



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:45 AM UTC

PDB ID : 5ZTX / pdb_00005ztx
Title : co-factor free Transaminase
Authors : Park, H.H.; Shin, Y.C.
Deposited on : 2018-05-05
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 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

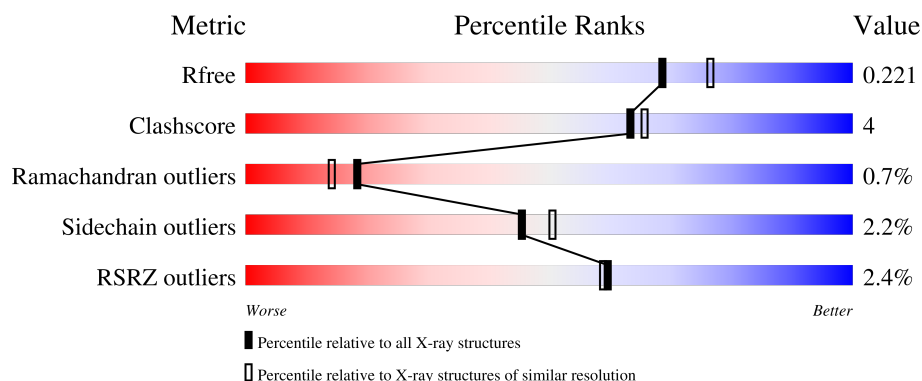
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (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 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	469	% 84% 6% 10%
1	B	469	4% 82% 10% 7%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7304 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 transaminase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	424	Total	C	N	O	S	0	0	0
			3287	2105	558	605	19			
1	B	435	Total	C	N	O	S	0	0	0
			3401	2178	577	628	18			

There are 32 discrepancies between the modelled and reference sequences:

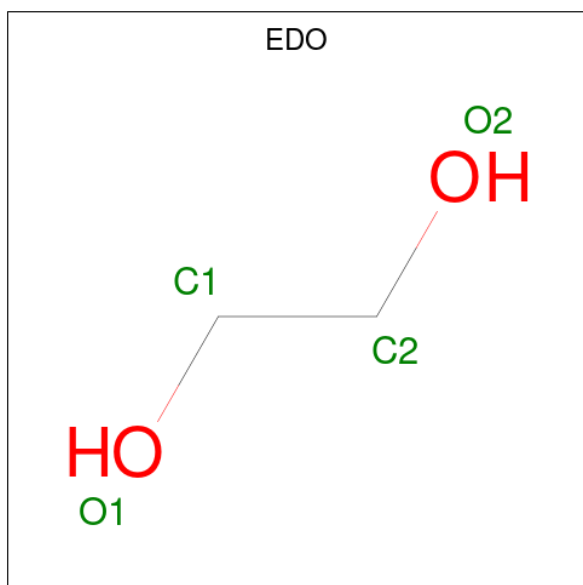
Chain	Residue	Modelled	Actual	Comment	Reference
A	-13	MET	-	initiating methionine	UNP F2XBU9
A	-12	ALA	-	expression tag	UNP F2XBU9
A	-11	SER	-	expression tag	UNP F2XBU9
A	-10	MET	-	expression tag	UNP F2XBU9
A	-9	THR	-	expression tag	UNP F2XBU9
A	-8	GLY	-	expression tag	UNP F2XBU9
A	-7	GLY	-	expression tag	UNP F2XBU9
A	-6	GLN	-	expression tag	UNP F2XBU9
A	-5	GLN	-	expression tag	UNP F2XBU9
A	-4	MET	-	expression tag	UNP F2XBU9
A	-3	GLY	-	expression tag	UNP F2XBU9
A	-2	ARG	-	expression tag	UNP F2XBU9
A	-1	GLY	-	expression tag	UNP F2XBU9
A	0	SER	-	expression tag	UNP F2XBU9
A	454	ALA	-	expression tag	UNP F2XBU9
A	455	ALA	-	expression tag	UNP F2XBU9
B	-13	MET	-	initiating methionine	UNP F2XBU9
B	-12	ALA	-	expression tag	UNP F2XBU9
B	-11	SER	-	expression tag	UNP F2XBU9
B	-10	MET	-	expression tag	UNP F2XBU9
B	-9	THR	-	expression tag	UNP F2XBU9
B	-8	GLY	-	expression tag	UNP F2XBU9
B	-7	GLY	-	expression tag	UNP F2XBU9
B	-6	GLN	-	expression tag	UNP F2XBU9
B	-5	GLN	-	expression tag	UNP F2XBU9

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-4	MET	-	expression tag	UNP F2XBU9
B	-3	GLY	-	expression tag	UNP F2XBU9
B	-2	ARG	-	expression tag	UNP F2XBU9
B	-1	GLY	-	expression tag	UNP F2XBU9
B	0	SER	-	expression tag	UNP F2XBU9
B	454	ALA	-	expression tag	UNP F2XBU9
B	455	ALA	-	expression tag	UNP F2XBU9

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



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

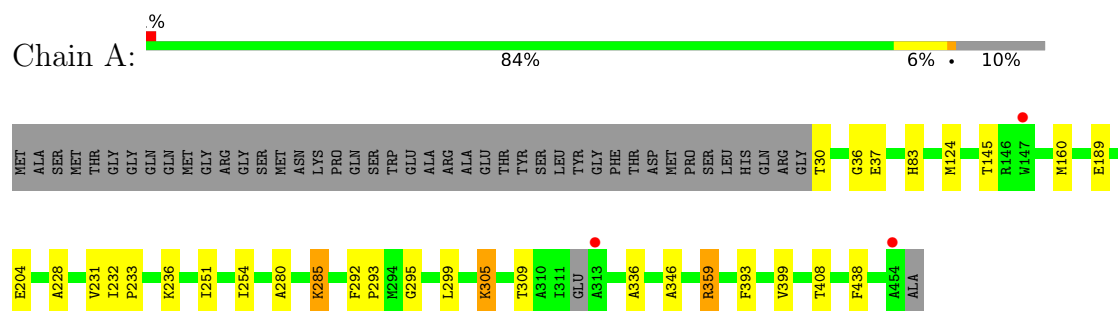
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	267	Total O 267 267	0	0
3	B	345	Total O 345 345	0	0

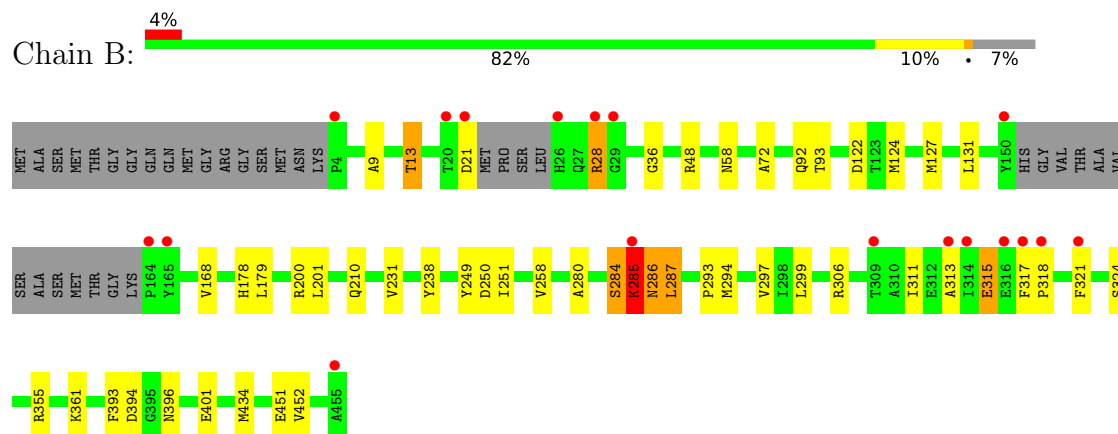
3 Residue-property plots [i](#)

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: transaminase



- Molecule 1: transaminase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	78.28Å 95.31Å 123.03Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.00 20.00 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.7 (20.00-2.00) 99.8 (20.00-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.26 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.8.0103	Depositor
R, R_{free}	0.177 , 0.230 (Not available) , 0.221	Depositor DCC
R_{free} test set	3193 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	20.9	Xtriage
Anisotropy	0.087	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 33.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7304	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.71% 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

5.1 Standard geometry

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	1.33	8/3369 (0.2%)	1.16	3/4567 (0.1%)
1	B	1.24	7/3488 (0.2%)	1.19	9/4726 (0.2%)
All	All	1.28	15/6857 (0.2%)	1.18	12/9293 (0.1%)

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	B	0	1

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	250	ASP	N-CA	7.15	1.57	1.46
1	B	251	ILE	N-CA	-6.76	1.40	1.46
1	A	292	PHE	CA-C	6.19	1.58	1.52
1	B	396	ASN	CA-C	-5.85	1.44	1.52
1	B	286	ASN	CA-C	5.85	1.60	1.52
1	B	122	ASP	N-CA	-5.81	1.39	1.46
1	A	295	GLY	N-CA	5.61	1.50	1.44
1	A	285	LYS	CA-C	5.59	1.60	1.52
1	B	72	ALA	N-CA	5.53	1.52	1.46
1	A	336	ALA	C-O	5.44	1.30	1.24
1	A	36	GLY	CA-C	-5.29	1.47	1.52
1	A	359	ARG	N-CA	-5.19	1.40	1.46
1	A	251	ILE	N-CA	5.16	1.50	1.46
1	A	346	ALA	N-CA	5.12	1.52	1.46
1	B	250	ASP	CB-CG	5.12	1.64	1.52

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	249	TYR	N-CA-C	7.97	125.57	113.61
1	B	285	LYS	N-CA-C	6.92	125.54	110.80
1	B	168	VAL	N-CA-C	-6.54	105.33	111.48
1	A	233	PRO	O-C-N	5.92	123.93	121.15
1	B	251	ILE	N-CA-C	-5.72	102.48	108.15
1	B	28	ARG	N-CA-C	5.62	118.25	111.40
1	B	92	GLN	CB-CA-C	-5.41	102.16	110.81
1	B	238	TYR	N-CA-C	5.30	117.49	111.02
1	A	232	ILE	CA-C-N	-5.25	115.88	119.66
1	A	232	ILE	C-N-CA	-5.25	115.88	119.66
1	B	36	GLY	N-CA-C	5.25	118.94	111.12
1	B	200	ARG	NE-CZ-NH2	5.12	123.81	119.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	284	SER	Peptide

5.2 Too-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 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	3287	0	3229	19	0
1	B	3401	0	3319	35	0
2	A	4	0	6	2	0
3	A	267	0	0	3	0
3	B	345	0	0	4	0
All	All	7304	0	6554	53	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 4.

All (53) 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:258:VAL:HG13	1:B:284:SER:HB3	1.60	0.84
1:B:127:MET:HE1	1:B:299:LEU:HD11	1.58	0.83
1:B:9:ALA:O	1:B:13:THR:HG23	1.77	0.83
1:B:258:VAL:HG13	1:B:284:SER:CB	2.12	0.80
1:A:189:GLU:OE1	3:A:601:HOH:O	2.07	0.71
1:B:124:MET:CE	1:B:297:VAL:HG13	2.21	0.71
1:B:284:SER:O	1:B:287:LEU:HB2	1.91	0.71
1:B:124:MET:HE2	1:B:297:VAL:HG13	1.74	0.69
1:B:394:ASP:HB3	3:B:678:HOH:O	1.92	0.69
1:A:124:MET:CE	1:A:280:ALA:CB	2.70	0.69
1:A:124:MET:HE3	1:A:280:ALA:HB3	1.75	0.68
1:B:58:ASN:HB3	1:B:285:LYS:CG	2.27	0.65
1:A:228:ALA:HB1	2:A:501:EDO:H22	1.79	0.64
1:B:58:ASN:HB3	1:B:285:LYS:CD	2.29	0.63
1:B:355:ARG:HG2	1:B:434:MET:HE1	1.81	0.62
1:B:58:ASN:HB3	1:B:285:LYS:HG3	1.82	0.62
1:B:355:ARG:CG	1:B:434:MET:HE1	2.31	0.60
1:A:305:LYS:O	1:A:309:THR:HG23	2.02	0.59
1:A:124:MET:HE3	1:A:280:ALA:CB	2.34	0.58
1:B:285:LYS:CG	1:B:286:ASN:N	2.67	0.57
1:B:311:ILE:O	1:B:315:GLU:HA	2.05	0.56
1:B:48:ARG:HD3	3:B:610:HOH:O	2.07	0.55
1:B:210:GLN:HG2	3:B:784:HOH:O	2.08	0.54
1:A:124:MET:CE	1:A:280:ALA:HB1	2.38	0.53
1:A:124:MET:HE1	1:A:280:ALA:CB	2.39	0.53
1:B:127:MET:HE1	1:B:299:LEU:HD21	1.90	0.52
1:A:124:MET:HE1	1:A:299:LEU:HD23	1.93	0.51
1:B:131:LEU:HD11	1:B:306:ARG:HB3	1.94	0.49
1:B:127:MET:HG2	1:B:321:PHE:CZ	2.48	0.48
1:B:124:MET:HE1	1:B:280:ALA:HB1	1.95	0.48
1:A:228:ALA:CB	2:A:501:EDO:H22	2.42	0.48
1:B:127:MET:HG2	1:B:321:PHE:CE2	2.49	0.48
1:B:127:MET:CE	1:B:299:LEU:HD11	2.40	0.46
1:B:127:MET:HE1	1:B:299:LEU:CD1	2.38	0.46
1:A:393:PHE:CG	1:A:399:VAL:HG21	2.52	0.45
1:A:293:PRO:HG2	1:B:293:PRO:HG2	1.99	0.44
1:B:311:ILE:HG12	1:B:318:PRO:HD3	2.00	0.44
1:B:393:PHE:CE1	1:B:452:VAL:HG11	2.53	0.43
1:A:83:HIS:HE1	3:A:819:HOH:O	2.01	0.43
1:A:359:ARG:HG2	1:A:438:PHE:CG	2.54	0.43
1:B:285:LYS:HG3	1:B:286:ASN:N	2.32	0.43
1:B:9:ALA:O	1:B:13:THR:CG2	2.58	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:93:THR:HG21	1:B:324:SER:HA	1.99	0.42
1:A:124:MET:HE1	1:A:280:ALA:HB1	2.00	0.42
1:A:145:THR:HG21	1:A:160:MET:HB3	2.02	0.42
1:B:93:THR:HG22	3:B:587:HOH:O	2.19	0.42
1:B:58:ASN:CB	1:B:285:LYS:HG3	2.48	0.41
1:B:285:LYS:CG	1:B:286:ASN:H	2.32	0.41
1:A:124:MET:HE3	1:A:254:ILE:HD12	2.01	0.41
1:A:359:ARG:HG2	1:A:438:PHE:CD2	2.56	0.41
1:B:178:HIS:CE1	1:B:179:LEU:O	2.74	0.41
1:B:285:LYS:HG2	1:B:286:ASN:H	1.87	0.40
1:A:204:GLU:CG	3:A:689:HOH:O	2.70	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	420/469 (90%)	404 (96%)	14 (3%)	2 (0%)	24	21
1	B	429/469 (92%)	408 (95%)	17 (4%)	4 (1%)	14	9
All	All	849/938 (90%)	812 (96%)	31 (4%)	6 (1%)	18	14

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	285	LYS
1	B	313	ALA
1	B	317	PHE
1	B	231	VAL
1	A	285	LYS
1	A	231	VAL

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	342/378 (90%)	337 (98%)	5 (2%)	57	64
1	B	353/378 (93%)	343 (97%)	10 (3%)	38	41
All	All	695/756 (92%)	680 (98%)	15 (2%)	45	50

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	THR
1	A	37	GLU
1	A	236	LYS
1	A	305	LYS
1	A	408	THR
1	B	13	THR
1	B	21	ASP
1	B	28	ARG
1	B	201	LEU
1	B	287	LEU
1	B	294	MET
1	B	315	GLU
1	B	361	LYS
1	B	401	GLU
1	B	451	GLU

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

Mol	Chain	Res	Type
1	A	83	HIS
1	A	148	ASN
1	A	178	HIS
1	B	45	ASN
1	B	53	ASN
1	B	139	GLN
1	B	178	HIS

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Mol	Chain	Res	Type
1	B	210	GLN
1	B	267	ASN
1	B	319	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is 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	A	501	-	3,3,3	0.97	0	2,2,2	1.12	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	501	-	-	0/1/1/1	-

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 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	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	424/469 (90%)	-0.45	3 (0%) 84 84	11, 18, 34, 56	0
1	B	435/469 (92%)	-0.10	18 (4%) 41 40	14, 23, 46, 83	0
All	All	859/938 (91%)	-0.27	21 (2%) 59 59	11, 21, 43, 83	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	454	ALA	4.4
1	B	150	TYR	4.3
1	B	313	ALA	3.9
1	B	164	PRO	3.9
1	B	455	ALA	3.4
1	B	4	PRO	3.1
1	B	314	ILE	3.1
1	B	26	HIS	3.0
1	B	317	PHE	3.0
1	B	318	PRO	2.8
1	B	285	LYS	2.7
1	B	29	GLY	2.5
1	B	316	GLU	2.5
1	A	147	TRP	2.4
1	B	309	THR	2.4
1	A	313	ALA	2.3
1	B	20	THR	2.2
1	B	21	ASP	2.2
1	B	321	PHE	2.1
1	B	28	ARG	2.1
1	B	165	TYR	2.1

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [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
2	EDO	A	501	4/4	0.74	0.14	35,35,36,42	0

6.5 Other polymers [i](#)

There are no such residues in this entry.