO	O1-C1-C2-O2
3	B	407	EDO	O1-C1-C2-O2
3	C	399	EDO	O1-C1-C2-O2
3	D	407	EDO	O1-C1-C2-O2
3	D	410	EDO	O1-C1-C2-O2
2	D	394	MLI	C3-C1-C2-O7
2	D	397	MLI	C3-C1-C2-O7
3	A	415	EDO	O1-C1-C2-O2
3	B	401	EDO	O1-C1-C2-O2
3	B	402	EDO	O1-C1-C2-O2
3	B	414	EDO	O1-C1-C2-O2
3	C	403	EDO	O1-C1-C2-O2
2	A	398	MLI	C2-C1-C3-O9
2	B	396	MLI	C2-C1-C3-O8
2	D	397	MLI	C3-C1-C2-O6
3	A	400	EDO	O1-C1-C2-O2
3	B	403	EDO	O1-C1-C2-O2
3	D	405	EDO	O1-C1-C2-O2
2	A	395	MLI	C2-C1-C3-O9
2	A	398	MLI	C2-C1-C3-O8
2	D	394	MLI	C3-C1-C2-O6
3	B	404	EDO	O1-C1-C2-O2
3	B	406	EDO	O1-C1-C2-O2
3	D	401	EDO	O1-C1-C2-O2
2	A	395	MLI	C3-C1-C2-O7

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Mol	Chain	Res	Type	Atoms
2	B	396	MLI	C2-C1-C3-O9
2	A	395	MLI	C3-C1-C2-O6
2	D	396	MLI	C2-C1-C3-O9
3	C	406	EDO	O1-C1-C2-O2
2	B	394	MLI	C3-C1-C2-O6
2	D	396	MLI	C2-C1-C3-O8
3	A	406	EDO	O1-C1-C2-O2
3	B	413	EDO	O1-C1-C2-O2
3	C	400	EDO	O1-C1-C2-O2
2	A	395	MLI	C2-C1-C3-O8
3	A	408	EDO	O1-C1-C2-O2
3	A	410	EDO	O1-C1-C2-O2
3	B	408	EDO	O1-C1-C2-O2
3	C	401	EDO	O1-C1-C2-O2
2	B	394	MLI	C3-C1-C2-O7
2	A	397	MLI	C2-C1-C3-O8
2	D	395	MLI	C2-C1-C3-O8
3	A	399	EDO	O1-C1-C2-O2
3	A	416	EDO	O1-C1-C2-O2
3	B	405	EDO	O1-C1-C2-O2
3	C	398	EDO	O1-C1-C2-O2
3	C	407	EDO	O1-C1-C2-O2
3	D	402	EDO	O1-C1-C2-O2
2	A	397	MLI	C2-C1-C3-O9
3	A	402	EDO	O1-C1-C2-O2
3	B	409	EDO	O1-C1-C2-O2
3	D	408	EDO	O1-C1-C2-O2
2	D	395	MLI	C2-C1-C3-O9
3	A	407	EDO	O1-C1-C2-O2
3	C	397	EDO	O1-C1-C2-O2
3	C	402	EDO	O1-C1-C2-O2
2	A	397	MLI	C3-C1-C2-O6
2	B	396	MLI	C3-C1-C2-O6
2	B	396	MLI	C3-C1-C2-O7

There are no ring outliers.

24 monomers are involved in 42 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	397	MLI	1	0
2	B	395	MLI	1	0
3	A	407	EDO	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	396	MLI	2	0
2	A	398	MLI	5	0
3	B	404	EDO	1	0
3	C	397	EDO	2	0
3	C	403	EDO	1	0
3	B	412	EDO	4	0
3	B	403	EDO	2	0
3	B	405	EDO	1	0
3	D	399	EDO	1	0
2	D	394	MLI	4	0
3	C	399	EDO	3	0
2	D	395	MLI	1	0
3	A	410	EDO	1	0
3	A	404	EDO	3	0
3	D	408	EDO	1	0
3	B	414	EDO	1	0
3	A	411	EDO	1	0
2	B	397	MLI	2	0
3	C	406	EDO	1	0
3	D	405	EDO	1	0
3	C	405	EDO	1	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 [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	389/393 (98%)	-0.56	2 (0%) 87 88	14, 22, 41, 64	0
1	B	389/393 (98%)	-0.37	1 (0%) 90 90	14, 27, 50, 63	0
1	C	392/393 (99%)	-0.43	4 (1%) 79 81	16, 25, 50, 66	0
1	D	389/393 (98%)	-0.42	1 (0%) 90 90	15, 26, 46, 61	0
All	All	1559/1572 (99%)	-0.45	8 (0%) 87 88	14, 25, 48, 66	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	4	LEU	4.0
1	C	4	LEU	2.6
1	C	2	ALA	2.5
1	C	393	HIS	2.5
1	B	4	LEU	2.5
1	A	7	LEU	2.3
1	C	281	ARG	2.1
1	D	7	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 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(Å ²)	Q<0.9
2	MLI	B	396	7/7	0.67	0.18	57,59,60,61	0
2	MLI	D	397	7/7	0.70	0.19	99,99,99,99	0
2	MLI	D	396	7/7	0.71	0.18	56,58,60,61	0
3	EDO	C	406	4/4	0.75	0.19	49,49,49,50	0
3	EDO	A	412	4/4	0.76	0.15	56,56,57,57	0
2	MLI	A	397	7/7	0.77	0.18	58,58,58,59	0
3	EDO	B	406	4/4	0.78	0.18	56,57,57,58	0
3	EDO	D	399	4/4	0.78	0.22	61,62,62,62	0
3	EDO	C	404	4/4	0.80	0.15	45,45,47,47	0
2	MLI	B	395	7/7	0.80	0.17	59,59,60,60	0
3	EDO	B	410	4/4	0.80	0.15	48,50,51,51	0
3	EDO	C	396	4/4	0.81	0.22	51,51,53,55	0
2	MLI	D	395	7/7	0.81	0.16	43,44,45,45	0
3	EDO	D	405	4/4	0.81	0.15	49,49,49,50	0
3	EDO	D	407	4/4	0.82	0.14	39,40,40,40	0
3	EDO	A	413	4/4	0.83	0.14	31,33,34,34	0
3	EDO	B	413	4/4	0.83	0.13	41,41,42,43	0
3	EDO	A	404	4/4	0.83	0.16	47,48,49,49	0
3	EDO	C	398	4/4	0.83	0.16	54,55,55,55	0
3	EDO	C	402	4/4	0.83	0.15	40,41,41,41	0
2	MLI	C	394	7/7	0.84	0.15	58,59,62,63	0
2	MLI	A	394	7/7	0.84	0.14	54,55,55,56	0
3	EDO	C	400	4/4	0.84	0.15	54,56,58,60	0
2	MLI	A	395	7/7	0.85	0.17	41,45,48,48	0
2	MLI	A	396	7/7	0.85	0.19	61,61,62,64	0
3	EDO	A	415	4/4	0.85	0.14	36,38,40,41	0
3	EDO	A	416	4/4	0.86	0.12	55,55,55,55	0
3	EDO	B	403	4/4	0.86	0.10	32,37,37,42	0
3	EDO	B	399	4/4	0.87	0.12	33,33,34,35	0
2	MLI	B	394	7/7	0.87	0.14	64,65,65,66	0
3	EDO	B	405	4/4	0.87	0.14	50,52,52,53	0
3	EDO	D	401	4/4	0.87	0.12	61,61,62,62	0
3	EDO	A	408	4/4	0.87	0.14	50,51,51,51	0
2	MLI	B	397	7/7	0.87	0.13	33,39,43,45	0
3	EDO	D	408	4/4	0.87	0.14	39,40,41,41	0
3	EDO	D	411	4/4	0.87	0.12	29,32,34,34	0
3	EDO	A	403	4/4	0.88	0.14	53,53,54,54	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	EDO	A	417	4/4	0.88	0.11	43,44,44,45	0
3	EDO	C	401	4/4	0.88	0.11	34,36,38,39	0
2	MLI	A	398	7/7	0.88	0.16	33,35,39,42	0
3	EDO	D	406	4/4	0.89	0.10	50,51,52,53	0
2	MLI	D	394	7/7	0.89	0.15	41,47,49,50	0
3	EDO	D	402	4/4	0.90	0.11	52,52,52,53	0
3	EDO	C	397	4/4	0.90	0.11	38,39,40,40	0
3	EDO	C	407	4/4	0.90	0.11	48,49,50,50	0
3	EDO	B	404	4/4	0.90	0.10	37,39,40,40	0
3	EDO	D	400	4/4	0.90	0.13	52,53,54,54	0
3	EDO	D	409	4/4	0.90	0.10	50,52,53,53	0
3	EDO	A	409	4/4	0.90	0.12	39,39,40,41	0
3	EDO	A	400	4/4	0.91	0.08	40,42,43,43	0
3	EDO	A	402	4/4	0.91	0.10	35,35,37,37	0
3	EDO	A	405	4/4	0.91	0.09	36,36,36,36	0
3	EDO	B	402	4/4	0.92	0.08	46,46,47,48	0
3	EDO	B	414	4/4	0.92	0.10	34,34,35,35	0
3	EDO	A	401	4/4	0.92	0.09	21,22,23,25	0
3	EDO	D	403	4/4	0.93	0.09	50,50,51,52	0
3	EDO	B	398	4/4	0.93	0.08	20,22,24,28	0
3	EDO	D	398	4/4	0.93	0.08	34,39,40,42	0
3	EDO	B	407	4/4	0.93	0.10	49,50,51,52	0
3	EDO	A	414	4/4	0.93	0.10	45,47,49,50	0
3	EDO	A	410	4/4	0.93	0.09	33,34,35,35	0
3	EDO	C	399	4/4	0.93	0.17	39,40,40,42	0
3	EDO	A	411	4/4	0.94	0.08	23,25,27,29	0
3	EDO	B	401	4/4	0.94	0.08	29,32,32,35	0
3	EDO	A	399	4/4	0.94	0.07	29,31,32,36	0
3	EDO	D	410	4/4	0.94	0.08	27,28,28,29	0
3	EDO	B	412	4/4	0.94	0.09	25,26,27,28	0
3	EDO	C	403	4/4	0.95	0.12	30,32,32,33	0
3	EDO	B	400	4/4	0.96	0.07	26,29,31,32	0
3	EDO	A	407	4/4	0.96	0.06	33,35,36,37	0
3	EDO	B	408	4/4	0.96	0.07	17,17,20,20	0
3	EDO	B	409	4/4	0.96	0.07	26,27,28,28	0
3	EDO	C	405	4/4	0.97	0.05	22,23,23,23	0
3	EDO	C	395	4/4	0.97	0.06	28,28,28,29	0
3	EDO	D	404	4/4	0.98	0.05	20,22,22,23	0
3	EDO	B	411	4/4	0.98	0.05	17,17,18,19	0
3	EDO	A	406	4/4	0.99	0.05	13,14,15,16	0

6.5 Other polymers [i](#)

There are no such residues in this entry.