



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 06:54 AM UTC

PDB ID : 6B4R / pdb_00006b4r
Title : The crystal structure of the aldehyde dehydrogenase KauB from *Pseudomonas aeruginosa*
Authors : Gonzalez-Segura, L.; Cardona-Cardona, Y.; Carrillo-Campos, J.; Munoz-Clares, R.A.
Deposited on : 2017-09-27
Resolution : 2.55 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 : 4-5-2 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

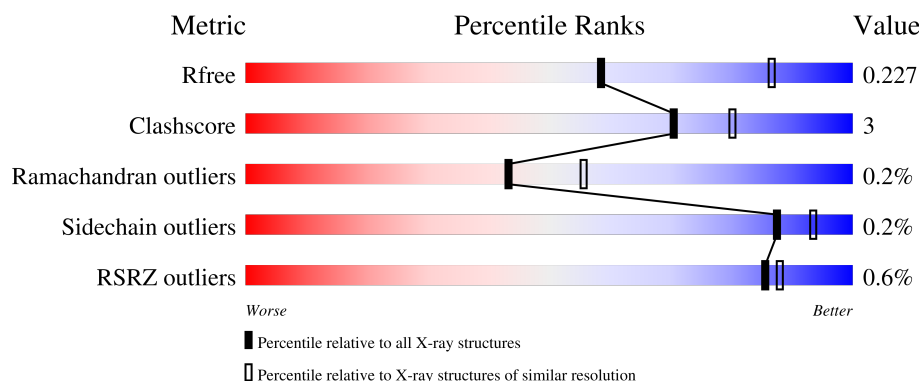
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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}	180053	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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	497	91% 9%
1	C	497	90% 9%
2	B	497	92% 8%
2	D	497	2% 90% 10%

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
1	CSD	C	302	-	-	X	-
5	TOE	C	503	-	-	X	-

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 15533 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.

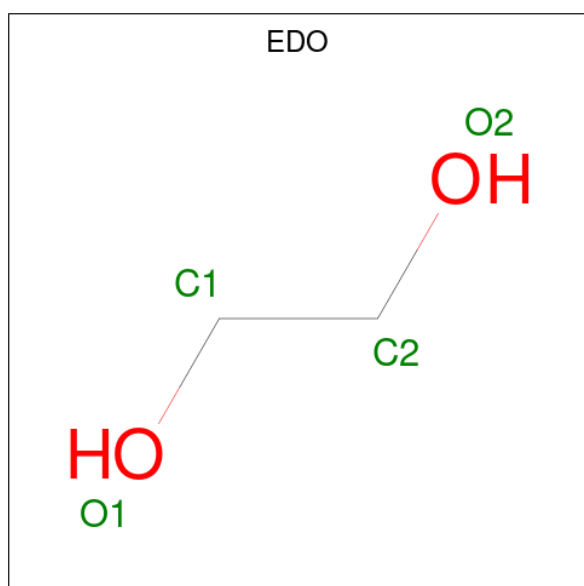
- Molecule 1 is a protein called KauB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	497	Total	C	N	O	S	0	1	0
			3752	2380	640	718	14			
1	C	495	Total	C	N	O	S	0	1	0
			3737	2371	638	715	13			

- Molecule 2 is a protein called KauB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	495	Total	C	N	O	S	0	1	0
			3735	2371	638	713	13			
2	D	495	Total	C	N	O	S	0	1	0
			3735	2371	638	713	13			

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).

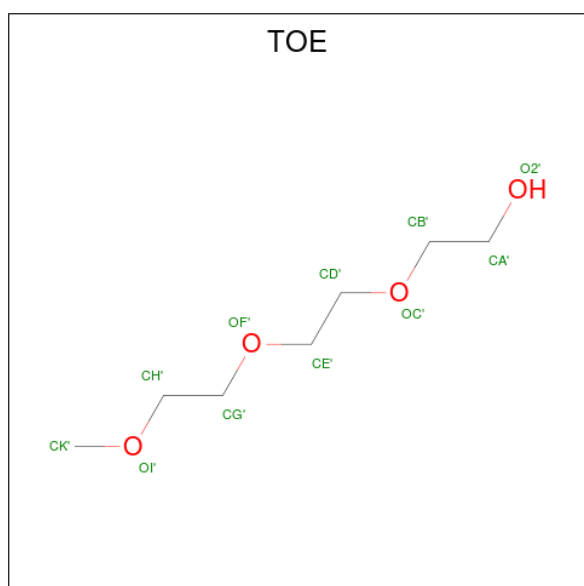


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	C	1	Total C O 4 2 2	0	0

- Molecule 4 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total K 1 1	0	0
4	B	1	Total K 1 1	0	0
4	C	1	Total K 1 1	0	0
4	D	1	Total K 1 1	0	0

- Molecule 5 is 2-[2-(2-METHOXY-ETHOXY)-ETHOXY]-ETHOXYL (CCD ID: TOE) (formula: C₇H₁₆O₄).



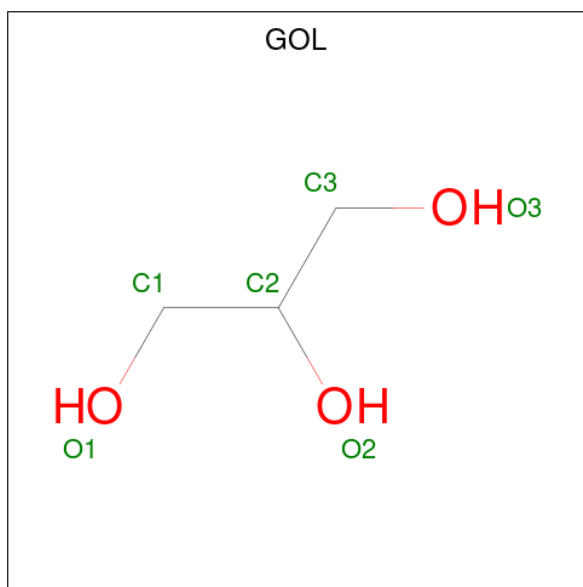
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 11 7 4	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			11	7	4		
5	C	1	Total	C	O	0	0
			10	6	4		

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



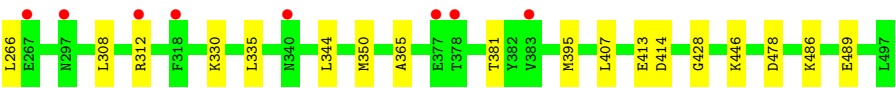
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	170	Total	O	0	0
			170	170		
7	B	146	Total	O	0	0
			146	146		
7	C	123	Total	O	0	0
			123	123		
7	D	77	Total	O	0	0
			77	77		

- Molecule 1: KauB





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	177.64Å 93.81Å 128.86Å 90.00° 92.28° 90.00°	Depositor
Resolution (Å)	46.12 – 2.55 46.12 – 2.55	Depositor EDS
% Data completeness (in resolution range)	99.4 (46.12-2.55) 99.4 (46.12-2.55)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.82 (at 2.54Å)	Xtriage
Refinement program	PHENIX 1.8.1_1168	Depositor
R, R_{free}	0.178 , 0.234 (Not available) , 0.227	Depositor DCC
R_{free} test set	3486 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	34.0	Xtriage
Anisotropy	0.177	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 39.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.020 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	15533	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.29% 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: K, CSD, EDO, TOE, GOL

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	0.29	0/3818	0.69	0/5186
1	C	0.28	0/3803	0.70	0/5166
2	B	0.28	0/3809	0.69	1/5174 (0.0%)
2	D	0.28	0/3809	0.69	0/5174
All	All	0.28	0/15239	0.69	1/20700 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	198	LEU	N-CA-C	5.23	116.66	111.07

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	3752	0	3760	27	0
1	C	3737	0	3741	33	0
2	B	3735	0	3741	24	0
2	D	3735	0	3741	28	0
3	A	8	0	12	2	0
3	B	4	0	6	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	4	0	6	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	11	0	16	0	0
5	B	11	0	16	0	0
5	C	10	0	13	8	0
6	B	6	0	8	0	0
7	A	170	0	0	0	1
7	B	146	0	0	1	0
7	C	123	0	0	0	0
7	D	77	0	0	0	0
All	All	15533	0	15060	101	1

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 3.

All (101) 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:C:302:CSD:CB	5:C:503:TOE:H3	1.88	1.04
1:C:302:CSD:HB3	5:C:503:TOE:H3	1.33	1.03
1:C:302:CSD:SG	5:C:503:TOE:H3	2.02	0.99
1:C:302:CSD:HB3	5:C:503:TOE:CA'	2.01	0.89
1:C:302:CSD:SG	5:C:503:TOE:CA'	2.80	0.70
2:D:350:MET:HE3	2:D:381:THR:HG23	1.75	0.69
2:D:312:ARG:HH12	2:D:413:GLU:HG3	1.58	0.68
2:D:365:ALA:HB2	2:D:395:MET:HE1	1.77	0.65
2:D:36:THR:HG21	2:D:50:LYS:HB3	1.80	0.64
1:A:142:VAL:HB	1:C:140:ASP:HB2	1.80	0.64
2:B:249:ALA:HB1	2:B:266:LEU:HD13	1.83	0.60
1:C:254:VAL:HG13	2:D:254:VAL:HG13	1.82	0.60
2:B:166:PRO:HD3	2:B:243:THR:HB	1.83	0.59
2:D:158:VAL:HG12	2:D:489:GLU:HG2	1.83	0.58
1:A:385:PRO:HA	1:A:405:PRO:HB2	1.86	0.58
1:C:242:PHE:HB3	1:C:266:LEU:HD23	1.86	0.57
1:C:302:CSD:CB	5:C:503:TOE:CA'	2.71	0.57
1:A:263:ARG:NH2	2:B:474:GLY:O	2.38	0.57
1:C:474:GLY:O	2:D:263:ARG:NH2	2.37	0.57
1:A:105:GLU:HG2	1:A:199:THR:HG21	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:140:ASP:HB2	1:C:142:VAL:HB	1.88	0.55
2:D:111:LYS:HD2	2:D:112:PRO:HD2	1.88	0.55
1:A:166:PRO:HD3	1:A:243:THR:HB	1.88	0.55
1:A:249:ALA:HB1	1:A:266:LEU:HD13	1.87	0.55
1:A:446:LYS:NZ	2:D:147:HIS:O	2.36	0.54
2:B:158:VAL:HG12	2:B:489:GLU:HG2	1.88	0.54
2:D:184:THR:OG1	2:D:186:ASN:ND2	2.40	0.54
1:C:158:VAL:HG12	1:C:489:GLU:HG2	1.89	0.54
2:B:140:ASP:HB2	2:D:142:VAL:HB	1.90	0.54
1:A:481:LEU:HD12	3:A:502:EDO:H22	1.88	0.53
2:B:315:LYS:HD2	2:B:411:GLU:HB3	1.89	0.53
2:B:169:PHE:HB2	2:B:173:MET:HG2	1.90	0.52
1:A:167:TRP:CD1	1:A:195:LYS:HD2	2.45	0.52
2:D:135:ILE:HG12	2:D:183:ALA:HB1	1.93	0.51
2:D:478:ASP:OD1	2:D:486:LYS:NZ	2.43	0.51
1:C:278:ALA:HA	1:C:314:ILE:HD13	1.93	0.51
2:B:100:GLU:H	3:B:501:EDO:H11	1.76	0.50
2:B:393:ASN:O	2:B:399:ARG:NH2	2.41	0.50
1:A:308:LEU:HB2	1:A:407:LEU:HD11	1.92	0.50
2:B:135:ILE:HG12	2:B:183:ALA:HB1	1.94	0.50
2:D:9:TRP:HB3	2:D:335:LEU:HD12	1.94	0.50
1:A:481:LEU:HB2	3:A:502:EDO:H22	1.93	0.50
1:C:135:ILE:HG12	1:C:183:ALA:HB1	1.93	0.50
2:B:317:LYS:NZ	7:B:602:HOH:O	2.45	0.49
2:D:55:ASP:CG	2:D:56:LEU:H	2.20	0.49
1:A:250:LYS:HG3	2:B:261:MET:HE2	1.95	0.49
2:D:242:PHE:HB3	2:D:266:LEU:HD23	1.94	0.49
2:B:141:GLU:HB2	2:B:153:VAL:HB	1.95	0.49
2:D:344:LEU:HB2	2:D:381:THR:HG22	1.95	0.49
1:A:240:LEU:HD11	1:A:252:LEU:HD13	1.95	0.49
1:C:468:GLY:HA3	1:C:477:ARG:HD3	1.95	0.48
2:D:308:LEU:HB2	2:D:407:LEU:HD11	1.95	0.48
2:B:241:VAL:HG12	2:B:265:TRP:HB2	1.94	0.48
1:A:104:LEU:O	1:A:108:ASP:HB2	2.13	0.48
2:B:301:VAL:HG12	2:B:303:THR:H	1.78	0.48
1:A:278:ALA:HA	1:A:314:ILE:HD13	1.94	0.48
1:A:169:PHE:HB2	1:A:173:MET:HG2	1.95	0.48
1:C:344:LEU:HD11	1:C:383:VAL:HG12	1.96	0.48
1:C:169:PHE:HB2	1:C:173:MET:HG2	1.96	0.47
1:C:41:SER:HB3	1:C:46:ARG:HG2	1.95	0.47
2:D:141:GLU:HB2	2:D:153:VAL:HB	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:147:HIS:O	2:D:446:LYS:NZ	2.46	0.46
1:C:302:CSD:OD1	5:C:503:TOE:CA'	2.63	0.46
2:B:142:VAL:HB	2:D:140:ASP:HB2	1.97	0.46
1:C:41:SER:HB2	1:C:48:LEU:HD21	1.96	0.46
1:C:164:ILE:HB	1:C:242:PHE:HD1	1.81	0.46
1:C:166:PRO:HD3	1:C:243:THR:HB	1.98	0.45
1:C:168:ASN:ND2	5:C:503:TOE:O2'	2.49	0.45
1:A:474:GLY:O	2:B:263:ARG:NH2	2.48	0.45
1:A:242:PHE:HB3	1:A:266:LEU:HD23	1.99	0.45
1:C:385:PRO:HA	1:C:405:PRO:HB2	1.98	0.45
1:A:478:ASP:OD1	1:A:486:LYS:NZ	2.48	0.44
1:C:478:ASP:OD1	1:C:486:LYS:NZ	2.42	0.44
2:D:312:ARG:NH1	2:D:413:GLU:HG3	2.29	0.44
2:B:278:ALA:HA	2:B:314:ILE:HD13	1.99	0.44
1:A:195:LYS:NZ	1:A:346:ASP:OD1	2.36	0.44
1:A:135:ILE:HG12	1:A:183:ALA:HB1	1.98	0.44
2:D:41:SER:HB3	2:D:46:ARG:HG2	1.99	0.44
1:C:93:LEU:HD12	1:C:210:ALA:HB2	2.00	0.44
2:B:314:ILE:HA	2:B:317:LYS:HE2	1.99	0.43
2:B:167:TRP:CD1	2:B:195:LYS:HD2	2.53	0.43
1:A:6:ARG:HD2	1:A:333:ASN:HD22	1.83	0.43
2:B:308:LEU:HB2	2:B:407:LEU:HD11	1.99	0.43
2:D:414:ASP:N	2:D:414:ASP:OD1	2.52	0.42
1:A:111:LYS:NZ	1:A:120:ASP:OD1	2.52	0.42
2:D:169:PHE:HB2	2:D:173:MET:HG2	2.02	0.42
2:D:104:LEU:O	2:D:108:ASP:HB2	2.20	0.41
1:C:141:GLU:HB2	1:C:153:VAL:HB	2.01	0.41
2:B:93:LEU:HD12	2:B:210:ALA:HB2	2.03	0.41
2:B:247:LYS:HE2	2:B:247:LYS:HB3	1.85	0.41
2:D:169:PHE:HB3	2:D:172:LEU:HB3	2.02	0.41
1:A:93:LEU:HD12	1:A:210:ALA:HB2	2.03	0.41
1:C:63:VAL:HG11	1:C:234:HIS:CE1	2.55	0.41
1:C:465:ALA:HA	1:C:466:PRO:HD2	1.95	0.41
2:D:67:ARG:NH2	2:D:238:ASP:OD2	2.49	0.41
1:A:412:PHE:CD2	1:A:418:ALA:HB2	2.55	0.41
1:C:347:THR:HA	1:C:381:THR:HG21	2.01	0.41
2:B:96:GLU:O	3:B:501:EDO:H22	2.21	0.41
1:C:315:LYS:HG3	1:C:411:GLU:HG2	2.02	0.41
1:C:109:MET:SD	1:C:111:LYS:HE3	2.61	0.40
1:C:280:ALA:HA	1:C:281:PRO:HD3	1.97	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:A:601:HOH:O	7:A:601:HOH:O[2_756]	2.17	0.03

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	495/497 (100%)	474 (96%)	20 (4%)	1 (0%)	43	56
1	C	493/497 (99%)	473 (96%)	19 (4%)	1 (0%)	43	56
2	B	494/497 (99%)	475 (96%)	18 (4%)	1 (0%)	43	56
2	D	494/497 (99%)	474 (96%)	19 (4%)	1 (0%)	43	56
All	All	1976/1988 (99%)	1896 (96%)	76 (4%)	4 (0%)	43	56

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	428	GLY
1	A	428	GLY
2	B	428	GLY
1	C	428	GLY

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	385/384 (100%)	385 (100%)	0	100	100
1	C	383/384 (100%)	383 (100%)	0	100	100
2	B	384/385 (100%)	384 (100%)	0	100	100
2	D	384/385 (100%)	381 (99%)	3 (1%)	73	84
All	All	1536/1538 (100%)	1533 (100%)	3 (0%)	87	93

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	D	5	THR
2	D	105	GLU
2	D	330	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	251	GLN
2	B	472	GLN
1	C	348	GLN
1	C	349	GLN
1	C	362	ASN
2	D	225	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 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	CSD	A	302	1	4,7,8	1.07	0	1,8,10	0.31	0
1	CSD	C	302	1	4,7,8	1.02	0	1,8,10	0.21	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
1	CSD	A	302	1	-	0/2/6/8	-
1	CSD	C	302	1	-	0/2/6/8	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	C	302	CSD	7	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 4 are monoatomic - leaving 8 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
3	EDO	A	502	-	3,3,3	0.44	0	2,2,2	0.31	0
3	EDO	B	501	-	3,3,3	0.44	0	2,2,2	0.34	0
3	EDO	C	501	-	3,3,3	0.44	0	2,2,2	0.31	0
5	TOE	B	504	-	10,10,10	0.45	0	9,9,9	0.46	0
3	EDO	A	501	-	3,3,3	0.43	0	2,2,2	0.41	0
6	GOL	B	503	-	5,5,5	0.38	0	5,5,5	0.20	0
5	TOE	C	503	-	9,9,10	0.51	0	8,8,9	0.58	0
5	TOE	A	504	-	10,10,10	0.46	0	9,9,9	0.29	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	502	-	-	0/1/1/1	-
3	EDO	B	501	-	-	1/1/1/1	-
3	EDO	C	501	-	-	0/1/1/1	-
5	TOE	B	504	-	-	2/8/8/8	-
3	EDO	A	501	-	-	0/1/1/1	-
6	GOL	B	503	-	-	2/4/4/4	-
5	TOE	C	503	-	-	4/7/7/8	-
5	TOE	A	504	-	-	1/8/8/8	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	B	503	GOL	O1-C1-C2-C3
5	C	503	TOE	OC'-CD'-CE'-OF'
6	B	503	GOL	O1-C1-C2-O2
5	B	504	TOE	OF'-CG'-CH'-OI'
5	A	504	TOE	O2'-CA'-CB'-OC'
5	C	503	TOE	OF'-CG'-CH'-OI'
5	B	504	TOE	OC'-CD'-CE'-OF'
5	C	503	TOE	CE'-CD'-OC'-CB'
5	C	503	TOE	CA'-CB'-OC'-CD'
3	B	501	EDO	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	502	EDO	2	0
3	B	501	EDO	2	0
5	C	503	TOE	8	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	496/497 (99%)	-0.37	0 100 100	16, 29, 41, 57	1 (0%)
1	C	494/497 (99%)	0.05	1 (0%) 91 93	20, 37, 53, 75	1 (0%)
2	B	495/497 (99%)	-0.28	0 100 100	18, 30, 49, 60	1 (0%)
2	D	495/497 (99%)	0.37	11 (2%) 62 63	19, 45, 68, 84	1 (0%)
All	All	1980/1988 (99%)	-0.06	12 (0%) 85 88	16, 34, 58, 84	4 (0%)

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	3	THR	2.9
2	D	45	GLY	2.6
2	D	267[A]	GLU	2.5
2	D	312	ARG	2.3
2	D	111	LYS	2.3
2	D	378	THR	2.2
2	D	383	VAL	2.2
1	C	3	THR	2.2
2	D	318	PHE	2.2
2	D	377	GLU	2.1
2	D	297	ASN	2.1
2	D	340	ASN	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($\text{\AA}^2$)	Q<0.9
1	CSD	C	302	8/9	0.85	0.17	45,50,65,66	0
1	CSD	A	302	8/9	0.87	0.12	27,33,50,63	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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($\text{\AA}^2$)	Q<0.9
5	TOE	B	504	11/11	0.71	0.21	51,55,61,69	0
5	TOE	A	504	11/11	0.77	0.18	44,51,58,59	0
6	GOL	B	503	6/6	0.77	0.28	32,40,49,51	0
5	TOE	C	503	10/11	0.78	0.17	49,54,63,67	0
3	EDO	B	501	4/4	0.84	0.14	34,42,45,45	0
3	EDO	C	501	4/4	0.84	0.14	38,41,44,44	0
3	EDO	A	501	4/4	0.85	0.15	38,41,46,50	0
4	K	D	501	1/1	0.86	0.10	58,58,58,58	0
3	EDO	A	502	4/4	0.91	0.18	26,33,35,41	0
4	K	C	502	1/1	0.94	0.07	48,48,48,48	0
4	K	A	503	1/1	0.96	0.05	32,32,32,32	0
4	K	B	502	1/1	0.98	0.04	29,29,29,29	0

6.5 Other polymers [i](#)

There are no such residues in this entry.