



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 12, 2024 – 01:28 AM EDT

PDB ID : 2QKP  
Title : Crystal structure of C-terminal domain of SMU\_1151c from Streptococcus mutans  
Authors : Ramagopal, U.A.; Toro, R.; Gilmore, M.; Wu, B.; Bain, K.; Gheyi, T.; Sauder, J.M.; Burley, S.K.; Almo, S.C.; New York SGX Research Center for Structural Genomics (NYSGXRC)  
Deposited on : 2007-07-11  
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	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.36.2
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)	:	2.36.2

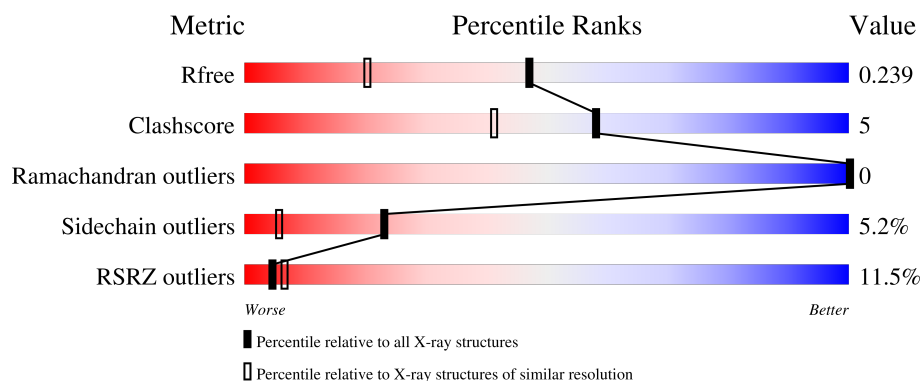
# 1 Overall quality at a glance

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	2340 (1.76-1.76)
Clashscore	141614	2466 (1.76-1.76)
Ramachandran outliers	138981	2437 (1.76-1.76)
Sidechain outliers	138945	2437 (1.76-1.76)
RSRZ outliers	127900	2298 (1.76-1.76)

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	151	<div> <div>11%</div> <div> <div></div> <div>83%</div> <div>8%</div> <div>9%</div> </div> </div>
1	B	151	<div> <div>15%</div> <div> <div></div> <div>74%</div> <div>16%</div> <div>9%</div> </div> </div>
1	C	151	<div> <div>8%</div> <div> <div></div> <div>82%</div> <div>8%</div> <div>9%</div> </div> </div>
1	D	151	<div> <div>8%</div> <div> <div></div> <div>78%</div> <div>10%</div> <div>11%</div> </div> </div>

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
3	GOL	D	601	-	-	X	-

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5020 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 Uncharacterized protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	138	Total	C	N	O	S	0	3	0
			1154	739	196	216	3			
1	B	138	Total	C	N	O	S	0	3	0
			1160	746	196	214	4			
1	C	138	Total	C	N	O	S	0	0	0
			1135	731	191	210	3			
1	D	134	Total	C	N	O	S	0	2	0
			1120	718	192	207	3			

There are 44 discrepancies between the modelled and reference sequences:

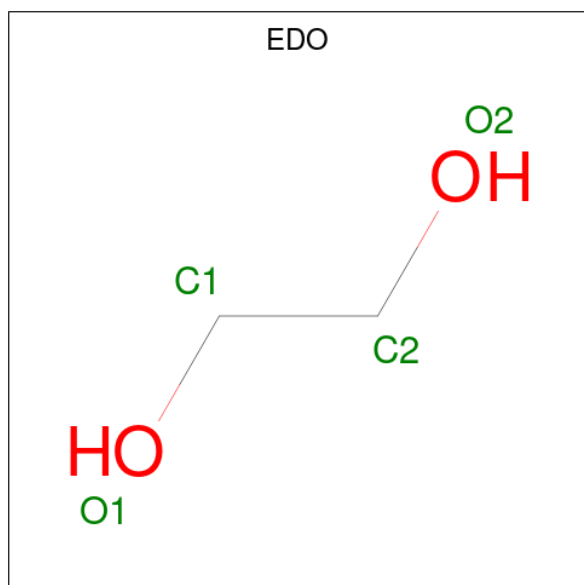
Chain	Residue	Modelled	Actual	Comment	Reference
A	312	MET	-	cloning artifact	UNP Q8DU08
A	313	SER	-	cloning artifact	UNP Q8DU08
A	314	LEU	-	cloning artifact	UNP Q8DU08
A	455	GLU	-	cloning artifact	UNP Q8DU08
A	456	GLY	-	cloning artifact	UNP Q8DU08
A	457	HIS	-	cloning artifact	UNP Q8DU08
A	458	HIS	-	cloning artifact	UNP Q8DU08
A	459	HIS	-	cloning artifact	UNP Q8DU08
A	460	HIS	-	cloning artifact	UNP Q8DU08
A	461	HIS	-	cloning artifact	UNP Q8DU08
A	462	HIS	-	cloning artifact	UNP Q8DU08
B	312	MET	-	cloning artifact	UNP Q8DU08
B	313	SER	-	cloning artifact	UNP Q8DU08
B	314	LEU	-	cloning artifact	UNP Q8DU08
B	455	GLU	-	cloning artifact	UNP Q8DU08
B	456	GLY	-	cloning artifact	UNP Q8DU08
B	457	HIS	-	cloning artifact	UNP Q8DU08
B	458	HIS	-	cloning artifact	UNP Q8DU08
B	459	HIS	-	cloning artifact	UNP Q8DU08
B	460	HIS	-	cloning artifact	UNP Q8DU08
B	461	HIS	-	cloning artifact	UNP Q8DU08

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Chain	Residue	Modelled	Actual	Comment	Reference
B	462	HIS	-	cloning artifact	UNP Q8DU08
C	312	MET	-	cloning artifact	UNP Q8DU08
C	313	SER	-	cloning artifact	UNP Q8DU08
C	314	LEU	-	cloning artifact	UNP Q8DU08
C	455	GLU	-	cloning artifact	UNP Q8DU08
C	456	GLY	-	cloning artifact	UNP Q8DU08
C	457	HIS	-	cloning artifact	UNP Q8DU08
C	458	HIS	-	cloning artifact	UNP Q8DU08
C	459	HIS	-	cloning artifact	UNP Q8DU08
C	460	HIS	-	cloning artifact	UNP Q8DU08
C	461	HIS	-	cloning artifact	UNP Q8DU08
C	462	HIS	-	cloning artifact	UNP Q8DU08
D	312	MET	-	cloning artifact	UNP Q8DU08
D	313	SER	-	cloning artifact	UNP Q8DU08
D	314	LEU	-	cloning artifact	UNP Q8DU08
D	455	GLU	-	cloning artifact	UNP Q8DU08
D	456	GLY	-	cloning artifact	UNP Q8DU08
D	457	HIS	-	cloning artifact	UNP Q8DU08
D	458	HIS	-	cloning artifact	UNP Q8DU08
D	459	HIS	-	cloning artifact	UNP Q8DU08
D	460	HIS	-	cloning artifact	UNP Q8DU08
D	461	HIS	-	cloning artifact	UNP Q8DU08
D	462	HIS	-	cloning artifact	UNP Q8DU08

- Molecule 2 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula:  $C_2H_6O_2$ ).



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

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



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

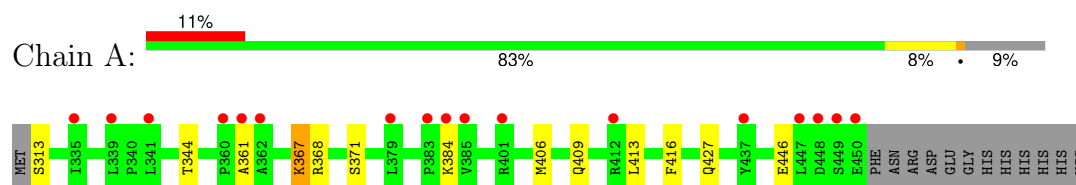
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	122	Total	O	0	0
			122	122		
4	B	97	Total	O	0	0
			97	97		
4	C	116	Total	O	0	0
			116	116		
4	D	102	Total	O	0	0
			102	102		

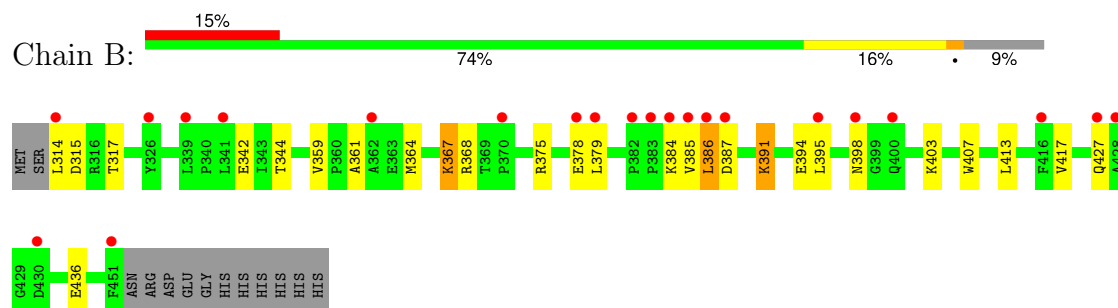
### 3 Residue-property plots [i](#)

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.

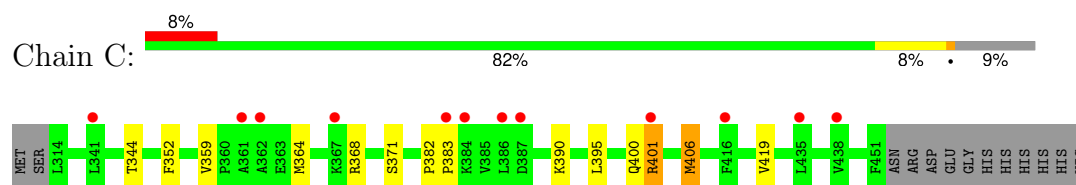
- Molecule 1: Uncharacterized protein



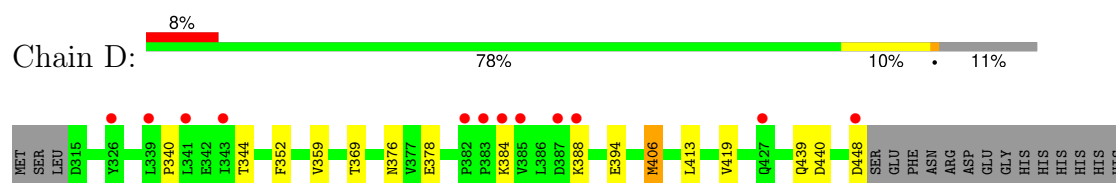
- Molecule 1: Uncharacterized protein



- Molecule 1: Uncharacterized protein



- Molecule 1: Uncharacterized protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	73.00Å 53.91Å 79.92Å 90.00° 106.05° 90.00°	Depositor
Resolution (Å)	29.40 – 1.75 29.40 – 1.75	Depositor EDS
% Data completeness (in resolution range)	99.7 (29.40-1.75) 99.7 (29.40-1.75)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	0.05	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.64 (at 1.75Å)	Xtriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.198 , 0.241 0.197 , 0.239	Depositor DCC
$R_{free}$ test set	3044 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	22.6	Xtriage
Anisotropy	0.067	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 52.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	5020	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 16.15% of the height of the origin peak. No significant pseudotranslation is detected.*

<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: EDO, 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	0.64	0/1180	0.68	0/1596
1	B	0.57	0/1190	0.65	0/1609
1	C	0.62	0/1162	0.70	0/1573
1	D	0.59	0/1146	0.67	0/1551
All	All	0.61	0/4678	0.67	0/6329

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	1154	0	1130	7	0
1	B	1160	0	1138	20	0
1	C	1135	0	1111	10	0
1	D	1120	0	1095	8	0
2	B	4	0	6	0	0
2	C	4	0	6	0	0
3	D	6	0	8	4	0
4	A	122	0	0	1	0
4	B	97	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	116	0	0	1	0
4	D	102	0	0	0	0
All	All	5020	0	4494	44	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 5.

All (44) 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:D:439:GLN:HA	3:D:601:GOL:H12	1.57	0.85
1:B:314:LEU:HD23	1:B:317:THR:HG23	1.68	0.76
1:B:344:THR:HG21	1:B:368:ARG:HE	1.53	0.73
1:C:401:ARG:HG3	1:C:401:ARG:HH11	1.54	0.72
1:B:378:GLU:CB	1:B:386:LEU:HD11	2.22	0.70
1:B:361:ALA:HA	1:B:364[B]:MET:HG3	1.74	0.69
1:C:401:ARG:HG3	1:C:401:ARG:NH1	2.07	0.69
1:D:406:MET:HG2	1:D:419:VAL:HB	1.76	0.67
1:D:340:PRO:O	3:D:601:GOL:H11	1.94	0.67
1:C:395:LEU:HD23	1:C:400:GLN:NE2	2.09	0.67
1:C:344:THR:HG23	1:C:352:PHE:CE2	2.30	0.66
1:C:401:ARG:HH11	1:C:401:ARG:CG	2.10	0.64
1:C:406:MET:HG2	1:C:419:VAL:HB	1.79	0.63
1:B:378:GLU:HB3	1:B:386:LEU:HD11	1.82	0.61
1:B:391:LYS:HE2	1:B:395:LEU:HG	1.83	0.59
1:C:344:THR:HG23	1:C:352:PHE:HE2	1.66	0.59
1:A:409:GLN:HG2	1:A:416:PHE:CE1	2.39	0.58
1:B:391:LYS:HA	1:B:391:LYS:HE3	1.85	0.58
1:C:364:MET:HE1	1:C:368:ARG:H	1.72	0.55
1:B:378:GLU:HB2	1:B:386:LEU:HD11	1.88	0.54
1:D:344:THR:HG23	1:D:352:PHE:CE2	2.43	0.54
1:A:344:THR:HG23	4:A:516:HOH:O	2.07	0.54
1:B:342:GLU:HG2	1:B:364[B]:MET:CE	2.39	0.53
1:D:344:THR:HG23	1:D:352:PHE:HE2	1.73	0.53
1:A:313:SER:HA	1:B:315:ASP:OD1	2.09	0.53
1:A:361:ALA:O	1:A:367:LYS:HE3	2.09	0.52
1:B:375:ARG:HD3	1:B:379:LEU:HD22	1.94	0.49
1:C:390:LYS:NZ	4:C:706:HOH:O	2.46	0.47
1:A:409:GLN:HG2	1:A:416:PHE:HE1	1.78	0.46
1:B:359:VAL:HB	1:B:364[B]:MET:CE	2.45	0.46
1:B:407:TRP:HA	1:B:417:VAL:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:376:ASN:OD1	1:D:378:GLU:HG3	2.16	0.46
1:D:440:ASP:H	3:D:601:GOL:C1	2.28	0.45
1:A:344:THR:HG21	1:A:368:ARG:HE	1.82	0.43
1:B:344:THR:OG1	1:B:436:GLU:HB3	2.19	0.43
1:B:394:GLU:HG2	1:B:398:ASN:HD21	1.84	0.43
1:D:440:ASP:H	3:D:601:GOL:H12	1.84	0.42
1:A:427[A]:GLN:HA	1:A:427[A]:GLN:OE1	2.18	0.42
1:B:391:LYS:HE3	1:B:391:LYS:CA	2.48	0.42
1:B:314:LEU:HD23	1:B:317:THR:CG2	2.43	0.42
1:B:359:VAL:HB	1:B:364[B]:MET:HE1	2.01	0.42
1:B:367:LYS:HE3	1:B:367:LYS:HA	2.00	0.41
1:C:382:PRO:HA	1:C:383:PRO:HD3	1.91	0.41
1:B:403:LYS:HE3	1:B:403:LYS:HB2	1.81	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	139/151 (92%)	135 (97%)	4 (3%)	0	100	100
1	B	139/151 (92%)	136 (98%)	3 (2%)	0	100	100
1	C	136/151 (90%)	134 (98%)	2 (2%)	0	100	100
1	D	134/151 (89%)	131 (98%)	3 (2%)	0	100	100
All	All	548/604 (91%)	536 (98%)	12 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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	128/137 (93%)	122 (95%)	6 (5%)	26	7
1	B	128/137 (93%)	120 (94%)	8 (6%)	18	3
1	C	125/137 (91%)	121 (97%)	4 (3%)	39	16
1	D	123/137 (90%)	115 (94%)	8 (6%)	17	3
All	All	504/548 (92%)	478 (95%)	26 (5%)	23	6

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	367	LYS
1	A	371	SER
1	A	384	LYS
1	A	406	MET
1	A	413	LEU
1	A	446	GLU
1	B	367	LYS
1	B	384	LYS
1	B	385	VAL
1	B	386	LEU
1	B	387	ASP
1	B	391	LYS
1	B	413	LEU
1	B	427	GLN
1	C	359	VAL
1	C	371	SER
1	C	401	ARG
1	C	406	MET
1	D	359	VAL
1	D	369	THR
1	D	384	LYS
1	D	388	LYS
1	D	394	GLU
1	D	406	MET

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Mol	Chain	Res	Type
1	D	413	LEU
1	D	448	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	320	GLN
1	A	405	ASN
1	A	409	GLN
1	B	320	GLN
1	B	398	ASN
1	B	405	ASN
1	C	400	GLN
1	D	405	ASN

### 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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

3 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	602	-	3,3,3	0.43	0	2,2,2	0.66	0
3	GOL	D	601	-	5,5,5	0.72	0	5,5,5	1.25	1 (20%)
2	EDO	B	603	-	3,3,3	0.37	0	2,2,2	0.63	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	602	-	-	0/1/1/1	-
3	GOL	D	601	-	-	2/4/4/4	-
2	EDO	B	603	-	-	0/1/1/1	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	601	GOL	O1-C1-C2	2.61	122.11	110.38

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	601	GOL	O1-C1-C2-C3
3	D	601	GOL	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	601	GOL	4	0

## 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	138/151 (91%)	0.54	17 (12%) <b>4</b> <b>6</b>	13, 22, 43, 56	0
1	B	138/151 (91%)	0.78	22 (15%) <b>1</b> <b>3</b>	13, 28, 49, 51	0
1	C	138/151 (91%)	0.49	12 (8%) <b>10</b> <b>13</b>	13, 23, 39, 45	0
1	D	134/151 (88%)	0.59	12 (8%) <b>9</b> <b>12</b>	13, 27, 43, 52	0
All	All	548/604 (90%)	0.60	63 (11%) <b>4</b> <b>6</b>	13, 25, 43, 56	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	385	VAL	4.7
1	B	386	LEU	4.5
1	D	384	LYS	4.4
1	D	382	PRO	4.1
1	A	449	SER	4.0
1	A	361	ALA	3.7
1	B	385	VAL	3.6
1	D	383	PRO	3.6
1	C	383	PRO	3.6
1	B	384	LYS	3.5
1	D	427	GLN	3.5
1	A	362	ALA	3.3
1	B	326	TYR	3.2
1	B	451	PHE	3.2
1	A	448	ASP	3.2
1	B	416	PHE	3.0
1	B	378	GLU	3.0
1	C	386	LEU	3.0
1	A	450	GLU	2.9
1	A	339	LEU	2.9
1	B	382	PRO	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	385	VAL	2.8
1	B	339	LEU	2.8
1	B	398	ASN	2.8
1	C	341	LEU	2.8
1	A	384	LYS	2.8
1	D	339	LEU	2.8
1	A	412	ARG	2.7
1	D	343	ILE	2.7
1	C	362	ALA	2.7
1	B	427	GLN	2.7
1	A	383	PRO	2.7
1	C	401	ARG	2.6
1	A	379	LEU	2.6
1	C	361	ALA	2.6
1	C	384	LYS	2.6
1	B	387	ASP	2.5
1	D	387[A]	ASP	2.5
1	A	341	LEU	2.5
1	A	335	ILE	2.5
1	A	447	LEU	2.5
1	B	314	LEU	2.5
1	D	326	TYR	2.4
1	C	367	LYS	2.4
1	D	341	LEU	2.4
1	B	341	LEU	2.3
1	C	435	LEU	2.3
1	B	379	LEU	2.3
1	B	395	LEU	2.2
1	B	428	ALA	2.2
1	C	387	ASP	2.2
1	C	438	VAL	2.1
1	A	437	TYR	2.1
1	B	430	ASP	2.1
1	A	360	PRO	2.1
1	A	401	ARG	2.1
1	B	400	GLN	2.1
1	D	388	LYS	2.1
1	B	370	PRO	2.0
1	B	362	ALA	2.0
1	C	416	PHE	2.0
1	D	448	ASP	2.0
1	B	383	PRO	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 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( $\text{\AA}^2$ )	Q<0.9
2	EDO	C	602	4/4	0.86	0.14	29,29,30,36	0
3	GOL	D	601	6/6	0.91	0.12	24,28,30,35	0
2	EDO	B	603	4/4	0.93	0.12	26,27,27,32	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.