



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 11:43 PM GMT

PDB ID : 1E3A  
Title : A SLOW PROCESSING PRECURSOR PENICILLIN ACYLASE FROM ES-  
CHERICHIA COLI  
Authors : Hewitt, L.; Kasche, V.; Lummer, K.; Lewis, R.J.; Murshudov, G.N.; Verma,  
C.S.; Dodson, G.G.; Wilson, K.S.  
Deposited on : 2000-06-07  
Resolution : 1.80 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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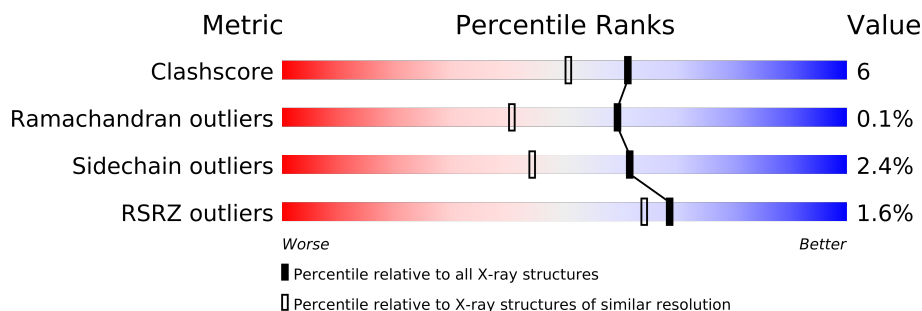
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance



The reported resolution of this entry is 1.80 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))
Clashscore	79885	4461 (1.80-1.80)
Ramachandran outliers	78287	4404 (1.80-1.80)
Sidechain outliers	78261	4403 (1.80-1.80)
RSRZ outliers	66119	3515 (1.80-1.80)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	260	
2	B	560	

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7742 atoms, of which 0 are hydrogen and 0 are deuterium.

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 PENICILLIN AMIDASE ALPHA SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	258	Total	C	N	O	S	0	4	0
			2036	1299	346	381	10			

- Molecule 2 is a protein called PENICILLIN AMIDASE BETA SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	560	Total	C	N	O	S	0	20	0
			4499	2854	777	858	10			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	263	GLY	THR	ENGINEERED MUTATION	UNP P06875

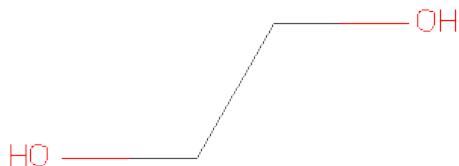
- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Ca	0	0
			1	1		

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Cl	0	0
			1	1		

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



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

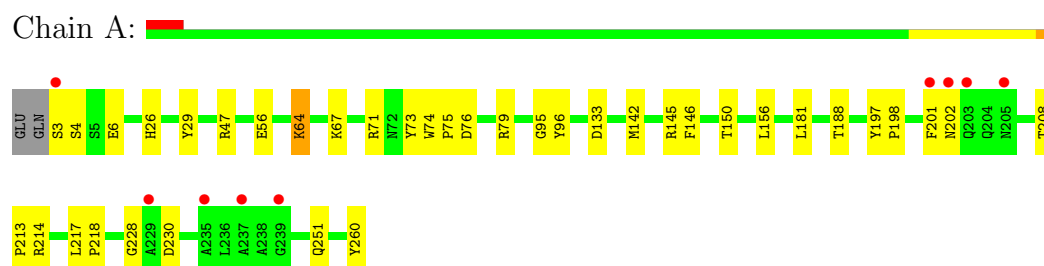
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	337	Total	O	0	0
			337	337		
6	B	856	Total	O	0	0
			856	856		

### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: PENICILLIN AMIDASE ALPHA SUBUNIT



#### • Molecule 2: PENICILLIN AMIDASE BETA SUBUNIT



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	50.71Å 64.27Å 72.00Å 66.14° 74.18° 74.23°	Depositor
Resolution (Å)	20.00 – 1.80 20.94 – 1.80	Depositor EDS
% Data completeness (in resolution range)	96.9 (20.00-1.80) 97.0 (20.94-1.80)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.23 (at 1.80Å)	Xtriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.158 , 0.197 0.155 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	10.9	Xtriage
Anisotropy	0.132	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 47.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 70641 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	7742	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.36% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 5 Model quality ⓘ

### 5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: CA, EDO, CL

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	0.49	0/2105	1.14	12/2856 (0.4%)
2	B	0.52	0/4716	1.23	37/6428 (0.6%)
All	All	0.51	0/6821	1.20	49/9284 (0.5%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
2	B	0	1
All	All	0	2

There are no bond length outliers.

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	734	ARG	NE-CZ-NH1	12.70	126.65	120.30
2	B	742	ARG	NE-CZ-NH2	-11.86	114.37	120.30
2	B	588[A]	ARG	NE-CZ-NH1	10.52	125.56	120.30
2	B	588[B]	ARG	NE-CZ-NH1	10.52	125.56	120.30
2	B	262	THR	CA-C-N	9.63	135.46	116.20
2	B	383	ARG	NE-CZ-NH1	9.04	124.82	120.30
1	A	71	ARG	CD-NE-CZ	8.24	135.14	123.60
2	B	383	ARG	NE-CZ-NH2	-8.19	116.20	120.30
1	A	145	ARG	NE-CZ-NH2	-8.18	116.21	120.30
2	B	404	ARG	NE-CZ-NH2	-7.93	116.34	120.30
2	B	700	ARG	NE-CZ-NH1	-7.84	116.38	120.30
2	B	588[A]	ARG	NE-CZ-NH2	-7.83	116.39	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	588[B]	ARG	NE-CZ-NH2	-7.83	116.39	120.30
2	B	588[A]	ARG	CD-NE-CZ	7.70	134.38	123.60
2	B	588[B]	ARG	CD-NE-CZ	7.70	134.38	123.60
2	B	718	ARG	NE-CZ-NH1	7.68	124.14	120.30
1	A	230	ASP	CB-CG-OD1	7.59	125.13	118.30
2	B	734	ARG	CD-NE-CZ	7.57	134.20	123.60
2	B	718	ARG	NE-CZ-NH2	-7.17	116.72	120.30
1	A	96	TYR	CB-CG-CD2	-6.64	117.02	121.00
2	B	462	ARG	NE-CZ-NH2	-6.62	116.99	120.30
2	B	734	ARG	NE-CZ-NH2	-6.46	117.07	120.30
2	B	339	ASP	CB-CG-OD2	6.37	124.03	118.30
2	B	532	ARG	NE-CZ-NH2	-6.36	117.12	120.30
2	B	262	THR	CA-C-O	-6.34	106.78	120.10
1	A	73	TYR	CB-CG-CD1	6.10	124.66	121.00
1	A	76	ASP	CB-CG-OD1	6.04	123.74	118.30
2	B	532	ARG	NE-CZ-NH1	5.96	123.28	120.30
2	B	796	ARG	NE-CZ-NH1	-5.91	117.34	120.30
2	B	287	PHE	CB-CG-CD1	-5.90	116.67	120.80
2	B	742	ARG	NH1-CZ-NH2	5.86	125.84	119.40
1	A	47	ARG	NE-CZ-NH2	-5.70	117.45	120.30
2	B	550	ARG	NE-CZ-NH2	-5.67	117.47	120.30
2	B	590	ASP	CB-CG-OD1	5.54	123.29	118.30
2	B	556	ASP	CB-CG-OD2	-5.54	113.31	118.30
2	B	718	ARG	CD-NE-CZ	5.47	131.26	123.60
2	B	757	ASP	CB-CG-OD1	5.43	123.19	118.30
2	B	554	ARG	NE-CZ-NH1	-5.41	117.59	120.30
2	B	515	ASP	CB-CG-OD1	5.32	123.09	118.30
1	A	73	TYR	CB-CG-CD2	-5.30	117.82	121.00
2	B	365	ARG	NE-CZ-NH1	5.30	122.95	120.30
1	A	79	ARG	NE-CZ-NH2	5.30	122.95	120.30
1	A	197	TYR	CB-CG-CD1	-5.29	117.83	121.00
1	A	145	ARG	NE-CZ-NH1	5.24	122.92	120.30
2	B	468	PRO	O-C-N	-5.23	114.33	122.70
2	B	560	ARG	NE-CZ-NH2	-5.12	117.74	120.30
2	B	508	GLN	N-CA-CB	-5.10	101.42	110.60
1	A	133	ASP	CB-CG-OD1	5.06	122.86	118.30
2	B	404	ARG	NE-CZ-NH1	5.03	122.82	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:



Mol	Chain	Res	Type	Group
1	A	228	GLY	Mainchain
2	B	366[B]	GLU	Mainchain

## 5.2 Close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2036	0	1998	35	0
2	B	4499	0	4334	61	0
3	B	1	0	0	0	0
4	B	1	0	0	0	0
5	B	12	0	18	1	0
6	A	337	0	0	2	0
6	B	856	0	0	10	1
All	All	7742	0	6350	75	1

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 6.

All (75) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:181[A]:LEU:CD2	2:B:469:ARG:CZ	2.43	0.95
1:A:181[A]:LEU:HD21	2:B:469:ARG:CZ	1.97	0.94
2:B:261:PRO:HD2	2:B:332:ALA:HB2	1.58	0.86
1:A:181[A]:LEU:HD21	2:B:469:ARG:NH2	1.91	0.84
1:A:181[A]:LEU:HD22	2:B:469:ARG:NH1	1.93	0.83
1:A:181[A]:LEU:HD22	2:B:469:ARG:CZ	2.10	0.80
2:B:325:VAL:HG21	2:B:424:LYS:HE3	1.72	0.72
2:B:378[A]:THR:HG23	6:B:2313:HOH:O	1.93	0.69
2:B:477:LYS:HD3	6:B:2421:HOH:O	1.94	0.68
1:A:181[A]:LEU:HD13	2:B:469:ARG:HD3	1.76	0.67
2:B:783:HIS:HD2	2:B:786:ASP:OD2	1.78	0.66
2:B:285:PRO:HB2	2:B:287:PHE:CE2	2.31	0.66
1:A:260:TYR:O	2:B:262:THR:HG21	1.98	0.64
1:A:3:SER:HB2	1:A:6:GLU:HB2	1.81	0.62
1:A:181[A]:LEU:HD13	2:B:469:ARG:CD	2.29	0.62

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:56[A]:GLU:HG2	2:B:370:THR:HB	1.81	0.61
1:A:156:LEU:HD11	1:A:212:LEU:HD13	1.82	0.61
2:B:783:HIS:HE1	2:B:811:GLU:OE2	1.85	0.59
2:B:368:THR:OG1	2:B:378[A]:THR:HG22	2.03	0.59
1:A:64:LYS:HA	1:A:64:LYS:HE3	1.85	0.58
1:A:142:MET:HE3	6:A:2204:HOH:O	2.05	0.56
1:A:212:LEU:HB3	6:B:2414:HOH:O	2.06	0.56
2:B:294:TYR:CE2	2:B:312:PRO:HB3	2.42	0.55
2:B:305:TYR:CZ	2:B:422:LYS:HE3	2.44	0.53
2:B:477:LYS:H	2:B:477:LYS:HZ3	1.56	0.53
1:A:146:PHE:HB3	2:B:261:PRO:CD	2.39	0.53
2:B:424:LYS:HG3	6:B:2362:HOH:O	2.08	0.52
1:A:156:LEU:CD1	1:A:212:LEU:HD13	2.38	0.52
2:B:477:LYS:H	2:B:477:LYS:NZ	2.07	0.52
2:B:391:GLN:NE2	6:B:2323:HOH:O	2.42	0.52
2:B:707:ASN:C	2:B:707:ASN:HD22	2.14	0.51
1:A:146:PHE:HB3	2:B:261:PRO:N	2.26	0.49
1:A:146:PHE:CD1	2:B:261:PRO:HA	2.48	0.49
2:B:396:THR:HG22	2:B:398:THR:OG1	2.12	0.49
2:B:261:PRO:HD2	2:B:332:ALA:CB	2.38	0.49
2:B:285:PRO:HG3	2:B:320:PHE:CZ	2.48	0.48
2:B:553[A]:SER:OG	2:B:742:ARG:HB3	2.14	0.48
2:B:305:TYR:CE2	2:B:422:LYS:HE3	2.48	0.48
2:B:519:PHE:CD2	5:B:905:EDO:H11	2.47	0.48
2:B:307:VAL:HG11	2:B:421:MET:HB3	1.96	0.47
1:A:181[A]:LEU:CD2	2:B:469:ARG:NH1	2.64	0.47
2:B:346:SER:HB2	2:B:359:TRP:CH2	2.51	0.46
1:A:213:PRO:HG2	6:B:2334:HOH:O	2.15	0.46
1:A:29:TYR:CZ	2:B:818[B]:VAL:HG11	2.51	0.45
1:A:181[B]:LEU:HG	1:A:201:PHE:HB3	1.97	0.45
1:A:29:TYR:HA	1:A:95:GLY:O	2.16	0.45
2:B:537:LYS:HD2	2:B:540:LEU:HG	1.98	0.45
1:A:56[A]:GLU:HG3	6:A:2096:HOH:O	2.17	0.45
1:A:217:LEU:HB3	1:A:218:PRO:HD2	1.98	0.45
1:A:4:SER:HB3	6:B:2845:HOH:O	2.16	0.45
2:B:389[A]:ILE:HD13	2:B:401:ALA:HB2	2.00	0.44
2:B:264[A]:SER:OG	2:B:285:PRO:HA	2.18	0.44
2:B:311:THR:HA	2:B:312:PRO:HD3	1.91	0.44
1:A:214:ARG:O	1:A:214:ARG:HG3	2.17	0.43
2:B:281:MET:O	2:B:748:MET:HA	2.18	0.43
2:B:431[A]:GLN:NE2	6:B:2375:HOH:O	2.52	0.43
2:B:767:ALA:HA	2:B:768:PRO:C	2.38	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:150:THR:HG21	2:B:340[B]:ILE:HD11	2.01	0.42
2:B:393:ASP:HB2	6:B:2324:HOH:O	2.19	0.42
1:A:74:TRP:HA	1:A:75:PRO:HD2	1.91	0.42
1:A:188:THR:HA	2:B:506:SER:O	2.18	0.42
2:B:263:GLY:O	2:B:264[A]:SER:OG	2.29	0.42
1:A:26:HIS:CD2	2:B:818[B]:VAL:HG23	2.54	0.42
2:B:427:GLN:HG2	6:B:2368:HOH:O	2.19	0.42
1:A:208:THR:HG21	2:B:472:VAL:HG12	2.01	0.42
2:B:343:GLU:O	2:B:398:THR:HA	2.20	0.42
1:A:146:PHE:HB3	2:B:261:PRO:HD3	2.01	0.41
1:A:67[A]:LYS:HG2	2:B:379:PHE:CE1	2.56	0.41
2:B:264[B]:SER:HB3	2:B:285:PRO:HA	2.03	0.41
2:B:431[A]:GLN:HE21	2:B:431[A]:GLN:HB3	1.69	0.41
2:B:311:THR:HB	2:B:318:LEU:HA	2.02	0.41
2:B:328:TRP:HA	2:B:443:TYR:O	2.21	0.41
2:B:262:THR:HA	2:B:334:PHE:HZ	1.85	0.40
2:B:623:VAL:HG13	2:B:631:ASP:HB2	2.03	0.40
1:A:198:PRO:HG2	2:B:488:MET:SD	2.60	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
6:B:2285:HOH:O	6:B:2686:HOH:O[1_465]	2.00	0.20

## 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	260/260 (100%)	255 (98%)	5 (2%)	0	100	100
2	B	578/560 (103%)	563 (97%)	13 (2%)	2 (0%)	50	31
All	All	838/820 (102%)	818 (98%)	18 (2%)	2 (0%)	59	38

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	264[A]	SER
2	B	264[B]	SER

### 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	213/211 (101%)	209 (98%)	4 (2%)	69	56
2	B	482/462 (104%)	468 (97%)	14 (3%)	55	37
All	All	695/673 (103%)	677 (97%)	18 (3%)	61	41

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	64	LYS
1	A	202	ASN
1	A	212	LEU
1	A	251	GLN
2	B	283	ASN
2	B	417	TRP
2	B	431[A]	GLN
2	B	431[B]	GLN
2	B	436	GLN
2	B	477	LYS
2	B	516[A]	LEU
2	B	516[B]	LEU
2	B	537	LYS
2	B	575	GLN
2	B	695	GLU
2	B	707	ASN
2	B	796	ARG
2	B	819	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	196	ASN

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Mol	Chain	Res	Type
1	A	202	ASN
1	A	205	ASN
2	B	356	ASN
2	B	427	GLN
2	B	508	GLN
2	B	575	GLN
2	B	581	GLN
2	B	707	ASN
2	B	783	HIS

### 5.3.3 RNA ⓘ

There are no RNA chains 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 carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 5 ligands modelled in this entry, 2 are monoatomic - leaving 3 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	$\# Z  > 2$
5	EDO	B	903	-	3,3,3	0.49	0	2,2,2	0.15	0
5	EDO	B	904	-	3,3,3	0.58	0	2,2,2	0.37	0
5	EDO	B	905	-	3,3,3	0.61	0	2,2,2	0.27	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
5	EDO	B	903	-	-	0/1/1/1	0/0/0/0
5	EDO	B	904	-	-	0/1/1/1	0/0/0/0
5	EDO	B	905	-	-	0/1/1/1	0/0/0/0

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.

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	258/260 (99%)	-0.24	11 (4%) 34 27	6, 13, 34, 50	0
2	B	560/560 (100%)	-0.60	2 (0%) 90 89	5, 10, 20, 44	0
All	All	818/820 (99%)	-0.49	13 (1%) 68 64	5, 11, 27, 50	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	3	SER	6.1
2	B	261	PRO	5.0
2	B	262	THR	4.4
1	A	202	ASN	3.2
1	A	229	ALA	2.7
1	A	203	GLN	2.7
1	A	239	GLY	2.6
1	A	201	PHE	2.6
1	A	237	ALA	2.4
1	A	212	LEU	2.4
1	A	205	ASN	2.3
1	A	235	ALA	2.2
1	A	210	ALA	2.1

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

### 6.3 Carbohydrates ⓘ

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
5	EDO	B	903	4/4	0.08	0.61	14,16,16,20	0
5	EDO	B	904	4/4	0.13	0.37	30,31,31,32	0
5	EDO	B	905	4/4	0.07	0.06	9,10,11,12	0
4	CL	B	902	1/1	0.06	-1.15	16,16,16,16	0
3	CA	B	901	1/1	0.04	-2.37	8,8,8,8	0

## 6.5 Other polymers

There are no such residues in this entry.