



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 31, 2023 – 12:41 PM EDT

PDB ID : 3STF  
Title : Crystal structure of a mutant (S211A) of 3-deoxy-D-manno-octulosonate 8-phosphate synthase (KDO8PS) from *Neisseria meningitidis*  
Authors : Allison, T.M.; Jameson, G.B.; Gloyne, B.J.; Parker, E.J.  
Deposited on : 2011-07-09  
Resolution : 1.90 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

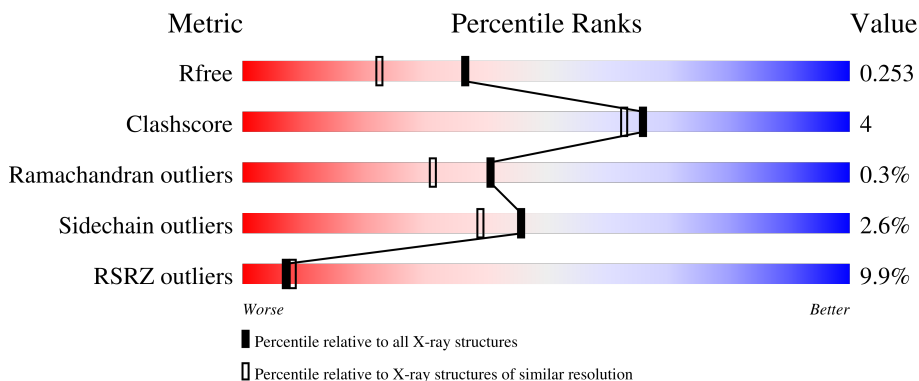
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.90 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	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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	280	 15% 80% 9% • 11%
1	B	280	 4% 79% 11% 10%
1	C	280	 4% 80% 9% • 10%
1	D	280	 13% 80% 9% • 10%

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8327 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 2-dehydro-3-deoxyphosphooctonate aldolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	250	Total 1926	C 1236	N 326	O 353	S 11	0	1	0
1	B	251	Total 1942	C 1250	N 327	O 354	S 11	0	3	0
1	C	252	Total 1950	C 1254	N 328	O 357	S 11	0	3	0
1	D	251	Total 1915	C 1225	N 321	O 358	S 11	0	4	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	211	ALA	SER	ENGINEERED MUTATION	UNP Q9JZ55
B	211	ALA	SER	ENGINEERED MUTATION	UNP Q9JZ55
C	211	ALA	SER	ENGINEERED MUTATION	UNP Q9JZ55
D	211	ALA	SER	ENGINEERED MUTATION	UNP Q9JZ55

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

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

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



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

- Molecule 4 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	C	1	Total Na 1 1	0	0

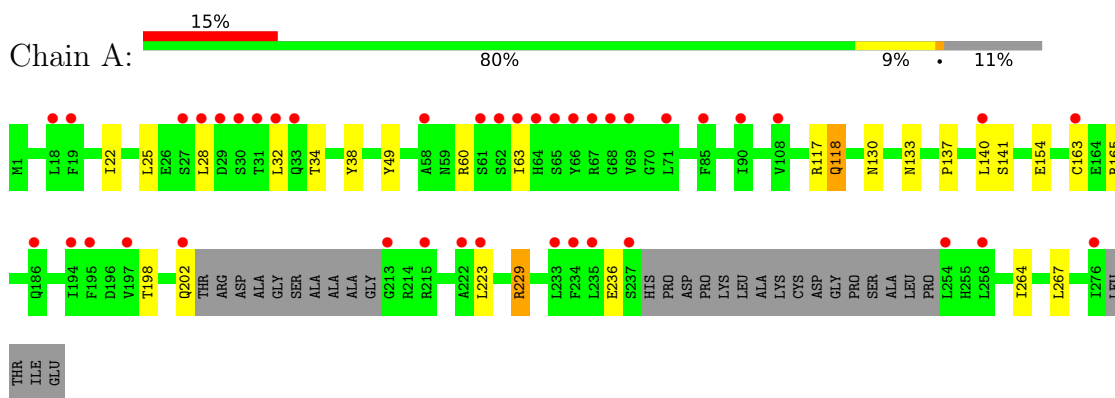
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	114	Total O 114 114	0	0
5	B	173	Total O 173 173	0	0
5	C	158	Total O 158 158	0	0
5	D	141	Total O 141 141	0	0

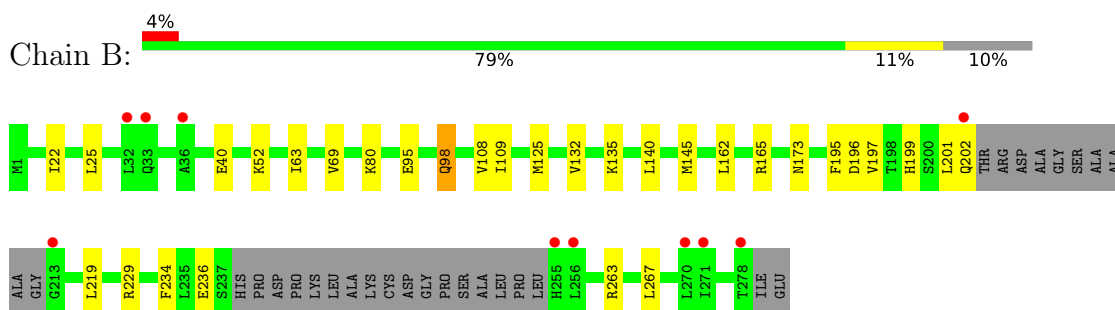
### 3 Residue-property plots [i](#)

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.

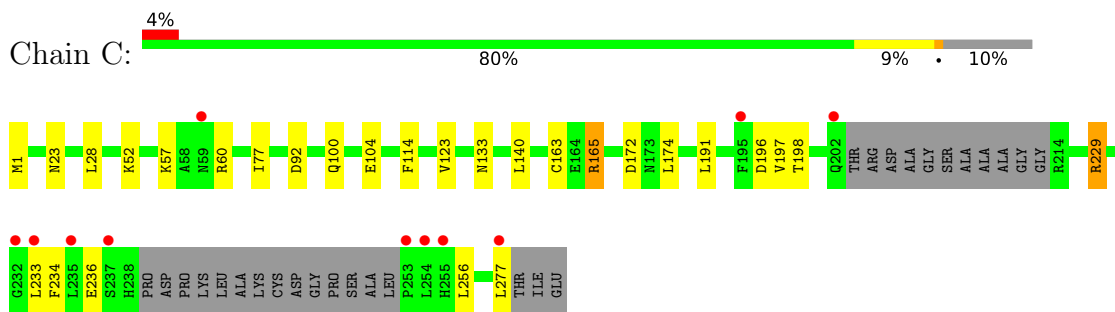
- Molecule 1: 2-dehydro-3-deoxyphosphooctonate aldolase



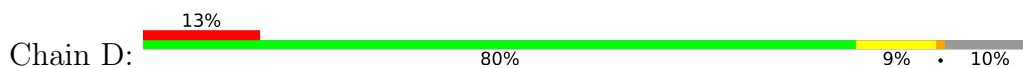
- Molecule 1: 2-dehydro-3-deoxyphosphooctonate aldolase

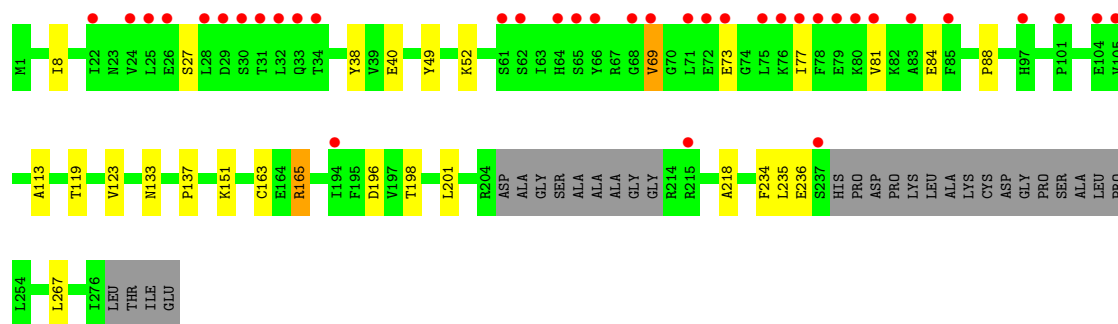


- Molecule 1: 2-dehydro-3-deoxyphosphooctonate aldolase



- Molecule 1: 2-dehydro-3-deoxyphosphooctonate aldolase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	82.19Å 85.83Å 163.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.87 – 1.90 36.87 – 1.90	Depositor EDS
% Data completeness (in resolution range)	93.7 (36.87-1.90) 93.8 (36.87-1.90)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.68 (at 1.89Å)	Xtrriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.218 , 0.258 0.211 , 0.253	Depositor DCC
$R_{free}$ test set	4322 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	33.1	Xtrriage
Anisotropy	0.403	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 41.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	0.022 for k,h,-l	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	8327	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	41.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.64% 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: NA, GOL, 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	0.52	0/1958	0.62	1/2645 (0.0%)
1	B	0.58	0/1980	0.65	0/2676
1	C	0.56	0/1989	0.65	1/2688 (0.0%)
1	D	0.53	0/1955	0.63	0/2648
All	All	0.55	0/7882	0.64	2/10657 (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	229	ARG	NE-CZ-NH2	-5.41	117.60	120.30
1	C	229	ARG	NE-CZ-NH2	-5.16	117.72	120.30

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	1926	0	1962	15	0
1	B	1942	0	1997	24	0
1	C	1950	0	1997	18	0
1	D	1915	0	1918	15	0
2	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	6	0	8	0	0
4	C	1	0	0	0	0
5	A	114	0	0	0	0
5	B	173	0	0	1	0
5	C	158	0	0	2	0
5	D	141	0	0	1	0
All	All	8327	0	7882	63	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 (63) 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:125:MET:HE1	1:B:132:VAL:HG13	1.62	0.82
1:B:22:ILE:HD11	1:B:25:LEU:HD23	1.64	0.78
1:A:267:LEU:HD13	1:B:267[B]:LEU:HD21	1.65	0.78
1:D:69:VAL:HG13	1:D:73:GLU:HB2	1.71	0.72
1:A:22:ILE:HD11	1:A:25:LEU:HD23	1.75	0.67
1:A:133:ASN:HD21	1:A:163:CYS:HB2	1.59	0.67
1:D:201:LEU:HD11	1:D:218:ALA:HA	1.79	0.63
1:B:201:LEU:O	1:B:202:GLN:CB	2.47	0.62
1:D:133:ASN:HD21	1:D:163:CYS:HB2	1.66	0.61
1:B:63:ILE:HD12	1:D:123:VAL:CG2	2.33	0.59
1:B:63:ILE:HD13	1:D:119:THR:HG23	1.84	0.59
1:B:135:LYS:HD2	1:B:165:ARG:HD2	1.84	0.58
1:A:267:LEU:CD1	1:B:267[B]:LEU:HD21	2.35	0.57
1:D:151:LYS:NZ	5:D:329:HOH:O	2.32	0.56
1:B:109:ILE:HG22	1:B:125:MET:HE3	1.87	0.56
1:C:165:ARG:O	1:C:165:ARG:NH1	2.39	0.55
1:A:137:PRO:HG2	1:A:140:LEU:HD12	1.89	0.55
1:A:267:LEU:HD13	1:B:267[B]:LEU:CD2	2.36	0.54
1:C:133:ASN:HD21	1:C:163:CYS:HB2	1.74	0.53
1:A:198:THR:HG21	1:A:236:GLU:HG3	1.92	0.51
1:D:38:TYR:HB3	1:D:49:TYR:CZ	2.46	0.51
1:B:219:LEU:HD21	1:B:263:ARG:HD3	1.93	0.51
1:B:135:LYS:HE3	5:B:425:HOH:O	2.11	0.50
1:C:1:MET:HE2	1:C:191:LEU:HD23	1.94	0.50
1:D:267:LEU:C	1:D:267:LEU:HD23	2.31	0.50
1:C:1:MET:CE	1:C:191:LEU:HD23	2.41	0.50
1:C:52:LYS:HD2	1:C:234:PHE:CZ	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:197[B]:VAL:CG1	1:C:233:LEU:HD11	2.42	0.49
1:C:100:GLN:O	1:C:104:GLU:HG3	2.13	0.49
1:C:197[B]:VAL:HG11	1:C:233:LEU:HD11	1.95	0.48
1:A:223:LEU:HD23	1:A:264:ILE:HD13	1.96	0.47
1:C:196:ASP:HA	1:C:234:PHE:HB3	1.96	0.47
1:A:117:ARG:NH2	1:A:140:LEU:HD21	2.30	0.47
1:C:172:ASP:OD2	5:C:465:HOH:O	2.21	0.46
1:D:165:ARG:O	1:D:165:ARG:HD3	2.15	0.46
1:B:199:HIS:HB2	1:B:236:GLU:OE2	2.16	0.46
1:B:52:LYS:HE2	1:B:234:PHE:CZ	2.51	0.45
1:B:196:ASP:HA	1:B:234:PHE:HB3	1.99	0.45
1:D:8:ILE:HD12	1:D:88:PRO:HG2	1.99	0.44
1:D:196:ASP:HA	1:D:234:PHE:HB3	1.99	0.44
1:D:198:THR:HG21	1:D:236:GLU:HB2	2.00	0.44
1:B:145:MET:CE	1:B:162:LEU:HD22	2.48	0.44
1:C:23:ASN:OD1	1:C:57:LYS:CE	2.66	0.44
1:A:28:LEU:HD11	1:A:32:LEU:HD11	2.00	0.44
1:C:277:LEU:HA	5:C:490:HOH:O	2.18	0.43
1:A:141:SER:HB3	1:C:172:ASP:OD2	2.19	0.43
1:B:145:MET:HE3	1:B:162:LEU:HD22	2.00	0.43
1:C:198:THR:HG21	1:C:236:GLU:HB2	2.00	0.43
1:A:38:TYR:HB3	1:A:49:TYR:CZ	2.54	0.43
1:B:63:ILE:HD12	1:D:123:VAL:HG22	2.01	0.42
1:D:77:ILE:O	1:D:81:VAL:HG23	2.20	0.42
1:D:113:ALA:O	1:D:137:PRO:HG3	2.20	0.42
1:B:195:PHE:CE2	1:B:197[B]:VAL:HG22	2.54	0.42
1:B:140:LEU:HD12	1:B:140:LEU:HA	1.94	0.42
1:B:95:GLU:O	1:B:98:GLN:HG3	2.19	0.42
1:B:108:VAL:O	1:B:109:ILE:HD13	2.20	0.41
1:A:22:ILE:HG23	1:A:34:THR:HG21	2.03	0.41
1:C:28:LEU:HD13	1:C:77:ILE:HG12	2.03	0.41
1:C:165:ARG:HH12	1:C:174:LEU:HD13	1.86	0.41
1:A:118:GLN:HG3	1:C:114:PHE:CE1	2.56	0.40
1:A:63:ILE:HD12	1:C:123:VAL:CG2	2.52	0.40
1:B:25:LEU:O	1:B:69:VAL:HG13	2.21	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	245/280 (88%)	239 (98%)	5 (2%)	1 (0%)	34	24
1	B	248/280 (89%)	241 (97%)	6 (2%)	1 (0%)	34	24
1	C	249/280 (89%)	245 (98%)	3 (1%)	1 (0%)	34	24
1	D	249/280 (89%)	245 (98%)	4 (2%)	0	100	100
All	All	991/1120 (88%)	970 (98%)	18 (2%)	3 (0%)	41	31

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	229	ARG
1	B	229	ARG
1	A	229	ARG

### 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	211/235 (90%)	205 (97%)	6 (3%)	43	36
1	B	215/235 (92%)	211 (98%)	4 (2%)	57	53
1	C	216/235 (92%)	211 (98%)	5 (2%)	50	45
1	D	209/235 (89%)	202 (97%)	7 (3%)	38	29
All	All	851/940 (90%)	829 (97%)	22 (3%)	46	39

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	60	ARG
1	A	118	GLN
1	A	130	ASN
1	A	154	GLU
1	A	165	ARG
1	A	202	GLN
1	B	40	GLU
1	B	80	LYS
1	B	98	GLN
1	B	173	ASN
1	C	60	ARG
1	C	92	ASP
1	C	140	LEU
1	C	165	ARG
1	C	256	LEU
1	D	27	SER
1	D	40	GLU
1	D	52	LYS
1	D	69	VAL
1	D	84	GLU
1	D	165	ARG
1	D	235	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	100	GLN
1	A	110	GLN
1	A	130	ASN
1	A	133	ASN
1	B	138	GLN
1	B	144	GLN
1	B	147	ASN
1	C	97	HIS
1	C	100	GLN
1	C	110	GLN
1	C	133	ASN
1	D	59	ASN
1	D	110	GLN
1	D	133	ASN
1	D	144	GLN
1	D	147	ASN
1	D	186	GLN

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Mol	Chain	Res	Type
1	D	202	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](#)

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](#)

Of 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 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	GOL	B	282	-	5,5,5	0.23	0	5,5,5	0.76	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	GOL	B	282	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	282	GOL	C1-C2-C3-O3
3	B	282	GOL	O2-C2-C3-O3

There are no ring outliers.

No monomer is involved in short contacts.

## 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	250/280 (89%)	0.83	41 (16%) <b>1</b> <b>1</b>	23, 40, 73, 100	1 (0%)
1	B	251/280 (89%)	0.28	10 (3%) 38 41	23, 33, 50, 90	1 (0%)
1	C	252/280 (90%)	0.29	11 (4%) 34 37	25, 34, 54, 91	0
1	D	251/280 (89%)	0.72	37 (14%) <b>2</b> <b>2</b>	24, 44, 86, 122	0
All	All	1004/1120 (89%)	0.53	99 (9%) <b>7</b> <b>8</b>	23, 37, 75, 122	2 (0%)

All (99) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	66	TYR	9.0
1	C	255	HIS	8.6
1	D	61	SER	8.5
1	A	235	LEU	5.7
1	D	24	VAL	5.6
1	A	215	ARG	5.5
1	C	202	GLN	5.4
1	C	254	LEU	5.1
1	A	64	HIS	5.0
1	A	61	SER	5.0
1	A	213	GLY	4.9
1	D	68	GLY	4.7
1	B	278	THR	4.7
1	C	253	PRO	4.7
1	A	69	VAL	4.7
1	D	62	SER	4.6
1	C	277	LEU	4.5
1	A	237	SER	4.4
1	D	65	SER	4.2
1	B	213	GLY	4.2
1	D	30[A]	SER	4.0

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	D	33	GLN	4.0
1	B	255	HIS	3.9
1	A	67	ARG	3.9
1	D	29	ASP	3.7
1	A	18	LEU	3.6
1	A	276	ILE	3.6
1	D	26	GLU	3.6
1	A	19	PHE	3.5
1	A	63	ILE	3.5
1	A	30	SER	3.5
1	A	90	ILE	3.5
1	A	202	GLN	3.5
1	D	31	THR	3.4
1	D	71	LEU	3.2
1	D	83	ALA	3.2
1	A	254	LEU	3.1
1	A	234	PHE	3.1
1	A	33	GLN	3.1
1	D	28	LEU	3.0
1	D	104	GLU	3.0
1	D	32	LEU	3.0
1	C	237	SER	3.0
1	A	68	GLY	3.0
1	A	29	ASP	3.0
1	D	97	HIS	3.0
1	A	195	PHE	2.9
1	B	36	ALA	2.9
1	D	81	VAL	2.8
1	D	79	GLU	2.8
1	D	34	THR	2.8
1	A	197	VAL	2.7
1	A	32	LEU	2.7
1	D	73	GLU	2.7
1	D	22	ILE	2.6
1	A	62	SER	2.6
1	D	237	SER	2.6
1	C	233	LEU	2.6
1	D	76	LYS	2.6
1	A	65	SER	2.6
1	A	27	SER	2.6
1	D	105	VAL	2.6
1	B	271	ILE	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	108	VAL	2.5
1	A	71	LEU	2.5
1	A	233	LEU	2.5
1	B	33	GLN	2.5
1	D	69	VAL	2.5
1	B	32	LEU	2.5
1	D	75	LEU	2.5
1	C	59	ASN	2.5
1	C	235	LEU	2.4
1	D	66	TYR	2.4
1	D	215	ARG	2.4
1	D	77	ILE	2.4
1	A	28	LEU	2.3
1	A	256	LEU	2.3
1	D	25	LEU	2.3
1	A	31	THR	2.3
1	A	194	ILE	2.3
1	A	186	GLN	2.3
1	C	232	GLY	2.3
1	A	223	LEU	2.3
1	B	256	LEU	2.3
1	D	64	HIS	2.3
1	D	72	GLU	2.2
1	C	195	PHE	2.2
1	D	85	PHE	2.2
1	B	270	LEU	2.2
1	A	222	ALA	2.1
1	A	58	ALA	2.1
1	A	85	PHE	2.1
1	D	101	PRO	2.1
1	D	78	PHE	2.1
1	B	202	GLN	2.0
1	D	80	LYS	2.0
1	A	140	LEU	2.0
1	D	194	ILE	2.0
1	A	163	CYS	2.0

## 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 [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
3	GOL	B	282	6/6	0.91	0.12	37,44,48,55	0
4	NA	C	281	1/1	0.92	0.09	32,32,32,32	0
2	CL	B	281	1/1	0.98	0.06	43,43,43,43	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.