



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 5, 2023 – 01:54 AM EST

PDB ID : 7KZ3
Title : Crystal structure of KabA from Bacillus cereus UW85 in complex with the internal aldimine
Authors : Prasertanan, T.; Palmer, D.R.J.; Sanders, D.A.R.
Deposited on : 2020-12-09
Resolution : 1.55 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
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) : 2.36

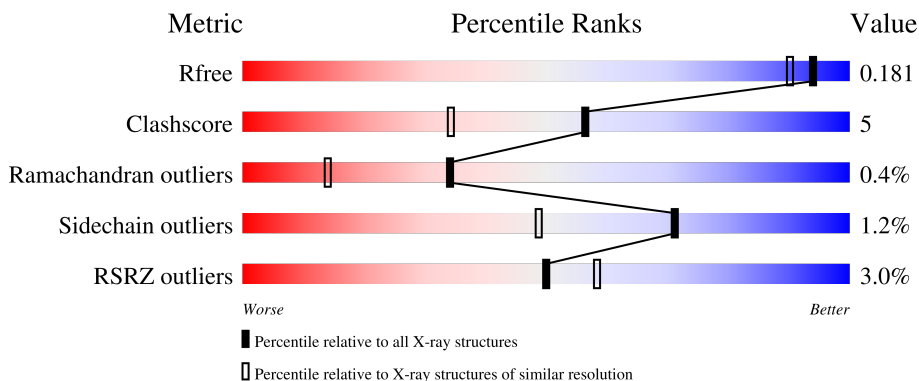
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.55 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	1483 (1.56-1.56)
Clashscore	141614	1529 (1.56-1.56)
Ramachandran outliers	138981	1498 (1.56-1.56)
Sidechain outliers	138945	1495 (1.56-1.56)
RSRZ outliers	127900	1465 (1.56-1.56)

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	445	
1	B	445	
1	C	445	
1	D	445	

2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 16029 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 Aminotransferase class I/II-fold pyridoxal phosphate-dependent enzyme.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	P	S			
1	A	439	Total 3583	C 2288	N 611	O 668	P 1	S 15	0	4	0
1	B	439	Total 3570	C 2280	N 609	O 665	P 1	S 15	0	2	0
1	C	439	Total 3569	C 2281	N 608	O 664	P 1	S 15	0	2	0
1	D	439	Total 3593	C 2298	N 611	O 668	P 1	S 15	0	5	0

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP C0JRF5
A	0	ALA	-	expression tag	UNP C0JRF5
A	1	MET	-	expression tag	UNP C0JRF5
A	2	ASP	-	expression tag	UNP C0JRF5
B	-1	GLY	-	expression tag	UNP C0JRF5
B	0	ALA	-	expression tag	UNP C0JRF5
B	1	MET	-	expression tag	UNP C0JRF5
B	2	ASP	-	expression tag	UNP C0JRF5
C	-1	GLY	-	expression tag	UNP C0JRF5
C	0	ALA	-	expression tag	UNP C0JRF5
C	1	MET	-	expression tag	UNP C0JRF5
C	2	ASP	-	expression tag	UNP C0JRF5
D	-1	GLY	-	expression tag	UNP C0JRF5
D	0	ALA	-	expression tag	UNP C0JRF5
D	1	MET	-	expression tag	UNP C0JRF5
D	2	ASP	-	expression tag	UNP C0JRF5

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



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

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

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

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

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	D	2	Total Na 2 2	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	407	Total O 407 407	0	0
4	B	340	Total O 340 340	0	0
4	C	301	Total O 301 301	0	0

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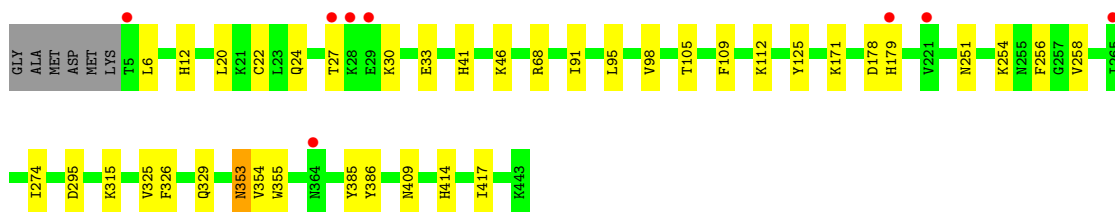
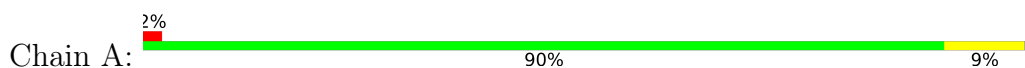
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	388	Total 388	O 388	0	0

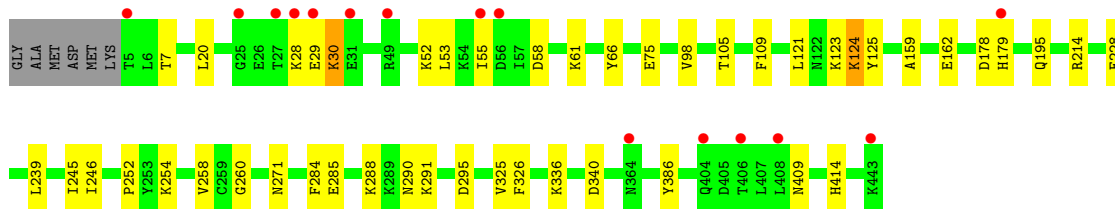
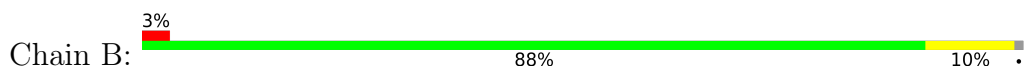
3 Residue-property plots [i](#)

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.

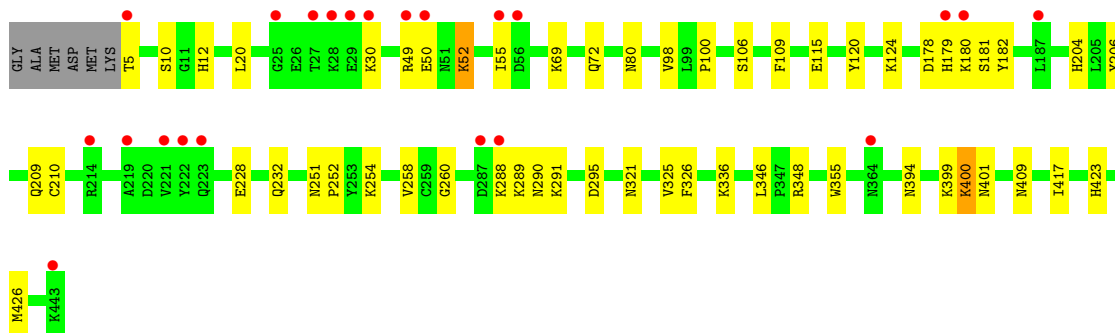
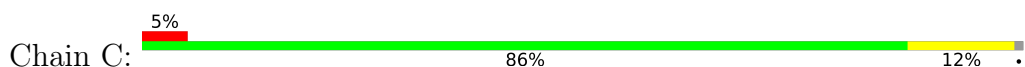
- Molecule 1: Aminotransferase class I/II-fold pyridoxal phosphate-dependent enzyme



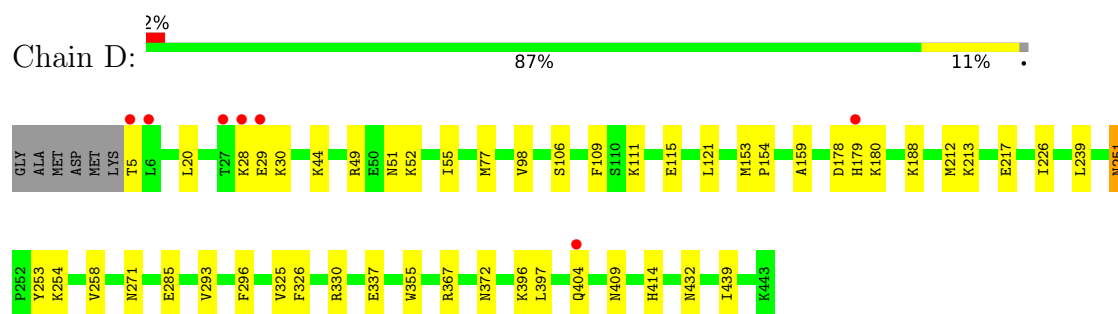
- Molecule 1: Aminotransferase class I/II-fold pyridoxal phosphate-dependent enzyme



- Molecule 1: Aminotransferase class I/II-fold pyridoxal phosphate-dependent enzyme



- Molecule 1: Aminotransferase class I/II-fold pyridoxal phosphate-dependent enzyme



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	66.46Å 66.68Å 111.83Å 81.08° 77.41° 88.05°	Depositor
Resolution (Å)	46.24 – 1.55 46.22 – 1.55	Depositor EDS
% Data completeness (in resolution range)	93.6 (46.24-1.55) 93.6 (46.22-1.55)	Depositor EDS
R_{merge}	0.03	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.52 (at 1.55Å)	Xtrriage
Refinement program	PHENIX dev_2398	Depositor
R, R_{free}	0.150 , 0.181 0.150 , 0.181	Depositor DCC
R_{free} test set	12572 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	17.3	Xtrriage
Anisotropy	0.172	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 52.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.015 for -k,-h,-l	Xtrriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	16029	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.47% 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, LLP, NA

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	0.84	0/3620	0.89	0/4880
1	B	0.79	0/3607	0.86	0/4862
1	C	0.77	0/3606	0.81	0/4861
1	D	0.84	0/3633	0.87	0/4897
All	All	0.81	0/14466	0.86	0/19500

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3583	0	3655	30	0
1	B	3570	0	3643	35	0
1	C	3569	0	3646	37	0
1	D	3593	0	3668	45	0
2	A	68	0	102	14	0
2	B	68	0	102	7	0
2	C	56	0	84	2	0
2	D	84	0	126	15	0
3	D	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	407	0	0	7	0
4	B	340	0	0	4	0
4	C	301	0	0	7	0
4	D	388	0	0	9	0
All	All	16029	0	15026	148	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 5.

All (148) 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:7:THR:OG1	1:B:30:LYS:HE2	1.55	1.07
1:B:105:THR:HG21	4:D:604:HOH:O	1.55	1.05
1:C:423:HIS:O	1:C:426:MET:HG2	1.58	1.03
1:D:330:ARG:HH22	1:D:432:ASN:HD21	1.13	0.94
1:B:121:LEU:HA	2:B:517:EDO:H22	1.53	0.91
1:C:295:ASP:H	2:C:507:EDO:H22	1.40	0.83
1:C:204:HIS:HD2	1:C:210:CYS:H	1.25	0.81
1:D:239:LEU:HB3	2:D:508:EDO:H11	1.62	0.81
1:A:105:THR:HB	2:A:506:EDO:H12	1.63	0.80
1:D:367:ARG:HH21	2:D:509:EDO:H11	1.48	0.78
1:D:121:LEU:HA	2:D:508:EDO:H12	1.66	0.78
1:C:204:HIS:CD2	1:C:210:CYS:H	2.04	0.75
1:D:325[A]:VAL:HG12	1:D:355:TRP:CD1	2.21	0.74
1:C:325[A]:VAL:HG12	1:C:355:TRP:HD1	1.54	0.72
1:D:296:PHE:HE1	2:D:501:EDO:H12	1.56	0.69
1:A:46:LYS:NZ	4:A:605:HOH:O	2.26	0.68
1:C:325[A]:VAL:HG12	1:C:355:TRP:CD1	2.28	0.68
1:C:400:LYS:HD3	1:C:401:ASN:OD1	1.96	0.66
1:D:271:ASN:HB2	4:D:951:HOH:O	1.97	0.65
1:D:296:PHE:CE1	2:D:501:EDO:H12	2.31	0.65
1:B:179:HIS:CE1	1:B:409[B]:ASN:HD21	2.15	0.64
1:D:251:ASN:ND2	4:D:604:HOH:O	2.29	0.64
4:C:611:HOH:O	1:D:49:ARG:HD3	1.97	0.64
1:D:111:LYS:NZ	1:D:115:GLU:HG3	2.13	0.63
1:D:325[A]:VAL:HG12	1:D:355:TRP:HD1	1.59	0.63
1:D:330:ARG:NH2	1:D:432:ASN:HD21	1.91	0.63
1:A:325[A]:VAL:HG12	1:A:355:TRP:HD1	1.65	0.62
1:A:178:ASP:O	1:A:409[B]:ASN:ND2	2.33	0.62
1:D:367:ARG:NH2	2:D:509:EDO:H11	2.13	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:204:HIS:HE1	1:C:228:GLU:OE1	1.83	0.61
1:A:329:GLN:NE2	4:A:601:HOH:O	2.23	0.60
1:A:325[A]:VAL:HG12	1:A:355:TRP:CD1	2.36	0.59
1:D:212:MET:H	2:D:514:EDO:H21	1.67	0.59
1:B:58:ASP:CG	1:B:61:LYS:HE2	2.24	0.59
1:C:206:TYR:H	1:C:232:GLN:NE2	2.01	0.58
1:A:353:ASN:ND2	1:A:355:TRP:H	2.00	0.58
1:D:98:VAL:HG22	1:D:109:PHE:CE1	2.38	0.58
1:D:111:LYS:HZ1	1:D:115:GLU:HG3	1.68	0.57
1:A:125:TYR:HA	2:A:516:EDO:H21	1.87	0.57
1:B:179:HIS:CE1	1:B:409[B]:ASN:ND2	2.72	0.57
1:A:385:TYR:CE2	2:A:509:EDO:H12	2.41	0.56
1:A:24:GLN:HE22	1:A:329:GLN:HE22	1.53	0.56
1:D:285:GLU:HB2	1:D:293[B]:VAL:CG2	2.36	0.56
1:B:125:TYR:HA	2:B:513:EDO:H12	1.87	0.55
1:A:325[A]:VAL:HG11	4:A:759:HOH:O	2.06	0.55
1:B:121:LEU:HA	2:B:517:EDO:C2	2.31	0.55
1:D:49:ARG:HA	2:D:507:EDO:H12	1.89	0.55
1:B:98:VAL:HG22	1:B:109:PHE:CE1	2.42	0.55
1:C:52:LYS:HD2	4:C:619:HOH:O	2.07	0.54
1:A:91:ILE:O	1:A:95:LEU:HG	2.08	0.54
1:A:274:ILE:HG21	2:A:513:EDO:H21	1.90	0.54
1:C:182:TYR:CE1	1:C:417:ILE:HD12	2.43	0.54
1:B:325:VAL:HG21	4:B:769:HOH:O	2.06	0.54
1:D:285:GLU:HB2	1:D:293[B]:VAL:HG21	1.90	0.54
1:B:7:THR:CB	1:B:30:LYS:HE2	2.38	0.53
1:A:385:TYR:HE2	2:A:509:EDO:H12	1.73	0.53
1:B:336:LYS:HE3	1:B:340:ASP:OD2	2.08	0.53
2:A:517:EDO:H21	1:C:295:ASP:OD2	2.08	0.53
1:A:20:LEU:HD11	1:A:326:PHE:HD1	1.73	0.52
1:B:285:GLU:HB3	1:B:288:LYS:HB2	1.91	0.52
1:C:10:SER:OG	1:C:12:HIS:HD2	1.92	0.52
2:A:512:EDO:H21	4:A:648:HOH:O	2.07	0.52
1:D:285:GLU:OE1	2:D:517:EDO:H12	2.09	0.52
1:C:20:LEU:HD11	1:C:326:PHE:HD1	1.75	0.52
1:C:5:THR:N	4:C:607:HOH:O	2.43	0.51
1:A:6:LEU:HD21	1:A:33:GLU:HG3	1.93	0.51
1:A:179:HIS:CE1	1:A:409[A]:ASN:OD1	2.64	0.51
2:A:508:EDO:H22	1:C:100:PRO:HA	1.93	0.51
1:C:69:LYS:HD3	1:D:55:ILE:O	2.10	0.51
1:C:325[A]:VAL:HG11	4:C:706:HOH:O	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:171:LYS:HG3	4:A:869:HOH:O	2.10	0.50
1:B:295:ASP:H	2:B:514:EDO:H22	1.75	0.50
1:B:52:LYS:HD3	1:B:53:LEU:O	2.11	0.50
1:C:50:GLU:O	1:C:52:LYS:HD3	2.11	0.50
1:A:27:THR:HG22	1:A:30:LYS:HG3	1.94	0.49
1:D:178:ASP:O	1:D:409[B]:ASN:ND2	2.45	0.49
1:C:49:ARG:HH12	1:C:55:ILE:HG13	1.77	0.49
1:A:295:ASP:OD2	2:A:511:EDO:H22	2.12	0.49
1:A:6:LEU:HD11	1:A:33:GLU:HG3	1.94	0.49
1:C:252:PRO:HD3	1:C:260:GLY:O	2.13	0.49
1:A:98:VAL:HG22	1:A:109:PHE:CE1	2.47	0.48
1:D:77:MET:SD	1:D:253[B]:TYR:HD1	2.37	0.48
1:A:417:ILE:HD13	2:A:503:EDO:H11	1.95	0.48
1:B:66:TYR:CZ	2:B:509:EDO:H12	2.48	0.48
2:A:510:EDO:O2	1:C:290:ASN:HB3	2.13	0.48
1:B:239:LEU:HB3	2:B:517:EDO:H11	1.95	0.48
1:D:213:LYS:O	1:D:217:GLU:HG3	2.14	0.48
1:B:178:ASP:O	1:B:409[A]:ASN:ND2	2.46	0.48
1:C:346:LEU:HB2	1:C:348:ARG:HH22	1.78	0.47
1:C:12:HIS:HE1	1:C:120:TYR:O	1.98	0.47
1:A:41[A]:HIS:HB3	1:A:315:LYS:HB3	1.96	0.46
1:C:124:LYS:HB2	1:C:124:LYS:HE3	1.46	0.46
1:C:178:ASP:OD2	1:C:181:SER:OG	2.29	0.46
1:C:98:VAL:HG22	1:C:109:PHE:CE1	2.50	0.46
1:B:228:GLU:HB2	1:B:246:ILE:HD12	1.98	0.46
1:C:325[B]:VAL:HG21	4:C:706:HOH:O	2.15	0.46
2:A:506:EDO:H21	4:A:676:HOH:O	2.15	0.46
1:D:325[A]:VAL:HG11	4:D:691:HOH:O	2.15	0.46
1:D:414:HIS:ND1	2:D:518:EDO:H11	2.32	0.45
1:B:7:THR:HG1	1:B:30:LYS:HE2	1.71	0.44
1:C:80:ASN:OD1	1:D:49:ARG:CZ	2.65	0.44
1:A:353:ASN:HD22	1:A:354:VAL:N	2.15	0.44
1:C:321:ASN:O	1:C:325[B]:VAL:HG23	2.17	0.44
1:D:20:LEU:HD11	1:D:326:PHE:HD1	1.82	0.44
1:D:153:MET:HB2	1:D:154:PRO:HD2	1.99	0.44
1:D:212:MET:HE1	1:D:226:ILE:HG21	1.99	0.44
1:A:386:TYR:O	1:A:414:HIS:HE1	1.99	0.44
1:D:51:ASN:O	1:D:52:LYS:HE3	2.18	0.43
1:B:291:LYS:HG3	4:D:898:HOH:O	2.17	0.43
1:D:106:SER:HB3	4:D:923:HOH:O	2.19	0.43
1:D:111:LYS:HA	1:D:111:LYS:HD2	1.76	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:75:GLU:OE1	2:B:516:EDO:H21	2.18	0.43
1:B:179:HIS:HE1	1:B:409[B]:ASN:HD21	1.60	0.43
1:B:123:LYS:HE3	4:B:759:HOH:O	2.18	0.43
1:C:72:GLN:NE2	1:D:29:GLU:OE2	2.52	0.42
1:D:396:LYS:HE2	1:D:397:LEU:N	2.34	0.42
1:A:112:LYS:NZ	4:A:608:HOH:O	2.40	0.42
1:A:256:PHE:HA	2:A:505:EDO:H22	2.00	0.42
1:C:295:ASP:H	2:C:507:EDO:C2	2.19	0.42
1:B:28:LYS:HD3	1:B:28:LYS:HA	1.73	0.42
1:B:245:ILE:C	1:B:246:ILE:HD13	2.40	0.42
1:C:115:GLU:HG3	4:C:755:HOH:O	2.19	0.42
1:B:195:GLN:HG2	4:B:876:HOH:O	2.18	0.42
1:B:252:PRO:HD3	1:B:260:GLY:O	2.20	0.42
2:D:501:EDO:H11	2:D:516:EDO:O2	2.20	0.42
1:D:5:THR:HG23	1:D:30:LYS:HG3	2.02	0.42
1:B:105:THR:CG2	4:D:604:HOH:O	2.35	0.42
1:A:12:HIS:CD2	2:A:512:EDO:H11	2.55	0.41
1:D:179:HIS:NE2	1:D:409[A]:ASN:OD1	2.53	0.41
1:C:179:HIS:CE1	1:C:409:ASN:OD1	2.73	0.41
1:C:399:LYS:HB3	1:C:399:LYS:HE2	1.59	0.41
1:D:179:HIS:CE1	1:D:409[A]:ASN:OD1	2.73	0.41
1:D:44:LYS:HG2	2:D:507:EDO:H21	2.02	0.41
1:D:372:ASN:HD21	2:D:510:EDO:C2	2.33	0.41
1:C:106:SER:O	4:C:601:HOH:O	2.21	0.41
1:C:204:HIS:CD2	1:C:209:GLN:HA	2.55	0.41
1:B:124:LYS:HE2	1:B:124:LYS:HB2	1.51	0.41
1:B:271:ASN:ND2	4:B:612:HOH:O	2.53	0.41
1:D:77:MET:SD	1:D:253[B]:TYR:CD1	3.13	0.41
1:D:337:GLU:OE1	1:D:439:ILE:HD12	2.21	0.41
1:A:22:CYS:HA	1:A:27:THR:OG1	2.20	0.41
1:B:386:TYR:O	1:B:414:HIS:HE1	2.03	0.41
1:B:20:LEU:HD11	1:B:326:PHE:HD1	1.86	0.40
1:B:284:PHE:CE1	1:B:290:ASN:HA	2.55	0.40
1:D:159:ALA:HB2	2:D:506:EDO:H22	2.02	0.40
1:B:159:ALA:O	1:B:162:GLU:HG2	2.21	0.40
1:D:188:LYS:HE3	4:D:631:HOH:O	2.21	0.40
2:D:516:EDO:H12	4:D:722:HOH:O	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	440/445 (99%)	426 (97%)	13 (3%)	1 (0%)	47	23
1	B	438/445 (98%)	422 (96%)	15 (3%)	1 (0%)	47	23
1	C	438/445 (98%)	427 (98%)	9 (2%)	2 (0%)	29	9
1	D	441/445 (99%)	425 (96%)	13 (3%)	3 (1%)	22	5
All	All	1757/1780 (99%)	1700 (97%)	50 (3%)	7 (0%)	34	14

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	28	LYS
1	C	180	LYS
1	D	180	LYS
1	C	258	VAL
1	D	258	VAL
1	A	258	VAL
1	B	258	VAL

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	400/400 (100%)	397 (99%)	3 (1%)	81	66
1	B	398/400 (100%)	393 (99%)	5 (1%)	69	44
1	C	398/400 (100%)	389 (98%)	9 (2%)	50	21

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	401/400 (100%)	399 (100%)	2 (0%)	88	78
All	All	1597/1600 (100%)	1578 (99%)	19 (1%)	71	49

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	68	ARG
1	A	251	ASN
1	A	353	ASN
1	B	29	GLU
1	B	30	LYS
1	B	55	ILE
1	B	124	LYS
1	B	214	ARG
1	C	30	LYS
1	C	52	LYS
1	C	251	ASN
1	C	288	LYS
1	C	289	LYS
1	C	291	LYS
1	C	336	LYS
1	C	394	ASN
1	C	400	LYS
1	D	251	ASN
1	D	404	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	ASN
1	A	103	GLN
1	A	179	HIS
1	A	278	GLN
1	A	329	GLN
1	A	353	ASN
1	A	372	ASN
1	A	440	ASN
1	B	103	GLN
1	B	179	HIS
1	C	12	HIS
1	C	16	ASN

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Mol	Chain	Res	Type
1	C	103	GLN
1	C	204	HIS
1	C	232	GLN
1	C	394	ASN
1	D	16	ASN
1	D	372	ASN
1	D	432	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](#)

4 non-standard protein/DNA/RNA residues 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
1	LLP	D	254	1	23,24,25	2.21	6 (26%)	25,32,34	2.60	8 (32%)
1	LLP	C	254	1	23,24,25	1.99	6 (26%)	25,32,34	2.54	8 (32%)
1	LLP	B	254	1	23,24,25	1.95	5 (21%)	25,32,34	2.40	10 (40%)
1	LLP	A	254	1	23,24,25	1.77	6 (26%)	25,32,34	2.72	8 (32%)

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
1	LLP	D	254	1	-	4/16/17/19	0/1/1/1
1	LLP	C	254	1	-	4/16/17/19	0/1/1/1
1	LLP	B	254	1	-	5/16/17/19	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	LLP	A	254	1	-	4/16/17/19	0/1/1/1

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	254	LLP	C4-C4'	6.47	1.58	1.46
1	D	254	LLP	C4-C4'	5.96	1.58	1.46
1	B	254	LLP	C4-C4'	5.87	1.57	1.46
1	D	254	LLP	C4-C5	-4.97	1.35	1.42
1	A	254	LLP	C4-C4'	4.36	1.54	1.46
1	B	254	LLP	C2'-C2	3.35	1.56	1.50
1	D	254	LLP	C2-N1	-3.31	1.27	1.33
1	C	254	LLP	CE-NZ	2.95	1.53	1.46
1	B	254	LLP	C4-C5	-2.90	1.38	1.42
1	C	254	LLP	C2-N1	-2.75	1.28	1.33
1	C	254	LLP	C4'-NZ	2.72	1.36	1.27
1	A	254	LLP	C4-C5	-2.68	1.38	1.42
1	B	254	LLP	C4'-NZ	2.55	1.35	1.27
1	A	254	LLP	C2'-C2	2.48	1.54	1.50
1	A	254	LLP	CD-CE	2.45	1.60	1.51
1	D	254	LLP	C4'-NZ	2.39	1.35	1.27
1	B	254	LLP	CE-NZ	2.38	1.52	1.46
1	A	254	LLP	CE-NZ	2.33	1.51	1.46
1	C	254	LLP	CD-CE	2.23	1.59	1.51
1	D	254	LLP	CD-CE	2.20	1.59	1.51
1	C	254	LLP	C2'-C2	2.16	1.54	1.50
1	A	254	LLP	C4'-NZ	2.10	1.34	1.27
1	D	254	LLP	CE-NZ	2.04	1.51	1.46

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	254	LLP	CD-CE-NZ	8.20	131.03	110.93
1	A	254	LLP	CD-CE-NZ	8.03	130.61	110.93
1	D	254	LLP	CD-CE-NZ	7.48	129.25	110.93
1	A	254	LLP	CE-NZ-C4'	5.62	136.17	118.90
1	C	254	LLP	CE-NZ-C4'	5.55	135.95	118.90
1	B	254	LLP	CE-NZ-C4'	5.46	135.68	118.90
1	D	254	LLP	CE-NZ-C4'	5.19	134.83	118.90
1	C	254	LLP	C3-C4-C5	-4.83	114.56	118.26
1	B	254	LLP	CD-CE-NZ	4.73	122.51	110.93
1	D	254	LLP	C5-C4-C4'	4.69	129.27	121.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	254	LLP	C4-C4'-NZ	-4.22	104.92	124.31
1	A	254	LLP	C3-C4-C5	-4.19	115.04	118.26
1	B	254	LLP	C3-C4-C5	-4.19	115.05	118.26
1	D	254	LLP	C3-C4-C4'	-3.93	113.09	120.41
1	A	254	LLP	C5-C4-C4'	3.79	127.78	121.56
1	D	254	LLP	C4-C4'-NZ	-3.69	107.35	124.31
1	A	254	LLP	C5'-C5-C6	-3.49	113.63	119.37
1	A	254	LLP	C4-C4'-NZ	-3.45	108.48	124.31
1	D	254	LLP	C5'-C5-C6	-3.14	114.21	119.37
1	B	254	LLP	C5-C6-N1	-3.12	118.62	123.82
1	B	254	LLP	OP3-P-OP4	-2.87	99.10	106.73
1	B	254	LLP	C5-C4-C4'	2.83	126.21	121.56
1	A	254	LLP	O3-C3-C2	2.76	123.51	117.49
1	B	254	LLP	CD-CG-CB	2.73	123.29	113.62
1	C	254	LLP	C4-C3-C2	2.63	121.82	120.19
1	C	254	LLP	CG-CD-CE	2.56	122.49	113.57
1	D	254	LLP	C5-C6-N1	-2.54	119.59	123.82
1	B	254	LLP	C6-C5-C4	2.38	122.53	118.15
1	D	254	LLP	O3-C3-C2	2.25	122.40	117.49
1	C	254	LLP	OP4-P-OP1	-2.18	100.36	106.47
1	A	254	LLP	OP3-P-OP4	-2.13	101.08	106.73
1	C	254	LLP	OP4-C5'-C5	2.10	113.35	109.35
1	B	254	LLP	C6-N1-C2	2.09	123.04	119.17
1	C	254	LLP	CD-CG-CB	2.04	120.86	113.62

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	C	254	LLP	C4-C4'-NZ-CE
1	A	254	LLP	CG-CD-CE-NZ
1	B	254	LLP	CG-CD-CE-NZ
1	D	254	LLP	CG-CD-CE-NZ
1	D	254	LLP	C3-C4-C4'-NZ
1	C	254	LLP	CD-CE-NZ-C4'
1	B	254	LLP	CD-CE-NZ-C4'
1	A	254	LLP	CD-CE-NZ-C4'
1	D	254	LLP	CD-CE-NZ-C4'
1	A	254	LLP	C3-C4-C4'-NZ
1	B	254	LLP	C3-C4-C4'-NZ
1	C	254	LLP	C3-C4-C4'-NZ
1	A	254	LLP	C5-C4-C4'-NZ

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Mol	Chain	Res	Type	Atoms
1	B	254	LLP	CE-CD-CG-CB
1	B	254	LLP	C5-C4-C4'-NZ
1	C	254	LLP	C5-C4-C4'-NZ
1	D	254	LLP	C5-C4-C4'-NZ

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 71 ligands modelled in this entry, 2 are monoatomic - leaving 69 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	A	510	-	3,3,3	0.67	0	2,2,2	0.29	0
2	EDO	B	517	-	3,3,3	0.88	0	2,2,2	1.67	1 (50%)
2	EDO	A	517	-	3,3,3	0.74	0	2,2,2	0.63	0
2	EDO	C	507	-	3,3,3	0.23	0	2,2,2	0.27	0
2	EDO	D	516	-	3,3,3	0.58	0	2,2,2	0.35	0
2	EDO	B	511	-	3,3,3	0.28	0	2,2,2	0.25	0
2	EDO	A	515	-	3,3,3	0.43	0	2,2,2	0.08	0
2	EDO	B	505	-	3,3,3	0.65	0	2,2,2	0.40	0
2	EDO	B	514	-	3,3,3	0.27	0	2,2,2	0.56	0
2	EDO	D	513	-	3,3,3	0.34	0	2,2,2	0.44	0
2	EDO	A	511	-	3,3,3	0.32	0	2,2,2	0.10	0
2	EDO	C	511	-	3,3,3	0.28	0	2,2,2	0.78	0
2	EDO	D	504	-	3,3,3	0.51	0	2,2,2	0.10	0
2	EDO	B	509	-	3,3,3	0.25	0	2,2,2	0.88	0
2	EDO	A	509	-	3,3,3	0.23	0	2,2,2	1.28	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	EDO	D	510	-	3,3,3	0.34	0	2,2,2	0.65	0
2	EDO	D	509	-	3,3,3	0.32	0	2,2,2	0.20	0
2	EDO	C	502	-	3,3,3	0.36	0	2,2,2	0.19	0
2	EDO	B	513	-	3,3,3	0.30	0	2,2,2	0.72	0
2	EDO	C	506	-	3,3,3	0.46	0	2,2,2	0.30	0
2	EDO	C	513	-	3,3,3	0.36	0	2,2,2	0.50	0
2	EDO	D	515	-	3,3,3	0.52	0	2,2,2	0.66	0
2	EDO	A	506	-	3,3,3	0.46	0	2,2,2	1.00	0
2	EDO	D	503	-	3,3,3	0.36	0	2,2,2	0.55	0
2	EDO	A	504	-	3,3,3	0.25	0	2,2,2	0.53	0
2	EDO	D	505	-	3,3,3	0.50	0	2,2,2	0.63	0
2	EDO	B	502	-	3,3,3	0.52	0	2,2,2	0.39	0
2	EDO	D	512	-	3,3,3	0.22	0	2,2,2	0.66	0
2	EDO	D	508	-	3,3,3	0.38	0	2,2,2	0.50	0
2	EDO	C	503	-	3,3,3	0.27	0	2,2,2	0.61	0
2	EDO	A	503	-	3,3,3	0.47	0	2,2,2	0.73	0
2	EDO	C	512	-	3,3,3	0.35	0	2,2,2	0.39	0
2	EDO	A	505	-	3,3,3	0.84	0	2,2,2	0.32	0
2	EDO	A	516	-	3,3,3	0.42	0	2,2,2	1.04	0
2	EDO	B	503	-	3,3,3	0.31	0	2,2,2	0.74	0
2	EDO	B	504	-	3,3,3	0.44	0	2,2,2	0.54	0
2	EDO	C	505	-	3,3,3	0.77	0	2,2,2	0.81	0
2	EDO	B	512	-	3,3,3	0.33	0	2,2,2	0.45	0
2	EDO	A	508	-	3,3,3	0.40	0	2,2,2	0.59	0
2	EDO	C	508	-	3,3,3	0.42	0	2,2,2	0.52	0
2	EDO	D	511	-	3,3,3	0.23	0	2,2,2	0.66	0
2	EDO	D	502	-	3,3,3	0.26	0	2,2,2	0.60	0
2	EDO	D	506	-	3,3,3	0.47	0	2,2,2	0.40	0
2	EDO	B	516	-	3,3,3	0.29	0	2,2,2	0.78	0
2	EDO	D	518	-	3,3,3	0.44	0	2,2,2	0.39	0
2	EDO	B	508	-	3,3,3	0.44	0	2,2,2	0.52	0
2	EDO	D	514	-	3,3,3	0.40	0	2,2,2	0.17	0
2	EDO	D	520	-	3,3,3	0.38	0	2,2,2	0.40	0
2	EDO	C	514	-	3,3,3	0.24	0	2,2,2	0.55	0
2	EDO	D	501	-	3,3,3	0.36	0	2,2,2	0.28	0
2	EDO	D	519	-	3,3,3	0.35	0	2,2,2	0.18	0
2	EDO	C	501	-	3,3,3	0.39	0	2,2,2	0.53	0
2	EDO	A	514	-	3,3,3	0.33	0	2,2,2	0.69	0
2	EDO	A	513	-	3,3,3	0.58	0	2,2,2	0.47	0
2	EDO	D	521	-	3,3,3	0.57	0	2,2,2	0.52	0
2	EDO	B	506	-	3,3,3	0.36	0	2,2,2	0.17	0
2	EDO	D	517	-	3,3,3	0.32	0	2,2,2	0.16	0
2	EDO	C	509	-	3,3,3	0.28	0	2,2,2	0.89	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	EDO	C	510	-	3,3,3	0.29	0	2,2,2	0.58	0
2	EDO	C	504	-	3,3,3	0.29	0	2,2,2	0.98	0
2	EDO	A	502	-	3,3,3	0.33	0	2,2,2	1.43	0
2	EDO	A	507	-	3,3,3	0.31	0	2,2,2	0.47	0
2	EDO	B	515	-	3,3,3	0.37	0	2,2,2	0.74	0
2	EDO	B	507	-	3,3,3	0.39	0	2,2,2	0.43	0
2	EDO	D	507	-	3,3,3	0.37	0	2,2,2	0.23	0
2	EDO	A	501	-	3,3,3	0.41	0	2,2,2	0.30	0
2	EDO	A	512	-	3,3,3	0.50	0	2,2,2	0.73	0
2	EDO	B	501	-	3,3,3	0.22	0	2,2,2	1.36	0
2	EDO	B	510	-	3,3,3	0.33	0	2,2,2	0.45	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	A	510	-	-	0/1/1/1	-
2	EDO	B	517	-	-	1/1/1/1	-
2	EDO	A	517	-	-	0/1/1/1	-
2	EDO	C	507	-	-	0/1/1/1	-
2	EDO	D	516	-	-	0/1/1/1	-
2	EDO	B	511	-	-	0/1/1/1	-
2	EDO	A	515	-	-	1/1/1/1	-
2	EDO	B	505	-	-	0/1/1/1	-
2	EDO	B	514	-	-	0/1/1/1	-
2	EDO	D	513	-	-	0/1/1/1	-
2	EDO	A	511	-	-	1/1/1/1	-
2	EDO	C	511	-	-	0/1/1/1	-
2	EDO	D	504	-	-	0/1/1/1	-
2	EDO	B	509	-	-	1/1/1/1	-
2	EDO	A	509	-	-	1/1/1/1	-
2	EDO	D	510	-	-	1/1/1/1	-
2	EDO	D	509	-	-	1/1/1/1	-
2	EDO	C	502	-	-	0/1/1/1	-
2	EDO	B	513	-	-	1/1/1/1	-
2	EDO	C	506	-	-	0/1/1/1	-
2	EDO	C	513	-	-	0/1/1/1	-
2	EDO	D	515	-	-	1/1/1/1	-
2	EDO	A	506	-	-	0/1/1/1	-
2	EDO	D	503	-	-	0/1/1/1	-
2	EDO	A	504	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	EDO	D	505	-	-	0/1/1/1	-
2	EDO	B	502	-	-	0/1/1/1	-
2	EDO	D	512	-	-	1/1/1/1	-
2	EDO	D	508	-	-	1/1/1/1	-
2	EDO	C	503	-	-	1/1/1/1	-
2	EDO	A	503	-	-	0/1/1/1	-
2	EDO	C	512	-	-	0/1/1/1	-
2	EDO	A	505	-	-	0/1/1/1	-
2	EDO	A	516	-	-	1/1/1/1	-
2	EDO	B	503	-	-	0/1/1/1	-
2	EDO	B	504	-	-	0/1/1/1	-
2	EDO	C	505	-	-	1/1/1/1	-
2	EDO	B	512	-	-	1/1/1/1	-
2	EDO	A	508	-	-	0/1/1/1	-
2	EDO	C	508	-	-	1/1/1/1	-
2	EDO	D	511	-	-	0/1/1/1	-
2	EDO	D	502	-	-	1/1/1/1	-
2	EDO	D	506	-	-	1/1/1/1	-
2	EDO	B	516	-	-	1/1/1/1	-
2	EDO	D	518	-	-	1/1/1/1	-
2	EDO	B	508	-	-	0/1/1/1	-
2	EDO	D	514	-	-	0/1/1/1	-
2	EDO	D	520	-	-	1/1/1/1	-
2	EDO	C	514	-	-	0/1/1/1	-
2	EDO	D	501	-	-	1/1/1/1	-
2	EDO	D	519	-	-	0/1/1/1	-
2	EDO	C	501	-	-	1/1/1/1	-
2	EDO	A	514	-	-	1/1/1/1	-
2	EDO	A	513	-	-	1/1/1/1	-
2	EDO	D	521	-	-	1/1/1/1	-
2	EDO	B	506	-	-	1/1/1/1	-
2	EDO	D	517	-	-	1/1/1/1	-
2	EDO	C	509	-	-	1/1/1/1	-
2	EDO	C	510	-	-	1/1/1/1	-
2	EDO	C	504	-	-	0/1/1/1	-
2	EDO	A	502	-	-	0/1/1/1	-
2	EDO	A	507	-	-	0/1/1/1	-
2	EDO	B	515	-	-	0/1/1/1	-
2	EDO	B	507	-	-	0/1/1/1	-
2	EDO	D	507	-	-	1/1/1/1	-
2	EDO	A	501	-	-	0/1/1/1	-
2	EDO	A	512	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	EDO	B	501	-	-	0/1/1/1	-
2	EDO	B	510	-	-	1/1/1/1	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	517	EDO	O1-C1-C2	2.14	127.33	111.91

There are no chirality outliers.

All (34) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	513	EDO	O1-C1-C2-O2
2	B	509	EDO	O1-C1-C2-O2
2	D	521	EDO	O1-C1-C2-O2
2	A	512	EDO	O1-C1-C2-O2
2	C	508	EDO	O1-C1-C2-O2
2	C	510	EDO	O1-C1-C2-O2
2	D	509	EDO	O1-C1-C2-O2
2	D	518	EDO	O1-C1-C2-O2
2	D	501	EDO	O1-C1-C2-O2
2	A	516	EDO	O1-C1-C2-O2
2	B	517	EDO	O1-C1-C2-O2
2	C	503	EDO	O1-C1-C2-O2
2	D	507	EDO	O1-C1-C2-O2
2	D	517	EDO	O1-C1-C2-O2
2	A	504	EDO	O1-C1-C2-O2
2	A	509	EDO	O1-C1-C2-O2
2	A	515	EDO	O1-C1-C2-O2
2	B	506	EDO	O1-C1-C2-O2
2	B	513	EDO	O1-C1-C2-O2
2	C	505	EDO	O1-C1-C2-O2
2	D	515	EDO	O1-C1-C2-O2
2	A	511	EDO	O1-C1-C2-O2
2	D	506	EDO	O1-C1-C2-O2
2	D	508	EDO	O1-C1-C2-O2
2	A	514	EDO	O1-C1-C2-O2
2	B	510	EDO	O1-C1-C2-O2
2	D	510	EDO	O1-C1-C2-O2
2	B	512	EDO	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
2	B	516	EDO	O1-C1-C2-O2
2	C	501	EDO	O1-C1-C2-O2
2	C	509	EDO	O1-C1-C2-O2
2	D	502	EDO	O1-C1-C2-O2
2	D	512	EDO	O1-C1-C2-O2
2	D	520	EDO	O1-C1-C2-O2

There are no ring outliers.

27 monomers are involved in 38 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	510	EDO	1	0
2	B	517	EDO	3	0
2	A	517	EDO	1	0
2	C	507	EDO	2	0
2	D	516	EDO	2	0
2	B	514	EDO	1	0
2	A	511	EDO	1	0
2	B	509	EDO	1	0
2	A	509	EDO	2	0
2	D	510	EDO	1	0
2	D	509	EDO	2	0
2	B	513	EDO	1	0
2	A	506	EDO	2	0
2	D	508	EDO	2	0
2	A	503	EDO	1	0
2	A	505	EDO	1	0
2	A	516	EDO	1	0
2	A	508	EDO	1	0
2	D	506	EDO	1	0
2	B	516	EDO	1	0
2	D	518	EDO	1	0
2	D	514	EDO	1	0
2	D	501	EDO	3	0
2	A	513	EDO	1	0
2	D	517	EDO	1	0
2	D	507	EDO	2	0
2	A	512	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 [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	438/445 (98%)	-0.10	8 (1%) 68 74	11, 19, 37, 76	0
1	B	438/445 (98%)	-0.15	15 (3%) 45 52	12, 22, 43, 84	0
1	C	438/445 (98%)	-0.05	22 (5%) 28 33	12, 24, 45, 87	0
1	D	438/445 (98%)	-0.22	7 (1%) 72 77	11, 19, 40, 77	0
All	All	1752/1780 (98%)	-0.13	52 (2%) 50 58	11, 21, 42, 87	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	55	ILE	5.6
1	C	27	THR	5.1
1	C	29	GLU	5.0
1	B	29	GLU	4.8
1	D	5	THR	4.7
1	C	179	HIS	4.6
1	B	28	LYS	4.5
1	A	5	THR	4.5
1	C	221	VAL	4.3
1	C	28	LYS	4.2
1	B	179	HIS	3.8
1	A	27	THR	3.7
1	C	56	ASP	3.7
1	C	55	ILE	3.6
1	D	29	GLU	3.5
1	C	5	THR	3.4
1	B	27	THR	3.3
1	C	187	LEU	3.3
1	D	179	HIS	3.2
1	B	56	ASP	3.1
1	A	179	HIS	3.1

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Mol	Chain	Res	Type	RSRZ
1	C	223	GLN	3.0
1	B	404	GLN	3.0
1	C	222	TYR	3.0
1	A	28	LYS	2.9
1	C	30	LYS	2.8
1	A	29	GLU	2.8
1	C	49	ARG	2.8
1	D	27	THR	2.8
1	D	28	LYS	2.7
1	B	408	LEU	2.7
1	C	288	LYS	2.7
1	A	364	ASN	2.7
1	B	406	THR	2.7
1	B	443	LYS	2.6
1	C	219	ALA	2.5
1	D	6	LEU	2.5
1	B	364	ASN	2.5
1	C	214	ARG	2.4
1	C	364	ASN	2.4
1	D	404	GLN	2.4
1	B	49	ARG	2.4
1	C	50	GLU	2.4
1	C	180	LYS	2.4
1	C	25	GLY	2.4
1	A	265	ILE	2.2
1	C	287	ASP	2.2
1	B	25	GLY	2.1
1	B	5	THR	2.1
1	B	31	GLU	2.0
1	A	221	VAL	2.0
1	C	443	LYS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	LLP	D	254	24/25	0.97	0.12	10,13,32,41	0
1	LLP	B	254	24/25	0.98	0.10	11,15,24,41	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
1	LLP	C	254	24/25	0.98	0.10	13,15,24,45	0
1	LLP	A	254	24/25	0.98	0.11	11,13,28,36	0

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(\AA^2)	Q<0.9
2	EDO	D	503	4/4	0.56	0.23	20,28,31,41	0
2	EDO	D	521	4/4	0.68	0.25	20,20,20,20	0
2	EDO	B	507	4/4	0.69	0.18	33,40,40,45	0
2	EDO	D	501	4/4	0.70	0.21	45,47,50,52	0
2	EDO	D	510	4/4	0.73	0.11	48,53,55,57	0
2	EDO	A	506	4/4	0.73	0.21	23,24,31,37	0
2	EDO	B	510	4/4	0.79	0.14	49,49,50,51	0
2	EDO	B	512	4/4	0.79	0.16	68,68,68,70	0
2	EDO	A	517	4/4	0.79	0.15	31,33,35,40	0
2	EDO	B	513	4/4	0.81	0.18	52,55,56,58	0
2	EDO	C	504	4/4	0.81	0.13	26,27,30,33	0
2	EDO	A	516	4/4	0.81	0.17	30,32,37,41	0
2	EDO	A	515	4/4	0.82	0.17	30,47,52,60	0
2	EDO	B	508	4/4	0.82	0.15	38,42,45,47	0
2	EDO	A	507	4/4	0.82	0.28	54,60,61,66	0
2	EDO	A	513	4/4	0.82	0.20	33,35,42,44	0
2	EDO	B	501	4/4	0.82	0.14	24,27,32,37	0
2	EDO	D	508	4/4	0.83	0.25	26,28,30,42	0
2	EDO	A	502	4/4	0.83	0.11	24,27,29,31	0
2	EDO	C	513	4/4	0.83	0.18	27,40,40,51	0
2	EDO	C	508	4/4	0.84	0.10	28,37,39,43	0
2	EDO	D	517	4/4	0.85	0.18	28,44,46,47	0
2	EDO	D	520	4/4	0.85	0.27	20,20,20,20	0
2	EDO	A	508	4/4	0.85	0.15	33,34,35,39	0
2	EDO	D	512	4/4	0.86	0.12	40,49,50,63	0
2	EDO	B	515	4/4	0.86	0.10	38,40,41,48	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	EDO	B	517	4/4	0.86	0.17	26,28,34,38	0
2	EDO	D	502	4/4	0.86	0.12	27,33,34,35	0
2	EDO	C	506	4/4	0.87	0.13	26,33,43,46	0
2	EDO	C	512	4/4	0.87	0.14	35,48,48,52	0
2	EDO	D	509	4/4	0.88	0.20	37,42,47,53	0
2	EDO	A	511	4/4	0.88	0.14	44,46,52,52	0
2	EDO	C	510	4/4	0.88	0.15	37,41,50,58	0
2	EDO	D	514	4/4	0.89	0.26	33,45,51,57	0
2	EDO	B	511	4/4	0.89	0.24	30,38,41,42	0
2	EDO	C	514	4/4	0.90	0.20	32,43,50,52	0
2	EDO	D	518	4/4	0.90	0.16	31,38,39,46	0
2	EDO	C	505	4/4	0.90	0.09	21,24,27,28	0
2	EDO	C	511	4/4	0.90	0.09	28,33,34,40	0
2	EDO	B	506	4/4	0.91	0.14	31,38,53,55	0
2	EDO	D	515	4/4	0.91	0.11	18,27,31,31	0
2	EDO	A	512	4/4	0.91	0.17	30,37,46,54	0
2	EDO	B	516	4/4	0.91	0.20	33,44,50,50	0
2	EDO	B	503	4/4	0.91	0.12	17,29,32,35	0
2	EDO	D	513	4/4	0.91	0.19	33,34,38,48	0
2	EDO	A	514	4/4	0.93	0.17	28,35,39,49	0
2	EDO	D	504	4/4	0.93	0.09	24,26,27,32	0
2	EDO	D	507	4/4	0.93	0.17	26,33,36,39	0
2	EDO	D	516	4/4	0.93	0.17	18,24,41,44	0
2	EDO	A	505	4/4	0.93	0.08	17,23,24,27	0
2	EDO	B	505	4/4	0.93	0.08	22,24,27,29	0
2	EDO	C	503	4/4	0.93	0.10	32,38,41,44	0
2	EDO	B	509	4/4	0.93	0.07	32,39,41,44	0
2	EDO	A	503	4/4	0.94	0.09	23,23,25,32	0
2	EDO	B	502	4/4	0.94	0.09	21,26,28,28	0
3	NA	D	522	1/1	0.94	0.12	28,28,28,28	0
2	EDO	C	509	4/4	0.95	0.12	29,35,40,49	0
2	EDO	C	502	4/4	0.95	0.08	23,26,28,33	0
2	EDO	B	514	4/4	0.95	0.16	19,31,40,44	0
2	EDO	C	507	4/4	0.95	0.16	18,29,45,46	0
2	EDO	A	509	4/4	0.95	0.10	24,34,37,38	0
2	EDO	D	505	4/4	0.95	0.08	22,25,26,28	0
2	EDO	D	506	4/4	0.95	0.15	19,24,33,42	0
2	EDO	B	504	4/4	0.96	0.06	26,28,34,37	0
2	EDO	D	511	4/4	0.96	0.09	26,27,28,40	0
2	EDO	A	501	4/4	0.96	0.07	23,24,27,29	0
2	EDO	C	501	4/4	0.96	0.06	27,29,34,40	0
2	EDO	A	510	4/4	0.96	0.09	16,25,26,31	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	EDO	A	504	4/4	0.96	0.09	25,29,33,42	0
3	NA	D	523	1/1	0.96	0.18	32,32,32,32	0
2	EDO	D	519	4/4	0.97	0.10	25,29,39,44	0

6.5 Other polymers [i](#)

There are no such residues in this entry.