



Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 08:01 am BST

PDB ID : 6PNU
Title : Crystal structure of native DauA
Authors : Pluvinae, B.; Boraston, A.B.
Deposited on : 2019-07-03
Resolution : 2.00 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

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

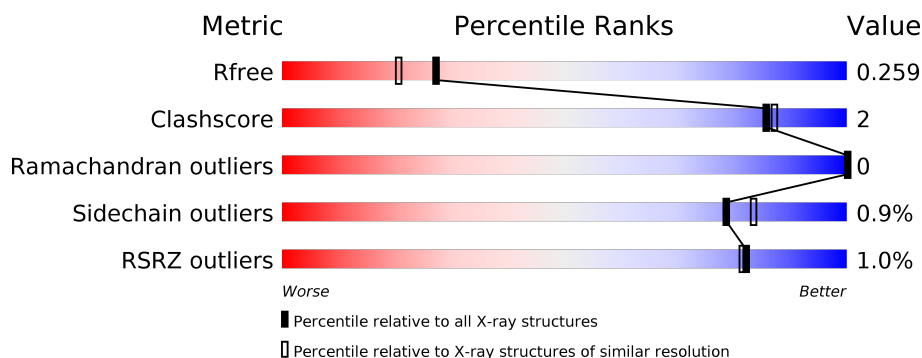
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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	497	<div> <div>%</div> <div> <div></div> <div>89%</div> <div>6%</div> <div>.</div> </div> </div>
1	B	497	<div> <div>2%</div> <div> <div></div> <div>92%</div> <div>5%</div> <div>.</div> </div> </div>
1	C	497	<div> <div>%</div> <div> <div></div> <div>92%</div> <div>5%</div> <div>.</div> </div> </div>
1	D	497	<div> <div>%</div> <div> <div></div> <div>90%</div> <div>6%</div> <div>.</div> </div> </div>

2 Entry composition [i](#)

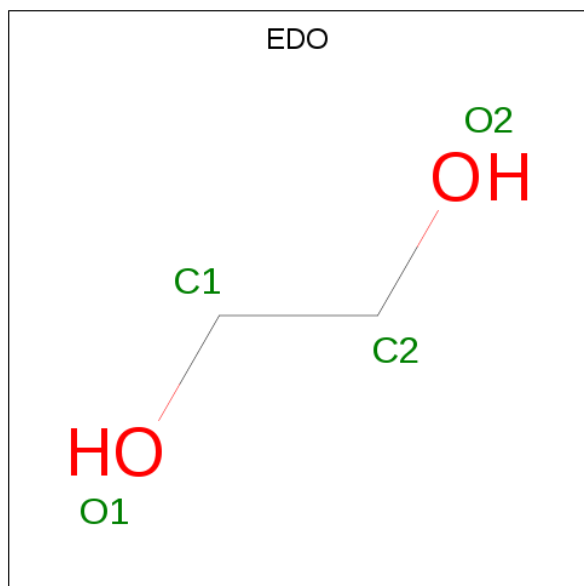
There are 3 unique types of molecules in this entry. The entry contains 15225 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Aldehyde dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	475	Total	C	N	O	S	0	1	0
			3561	2250	595	699	17			
1	B	482	Total	C	N	O	S	0	2	0
			3604	2278	601	708	17			
1	C	483	Total	C	N	O	S	0	0	0
			3593	2272	602	701	18			
1	D	482	Total	C	N	O	S	0	3	0
			3616	2284	604	710	18			

- 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		

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	159	Total	O	0	0
			159	159		

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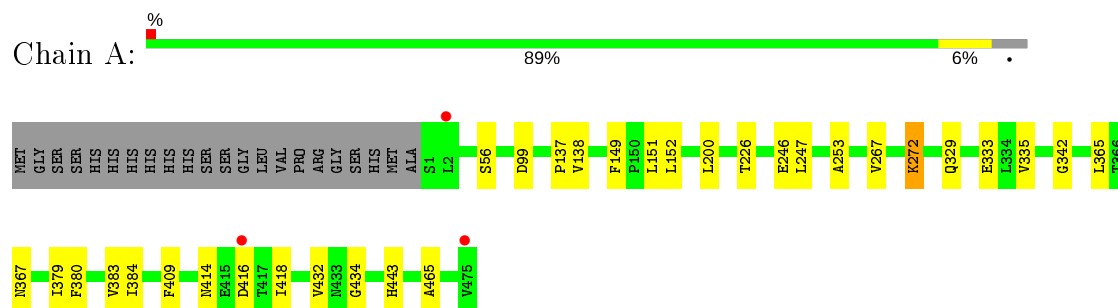
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	193	Total 193	O 193	0	0
3	C	209	Total 209	O 209	0	0
3	D	210	Total 210	O 210	0	0

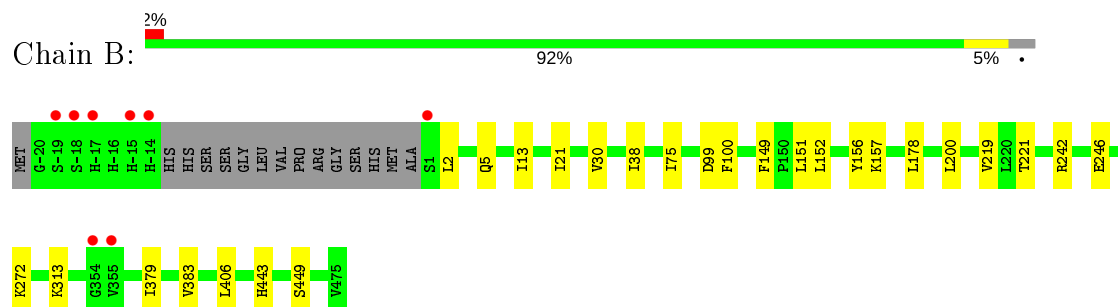
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

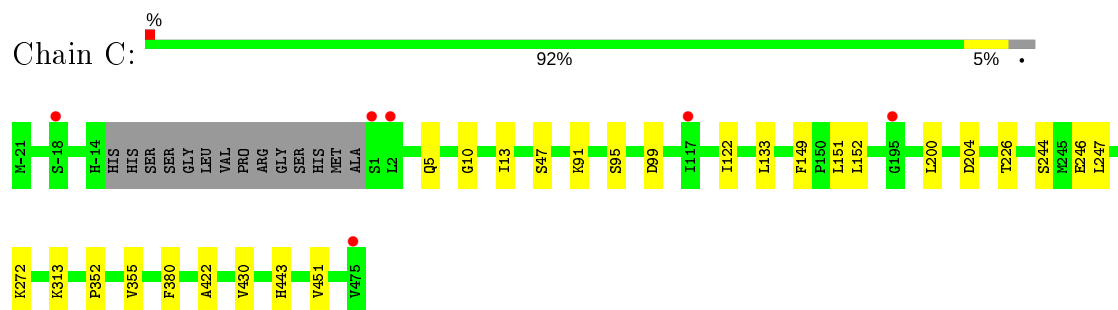
• Molecule 1: Aldehyde dehydrogenase



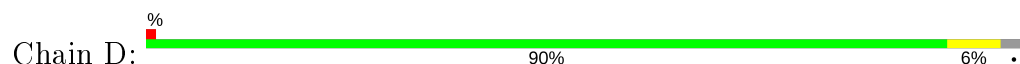
• Molecule 1: Aldehyde dehydrogenase

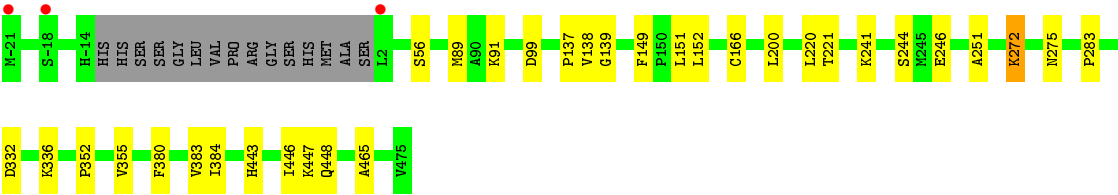


• Molecule 1: Aldehyde dehydrogenase



• Molecule 1: Aldehyde dehydrogenase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	82.88Å 137.80Å 108.87Å 90.00° 109.94° 90.00°	Depositor
Resolution (Å)	75.66 – 2.00 75.66 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.3 (75.66-2.00) 99.3 (75.66-2.00)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.28 (at 2.00Å)	Xtriage
Refinement program	PHENIX 1.15.2_3472	Depositor
R, R_{free}	0.222 , 0.259 0.222 , 0.259	Depositor DCC
R_{free} test set	7532 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å ²)	27.8	Xtriage
Anisotropy	0.508	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 43.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.000 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	15225	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 46.28 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.1705e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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

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	0.26	0/3625	0.42	0/4932
1	B	0.27	0/3667	0.43	0/4988
1	C	0.26	0/3656	0.43	0/4971
1	D	0.26	0/3679	0.43	0/5003
All	All	0.26	0/14627	0.43	0/19894

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	3561	0	3509	15	0
1	B	3604	0	3546	16	0
1	C	3593	0	3541	12	0
1	D	3616	0	3551	20	0
2	A	12	0	18	1	0
2	B	24	0	36	3	0
2	C	20	0	30	0	0
2	D	24	0	36	2	0
3	A	159	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	193	0	0	0	0
3	C	209	0	0	2	0
3	D	210	0	0	2	0
All	All	15225	0	14267	63	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 2.

All (63) 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:38:ILE:H	2:B:502:EDO:H12	1.52	0.74
1:D:99:ASP:HB3	1:D:151:LEU:HB2	1.74	0.70
1:B:99:ASP:HB3	1:B:151:LEU:HB2	1.72	0.69
1:C:99:ASP:HB3	1:C:151:LEU:HB2	1.77	0.67
1:A:149:PHE:HB3	1:A:152:LEU:HB3	1.77	0.66
1:D:149:PHE:HB3	1:D:152:LEU:HB3	1.80	0.63
1:B:149:PHE:HB3	1:B:152:LEU:HB3	1.82	0.61
1:C:149:PHE:HB3	1:C:152:LEU:HB3	1.80	0.61
1:B:21:ILE:HG13	1:B:178:LEU:HD21	1.84	0.60
1:A:99:ASP:HB3	1:A:151:LEU:HB2	1.84	0.59
1:C:246:GLU:OE2	1:C:443:HIS:NE2	2.35	0.59
1:D:89[A]:MET:SD	1:D:91:LYS:HE3	2.45	0.57
1:A:226:THR:HA	1:A:247:LEU:HD13	1.86	0.56
1:A:246:GLU:OE1	1:A:443:HIS:NE2	2.37	0.56
1:A:329:GLN:HE21	1:A:333:GLU:HG3	1.70	0.56
1:A:138:VAL:HG12	1:A:465:ALA:HB2	1.89	0.54
1:A:267:VAL:HG21	1:A:434:GLY:HA2	1.88	0.53
1:B:246:GLU:OE1	1:B:443:HIS:NE2	2.41	0.52
1:C:226:THR:HA	1:C:247:LEU:HD13	1.91	0.51
1:D:91:LYS:HE2	1:D:275:ASN:OD1	2.11	0.50
1:D:272:LYS:HE3	1:D:383:VAL:O	2.11	0.50
1:B:5:GLN:HB3	1:B:13:ILE:O	2.11	0.50
1:D:138:VAL:HG12	1:D:465:ALA:HB2	1.94	0.49
1:B:219:VAL:HG22	1:B:242:ARG:HB2	1.94	0.49
1:A:335:VAL:HG13	1:A:365:LEU:HD11	1.95	0.49
1:B:379:ILE:HG21	1:B:383:VAL:HB	1.95	0.49
1:A:253:ALA:HA	1:A:409:PHE:HB2	1.94	0.48
1:B:21:ILE:HG13	1:B:178:LEU:CD2	2.43	0.48
1:C:91:LYS:HG2	1:C:95:SER:HB2	1.95	0.48
1:B:406:LEU:HG	1:B:449:SER:HB3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:313:LYS:NZ	3:C:602:HOH:O	2.39	0.47
1:D:251:ALA:HB3	1:D:283:PRO:HA	1.96	0.47
1:D:272:LYS:NZ	3:D:606:HOH:O	2.42	0.47
1:D:272:LYS:HG3	1:D:384:ILE:HG12	1.96	0.46
1:A:414:ASN:O	1:A:418:ILE:HG13	2.16	0.46
1:A:56:SER:HB2	1:A:137:PRO:HG2	1.99	0.45
1:B:38:ILE:HG13	2:B:502:EDO:H21	1.99	0.45
1:A:379:ILE:HG21	1:A:383:VAL:HB	1.98	0.45
1:D:220:LEU:HG	1:D:241:LYS:HD3	1.99	0.45
1:D:352:PRO:HG2	1:D:355:VAL:HG21	1.99	0.44
1:C:244:SER:HA	1:C:451:VAL:HG23	1.99	0.44
1:C:5:GLN:HB3	1:C:13:ILE:O	2.17	0.44
1:D:447:LYS:NZ	3:D:604:HOH:O	2.41	0.44
1:D:221:THR:HA	1:D:244:SER:O	2.18	0.44
1:D:272:LYS:HE2	1:D:283:PRO:O	2.18	0.43
2:A:503:EDO:H11	1:B:313:LYS:HG3	2.00	0.43
1:C:122:ILE:HD11	1:C:133:LEU:HD22	1.99	0.43
1:D:446:ILE:HB	2:D:501:EDO:H11	2.01	0.43
1:D:246:GLU:OE2	1:D:443:HIS:NE2	2.47	0.43
1:C:352:PRO:HG2	1:C:355:VAL:HG11	2.01	0.42
1:B:156:TYR:HE2	2:B:506:EDO:H12	1.84	0.42
1:D:332:ASP:OD2	1:D:336:LYS:HE2	2.20	0.42
1:B:75[B]:ILE:HD13	1:B:100:PHE:HZ	1.84	0.42
1:B:2:LEU:HD12	1:B:30:VAL:O	2.20	0.41
1:C:10:GLY:N	3:C:613:HOH:O	2.49	0.41
1:D:139:GLY:O	1:D:166:CYS:HB3	2.20	0.41
1:D:448:GLN:HB2	2:D:502:EDO:H22	2.01	0.41
1:C:422:ALA:HA	1:C:430:VAL:HG11	2.02	0.41
1:D:56:SER:HB2	1:D:137:PRO:HG2	2.02	0.41
1:A:272:LYS:HD3	1:A:384:ILE:HG12	2.03	0.40
1:A:418:ILE:HD13	1:A:432:VAL:HG11	2.02	0.40
1:A:342:GLY:O	1:A:367:ASN:ND2	2.40	0.40
1:B:157:LYS:HD2	1:B:221:THR:HG21	2.03	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	474/497 (95%)	458 (97%)	16 (3%)	0	100	100
1	B	480/497 (97%)	465 (97%)	15 (3%)	0	100	100
1	C	479/497 (96%)	461 (96%)	18 (4%)	0	100	100
1	D	481/497 (97%)	464 (96%)	17 (4%)	0	100	100
All	All	1914/1988 (96%)	1848 (97%)	66 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	382/407 (94%)	378 (99%)	4 (1%)	76	81
1	B	385/407 (95%)	383 (100%)	2 (0%)	88	92
1	C	382/407 (94%)	377 (99%)	5 (1%)	69	74
1	D	386/407 (95%)	383 (99%)	3 (1%)	81	86
All	All	1535/1628 (94%)	1521 (99%)	14 (1%)	78	83

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	200	LEU
1	A	272	LYS

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Mol	Chain	Res	Type
1	A	380	PHE
1	A	416	ASP
1	B	200	LEU
1	B	272	LYS
1	C	47	SER
1	C	200	LEU
1	C	204	ASP
1	C	272	LYS
1	C	380	PHE
1	D	200	LEU
1	D	272	LYS
1	D	380	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	ASN
1	A	329	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

20 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond

length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	EDO	C	505	-	3,3,3	0.47	0	2,2,2	0.28	0
2	EDO	B	506	-	3,3,3	0.49	0	2,2,2	0.28	0
2	EDO	B	503	-	3,3,3	0.50	0	2,2,2	0.22	0
2	EDO	C	501	-	3,3,3	0.47	0	2,2,2	0.32	0
2	EDO	B	502	-	3,3,3	0.48	0	2,2,2	0.32	0
2	EDO	D	502	-	3,3,3	0.49	0	2,2,2	0.27	0
2	EDO	A	502	-	3,3,3	0.46	0	2,2,2	0.35	0
2	EDO	C	502	-	3,3,3	0.46	0	2,2,2	0.31	0
2	EDO	D	506	-	3,3,3	0.51	0	2,2,2	0.19	0
2	EDO	B	505	-	3,3,3	0.44	0	2,2,2	0.36	0
2	EDO	C	503	-	3,3,3	0.51	0	2,2,2	0.15	0
2	EDO	D	501	-	3,3,3	0.43	0	2,2,2	0.28	0
2	EDO	B	501	-	3,3,3	0.47	0	2,2,2	0.36	0
2	EDO	D	505	-	3,3,3	0.48	0	2,2,2	0.26	0
2	EDO	A	503	-	3,3,3	0.50	0	2,2,2	0.13	0
2	EDO	C	504	-	3,3,3	0.47	0	2,2,2	0.18	0
2	EDO	D	504	-	3,3,3	0.45	0	2,2,2	0.33	0
2	EDO	B	504	-	3,3,3	0.45	0	2,2,2	0.37	0
2	EDO	D	503	-	3,3,3	0.48	0	2,2,2	0.28	0
2	EDO	A	501	-	3,3,3	0.47	0	2,2,2	0.31	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	C	505	-	-	0/1/1/1	-
2	EDO	B	506	-	-	0/1/1/1	-
2	EDO	B	503	-	-	0/1/1/1	-
2	EDO	C	501	-	-	0/1/1/1	-
2	EDO	B	502	-	-	0/1/1/1	-
2	EDO	D	502	-	-	0/1/1/1	-
2	EDO	A	502	-	-	0/1/1/1	-
2	EDO	C	502	-	-	0/1/1/1	-
2	EDO	D	506	-	-	0/1/1/1	-
2	EDO	B	505	-	-	0/1/1/1	-
2	EDO	C	503	-	-	0/1/1/1	-
2	EDO	D	501	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	EDO	B	501	-	-	0/1/1/1	-
2	EDO	D	505	-	-	0/1/1/1	-
2	EDO	A	503	-	-	0/1/1/1	-
2	EDO	C	504	-	-	0/1/1/1	-
2	EDO	D	504	-	-	0/1/1/1	-
2	EDO	B	504	-	-	0/1/1/1	-
2	EDO	D	503	-	-	1/1/1/1	-
2	EDO	A	501	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	503	EDO	O1-C1-C2-O2
2	A	501	EDO	O1-C1-C2-O2

There are no ring outliers.

5 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	506	EDO	1	0
2	B	502	EDO	2	0
2	D	502	EDO	1	0
2	D	501	EDO	1	0
2	A	503	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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	475/497 (95%)	0.02	3 (0%) 89 88	17, 33, 52, 63	0
1	B	482/497 (96%)	-0.10	8 (1%) 70 68	17, 28, 43, 56	0
1	C	483/497 (97%)	-0.05	6 (1%) 79 78	18, 28, 44, 61	0
1	D	482/497 (96%)	-0.05	3 (0%) 89 88	19, 27, 41, 51	0
All	All	1922/1988 (96%)	-0.05	20 (1%) 82 81	17, 29, 47, 63	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1	SER	4.8
1	B	-18	SER	4.4
1	C	2	LEU	3.6
1	C	1	SER	3.6
1	D	-18	SER	3.5
1	A	2	LEU	3.5
1	A	475	VAL	3.1
1	B	-15	HIS	2.9
1	A	416	ASP	2.9
1	C	475	VAL	2.7
1	B	-17	HIS	2.7
1	B	-14	HIS	2.6
1	C	117	ILE	2.5
1	B	354	GLY	2.5
1	D	2	LEU	2.3
1	D	-21	MET	2.3
1	B	355	VAL	2.2
1	C	-18	SER	2.1
1	C	195	GLY	2.0
1	B	-19	SER	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	EDO	B	506	4/4	0.67	0.19	41,42,43,45	0
2	EDO	A	503	4/4	0.70	0.29	35,37,41,44	0
2	EDO	C	501	4/4	0.75	0.14	34,39,41,46	0
2	EDO	D	506	4/4	0.76	0.20	34,37,39,46	0
2	EDO	D	503	4/4	0.78	0.17	33,37,37,41	0
2	EDO	D	505	4/4	0.80	0.13	32,40,47,48	0
2	EDO	B	503	4/4	0.84	0.23	33,36,37,39	0
2	EDO	B	502	4/4	0.85	0.13	37,38,39,44	0
2	EDO	A	502	4/4	0.87	0.14	40,44,47,52	0
2	EDO	D	502	4/4	0.87	0.12	28,31,38,42	0
2	EDO	D	501	4/4	0.87	0.25	24,27,28,35	0
2	EDO	B	501	4/4	0.88	0.17	33,35,37,43	0
2	EDO	C	504	4/4	0.89	0.12	34,38,48,48	0
2	EDO	C	505	4/4	0.90	0.17	31,36,41,43	0
2	EDO	D	504	4/4	0.91	0.15	44,45,46,51	0
2	EDO	C	503	4/4	0.91	0.17	31,31,32,35	0
2	EDO	A	501	4/4	0.91	0.12	38,38,45,50	0
2	EDO	C	502	4/4	0.94	0.11	38,39,39,52	0
2	EDO	B	505	4/4	0.95	0.10	30,34,38,44	0
2	EDO	B	504	4/4	0.95	0.16	32,40,40,41	0

6.5 Other polymers [i](#)

There are no such residues in this entry.