



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 14, 2023 – 11:02 AM JST

PDB ID : 5ZPI  
Title : Copper amine oxidase from *Arthrobacter globiformis* anaerobically reduced by ethylamine at pH 6 at 293 K (3)  
Authors : Murakawa, T.; Baba, S.; Kawano, Y.; Hayashi, H.; Yano, T.; Tanizawa, K.; Kumasaka, T.; Yamamoto, M.; Okajima, T.  
Deposited on : 2018-04-16  
Resolution : 1.75 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](mailto: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>.

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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 : 1.8.5 (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

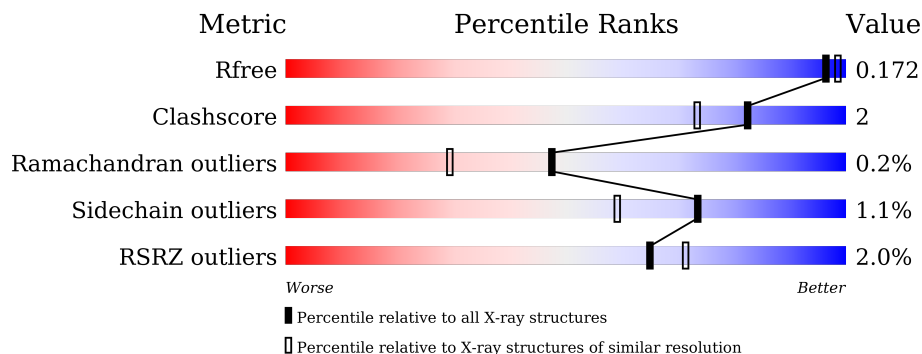
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.75 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	3764 (1.76-1.72)
Clashscore	141614	3923 (1.76-1.72)
Ramachandran outliers	138981	3878 (1.76-1.72)
Sidechain outliers	138945	3878 (1.76-1.72)
RSRZ outliers	127900	3705 (1.76-1.72)

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  $\geq 3$ , 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	620	 2% 93% 6%
1	B	620	 2% 95% 5%

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
4	EDO	A	1005	-	-	-	X
4	EDO	B	705	-	-	-	X

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 10757 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 Phenylethylamine oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	620	4939	3128	863	935	13	0	12	0
1	B	620	4919	3114	860	934	11	0	8	0

- Molecule 2 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total 1	Cu 1	0	0
2	B	1	Total 1	Cu 1	0	0

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total 1	Na 1	0	0
3	B	1	Total 1	Na 1	0	0

- Molecule 4 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
4	A	1	Total C O 4 2 2	0	0
4	A	1	Total C O 4 2 2	0	0
4	A	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0

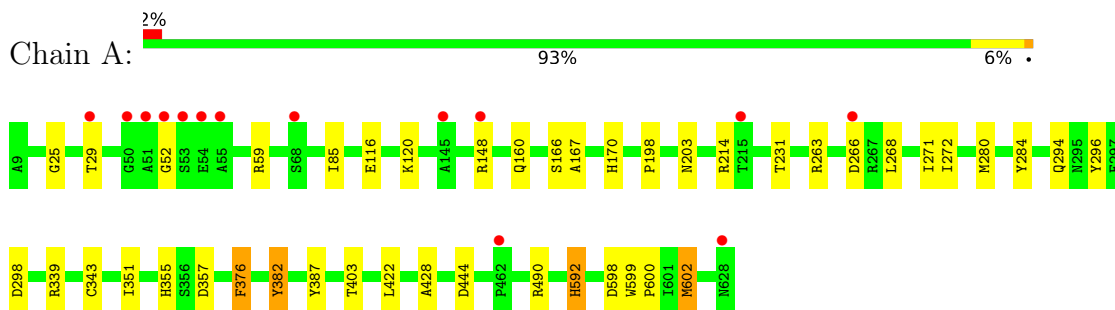
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	416	Total O 416 416	0	5
5	B	451	Total O 451 451	0	4

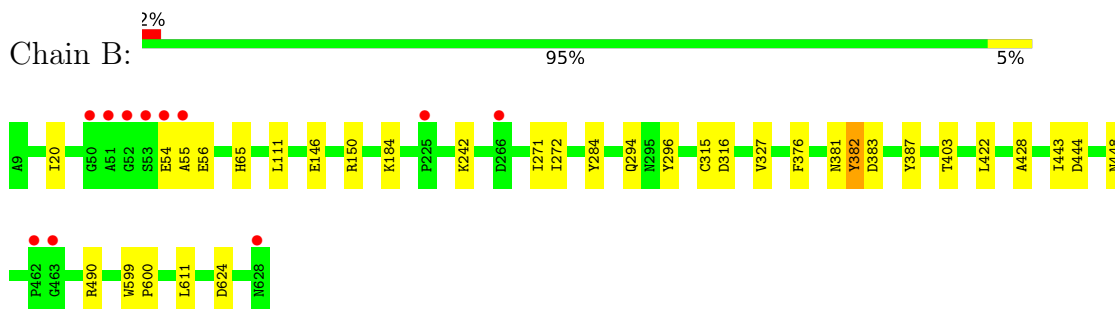
### 3 Residue-property plots [i](#)

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: Phenylethylamine oxidase



- Molecule 1: Phenylethylamine oxidase



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	193.13Å 64.22Å 158.78Å 90.00° 117.06° 90.00°	Depositor
Resolution (Å)	40.35 – 1.75 45.43 – 1.75	Depositor EDS
% Data completeness (in resolution range)	99.3 (40.35-1.75) 93.1 (45.43-1.75)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.21 (at 1.75Å)	Xtrriage
Refinement program	PHENIX 1.11.1_2575	Depositor
R, $R_{free}$	0.153 , 0.175 0.150 , 0.172	Depositor DCC
$R_{free}$ test set	8750 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	17.6	Xtrriage
Anisotropy	0.632	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 53.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.98	EDS
Total number of atoms	10757	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	25.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 78.77 % 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 6.4433e-07. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

<sup>2</sup>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: NA, TYQ, EDO, CU

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	0.34	0/5069	0.58	2/6900 (0.0%)
1	B	0.37	0/5037	0.58	0/6857
All	All	0.35	0/10106	0.58	2/13757 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	343[A]	CYS	CA-CB-SG	5.31	123.56	114.00
1	A	343[B]	CYS	CA-CB-SG	5.31	123.56	114.00

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	4939	0	4792	27	0
1	B	4919	0	4758	24	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	12	0	18	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	16	0	24	0	0
5	A	416	0	0	2	0
5	B	451	0	0	2	0
All	All	10757	0	9592	48	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 (48) 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:A:166:SER:H	1:A:203:ASN:HD21	1.24	0.86
1:A:490[A]:ARG:NH1	5:A:1101:HOH:O	2.16	0.76
1:A:355:HIS:HD2	1:B:383:ASP:OD1	1.74	0.70
1:B:443:ILE:H	1:B:448:ASN:HD21	1.40	0.69
1:A:263[A]:ARG:HG2	1:A:268:LEU:HD13	1.77	0.66
1:B:294:GLN:OE1	1:B:296[B]:TYR:OH	2.12	0.65
1:A:355:HIS:HE1	1:A:357:ASP:OD1	1.82	0.63
1:B:376:PHE:HE1	1:B:387:TYR:HH	1.44	0.62
1:B:20:ILE:HD12	1:B:327:VAL:HG12	1.82	0.62
1:A:170:HIS:HD2	1:A:198:PRO:O	1.82	0.62
1:A:116:GLU:O	1:A:120:LYS:HG2	2.01	0.60
1:B:599:TRP:CD2	1:B:600:PRO:HA	2.37	0.60
1:B:54:GLU:O	1:B:56:GLU:N	2.35	0.60
1:A:231:THR:HG22	4:A:1005:EDO:H11	1.84	0.59
1:A:599:TRP:CD2	1:A:600:PRO:HA	2.41	0.56
1:A:294:GLN:OE1	1:A:296[B]:TYR:OH	2.13	0.55
1:B:65:HIS:HD2	5:B:1225:HOH:O	1.88	0.55
1:A:166:SER:H	1:A:203:ASN:ND2	2.01	0.54
1:A:592:HIS:HE1	1:A:598:ASP:OD2	1.91	0.54
1:A:382[B]:TYQ:HD2	1:A:403:THR:O	2.09	0.53
4:A:1004:EDO:H11	5:A:1423:HOH:O	2.09	0.51
1:A:280[B]:MET:HG3	1:A:298:ASP:HB2	1.93	0.51
1:B:422:LEU:HD11	1:B:428:ALA:HB2	1.92	0.51
1:B:382[B]:TYQ:HD2	1:B:403:THR:O	2.12	0.49
4:A:1004:EDO:H11	1:B:490:ARG:HG3	1.95	0.49
1:B:146:GLU:O	1:B:150:ARG:HD3	2.12	0.49
1:B:296[A]:TYR:CZ	1:B:382[A]:TYQ:N5	2.81	0.48
1:A:160:GLN:HG3	1:A:167:ALA:HB2	1.96	0.47
1:B:381:ASN:OD1	1:B:382[A]:TYQ:N5	2.47	0.46
1:B:376:PHE:HE1	1:B:387:TYR:OH	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:284:TYR:OH	1:A:382[A]:TYQ:OH	2.16	0.46
1:A:382[B]:TYQ:N5	1:A:602[B]:MET:SD	2.89	0.45
1:B:284:TYR:OH	1:B:382[A]:TYQ:OH	2.20	0.45
1:A:166:SER:N	1:A:203:ASN:HD21	2.03	0.44
1:A:214:ARG:HG3	1:B:624:ASP:O	2.18	0.44
1:A:271:ILE:HG22	1:A:272:ILE:HG13	2.00	0.43
1:A:422:LEU:HD11	1:A:428:ALA:HB2	2.02	0.42
1:B:65:HIS:HE1	5:B:868:HOH:O	2.03	0.42
1:A:351[A]:ILE:HD11	1:B:316:ASP:HA	2.00	0.42
1:B:111:LEU:HD11	1:B:184:LYS:HA	2.01	0.41
1:A:376:PHE:HE1	1:A:387:TYR:OH	2.03	0.41
1:A:490[B]:ARG:NH1	4:A:1004:EDO:O2	2.53	0.41
1:A:25:GLY:O	1:A:29:THR:HG23	2.21	0.41
1:B:296[A]:TYR:CE2	1:B:382[A]:TYQ:N5	2.89	0.41
1:B:271:ILE:HG22	1:B:272:ILE:HG13	2.02	0.40
1:A:355:HIS:CD2	1:B:383:ASP:OD1	2.65	0.40
1:A:59:ARG:NH1	1:A:85:ILE:HD13	2.36	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	628/620 (101%)	607 (97%)	20 (3%)	1 (0%)	47 29
1	B	624/620 (101%)	599 (96%)	24 (4%)	1 (0%)	47 29
All	All	1252/1240 (101%)	1206 (96%)	44 (4%)	2 (0%)	47 29

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	55	ALA

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Mol	Chain	Res	Type
1	A	52	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	524/513 (102%)	516 (98%)	8 (2%)	65	47
1	B	520/513 (101%)	516 (99%)	4 (1%)	81	72
All	All	1044/1026 (102%)	1032 (99%)	12 (1%)	73	59

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	148	ARG
1	A	266	ASP
1	A	339	ARG
1	A	376	PHE
1	A	444	ASP
1	A	592	HIS
1	A	602[A]	MET
1	A	602[B]	MET
1	B	242	LYS
1	B	315	CYS
1	B	444	ASP
1	B	611	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	170	HIS
1	A	203	ASN
1	A	355	HIS
1	A	592	HIS
1	B	65	HIS

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Mol	Chain	Res	Type
1	B	110	GLN
1	B	259	ASN
1	B	448	ASN
1	B	519	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](#)

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	TYQ	A	382[B]	2	13,14,15	1.10	0	15,19,21	1.71	3 (20%)
1	TYQ	B	382[B]	2	13,14,15	0.97	0	15,19,21	1.66	3 (20%)
1	TYQ	A	382[A]	-	13,14,15	1.14	1 (7%)	15,19,21	1.93	3 (20%)
1	TYQ	B	382[A]	-	13,14,15	1.13	1 (7%)	15,19,21	1.85	4 (26%)

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. '2' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	TYQ	A	382[B]	2	-	3/5/6/8	0/1/1/1
1	TYQ	B	382[B]	2	-	3/5/6/8	0/1/1/1
1	TYQ	A	382[A]	-	-	2/5/6/8	0/1/1/1
1	TYQ	B	382[A]	-	-	2/5/6/8	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	382[A]	TYQ	CB-CG	-2.80	1.47	1.51
1	B	382[A]	TYQ	CB-CG	-2.59	1.48	1.51

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	382[A]	TYQ	CB-CA-C	-4.85	102.38	111.47
1	B	382[A]	TYQ	CB-CA-C	-4.23	103.54	111.47
1	B	382[B]	TYQ	CG-CB-CA	-3.86	108.56	114.53
1	A	382[B]	TYQ	OH-CZ-CE2	3.61	122.29	116.25
1	A	382[B]	TYQ	CG-CB-CA	-3.15	109.65	114.53
1	B	382[B]	TYQ	OH-CZ-CE2	2.95	121.19	116.25
1	B	382[A]	TYQ	OH-CZ-CE2	2.94	121.17	116.25
1	A	382[A]	TYQ	OH-CZ-CE2	2.74	120.84	116.25
1	B	382[A]	TYQ	CZ-CE2-N5	2.72	123.09	118.77
1	A	382[B]	TYQ	CB-CA-C	-2.36	107.05	111.47
1	B	382[B]	TYQ	CB-CA-C	-2.33	107.11	111.47
1	A	382[A]	TYQ	CB-CG-CD1	2.18	123.23	120.95
1	B	382[A]	TYQ	CB-CG-CD1	2.10	123.15	120.95

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	382[A]	TYQ	N-CA-CB-CG
1	A	382[A]	TYQ	C-CA-CB-CG
1	A	382[B]	TYQ	C-CA-CB-CG
1	B	382[A]	TYQ	N-CA-CB-CG
1	B	382[B]	TYQ	C-CA-CB-CG
1	A	382[B]	TYQ	N-CA-CB-CG
1	B	382[B]	TYQ	N-CA-CB-CG
1	B	382[A]	TYQ	C-CA-CB-CG
1	A	382[B]	TYQ	CA-CB-CG-CD2
1	B	382[B]	TYQ	CA-CB-CG-CD2

There are no ring outliers.

4 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	382[B]	TYQ	2	0
1	B	382[B]	TYQ	1	0
1	A	382[A]	TYQ	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	382[A]	TYQ	5	0

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 11 ligands modelled in this entry, 4 are monoatomic - leaving 7 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	$\# Z  > 2$
4	EDO	B	704	-	3,3,3	0.49	0	2,2,2	0.20	0
4	EDO	A	1003	-	3,3,3	0.50	0	2,2,2	0.25	0
4	EDO	A	1005	-	3,3,3	0.42	0	2,2,2	0.30	0
4	EDO	B	703	-	3,3,3	0.49	0	2,2,2	0.36	0
4	EDO	A	1004	-	3,3,3	0.42	0	2,2,2	0.11	0
4	EDO	B	705	-	3,3,3	0.47	0	2,2,2	0.33	0
4	EDO	B	706	-	3,3,3	0.52	0	2,2,2	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
4	EDO	B	704	-	-	1/1/1/1	-
4	EDO	A	1003	-	-	1/1/1/1	-
4	EDO	A	1005	-	-	0/1/1/1	-
4	EDO	B	703	-	-	0/1/1/1	-
4	EDO	A	1004	-	-	0/1/1/1	-
4	EDO	B	705	-	-	1/1/1/1	-
4	EDO	B	706	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	705	EDO	O1-C1-C2-O2
4	A	1003	EDO	O1-C1-C2-O2
4	B	704	EDO	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1005	EDO	1	0
4	A	1004	EDO	3	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, 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	619/620 (99%)	-0.28	14 (2%) 60 66	11, 22, 46, 157	0
1	B	619/620 (99%)	-0.36	11 (1%) 68 74	9, 19, 40, 172	0
All	All	1238/1240 (99%)	-0.32	25 (2%) 65 71	9, 21, 43, 172	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	51	ALA	21.2
1	A	51	ALA	19.9
1	A	52	GLY	13.5
1	B	52	GLY	10.3
1	B	55	ALA	8.6
1	A	53	SER	8.5
1	B	53	SER	8.1
1	B	54	GLU	7.4
1	A	54	GLU	6.8
1	A	55	ALA	6.4
1	A	50	GLY	5.2
1	B	628	ASN	4.4
1	A	628	ASN	4.2
1	B	50	GLY	3.4
1	A	266	ASP	3.4
1	B	462	PRO	3.1
1	A	462	PRO	2.9
1	B	266	ASP	2.6
1	B	225	PRO	2.4
1	B	463	GLY	2.4
1	A	29	THR	2.2
1	A	215	THR	2.2
1	A	145	ALA	2.1
1	A	68	SER	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	148	ARG	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, 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	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
1	TYQ	A	382[A]	14/15	0.95	0.14	15,33,43,47	11
1	TYQ	A	382[B]	14/15	0.95	0.14	15,28,36,37	11
1	TYQ	B	382[A]	14/15	0.97	0.12	12,26,30,33	11
1	TYQ	B	382[B]	14/15	0.97	0.12	12,21,33,42	11

## 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, 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	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
4	EDO	A	1005	4/4	0.63	0.66	78,79,79,81	0
4	EDO	B	705	4/4	0.68	0.50	78,79,79,79	0
4	EDO	A	1004	4/4	0.70	0.36	66,67,68,71	0
4	EDO	B	703	4/4	0.80	0.31	53,58,58,62	0
4	EDO	B	704	4/4	0.89	0.16	51,51,53,58	0
4	EDO	A	1003	4/4	0.93	0.27	50,53,54,60	0
4	EDO	B	706	4/4	0.96	0.07	16,31,33,34	0
3	NA	A	1002	1/1	0.99	0.07	14,14,14,14	0
3	NA	B	702	1/1	0.99	0.09	10,10,10,10	0
2	CU	A	1001	1/1	0.99	0.07	17,17,17,17	0
2	CU	B	701	1/1	1.00	0.06	15,15,15,15	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.