



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:20 PM UTC

PDB ID : 9G7P / pdb_00009g7p
Title : Crystal structure of Shewanella putrefaciens PE-like toxin, Spx
Authors : Masuyer, G.
Deposited on : 2024-07-22
Resolution : 1.86 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

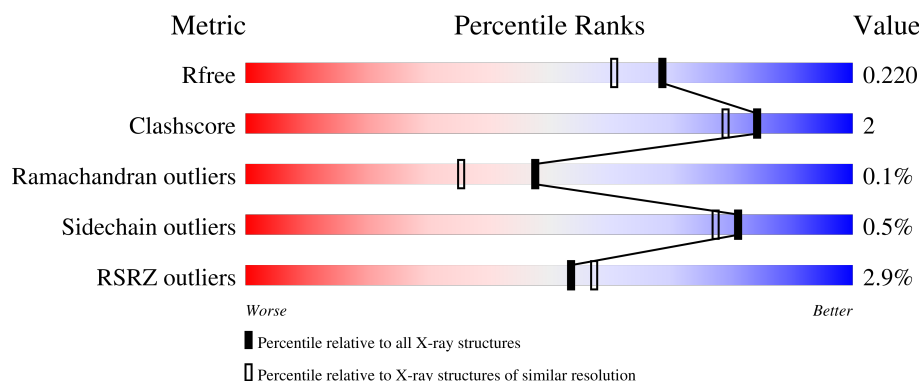
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.86 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	3428 (1.86-1.86)
Clashscore	190562	3579 (1.86-1.86)
Ramachandran outliers	187476	3553 (1.86-1.86)
Sidechain outliers	187428	3553 (1.86-1.86)
RSRZ outliers	180081	3429 (1.86-1.86)

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	AAA	613	2% 93% 5%
1	BBB	613	3% 91% 8%
1	CCC	613	3% 92% 6%

2 Entry composition [i](#)

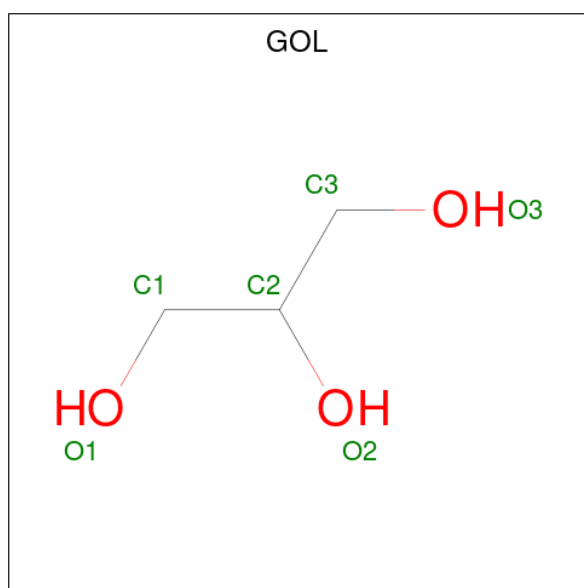
There are 4 unique types of molecules in this entry. The entry contains 28572 atoms, of which 13386 are hydrogens and 0 are deuteriums.

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 *Shewanella putrefaciens* PE-like toxin catalytically inactive mutant.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	AAA	603	Total	C	H	N	O	S	156	0	0
			9081	2897	4446	819	908	11			
1	BBB	603	Total	C	H	N	O	S	156	0	0
			9081	2897	4446	819	908	11			
1	CCC	603	Total	C	H	N	O	S	156	0	0
			9081	2897	4446	819	908	11			

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	AAA	1	Total	C	H	O	2	0
			14	3	8	3		
2	AAA	1	Total	C	H	O	2	0
			14	3	8	3		
2	BBB	1	Total	C	H	O	2	0
			14	3	8	3		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	BBB	1	Total	C	H	O	2	0
			14	3	8	3		
2	CCC	1	Total	C	H	O	2	0
			14	3	8	3		
2	CCC	1	Total	C	H	O	2	0
			14	3	8	3		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	AAA	1	Total	Mg	0	0
			1	1		
3	BBB	2	Total	Mg	0	0
			2	2		
3	CCC	1	Total	Mg	0	0
			1	1		

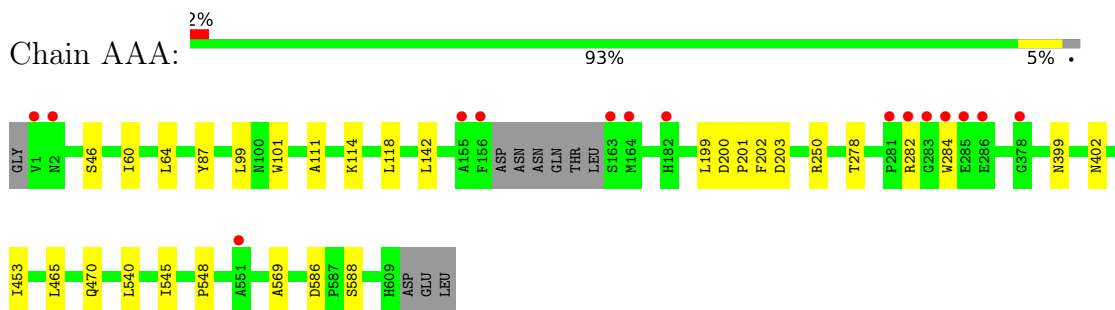
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	AAA	400	Total	O	0	0
			400	400		
4	BBB	380	Total	O	0	0
			380	380		
4	CCC	461	Total	O	0	0
			461	461		

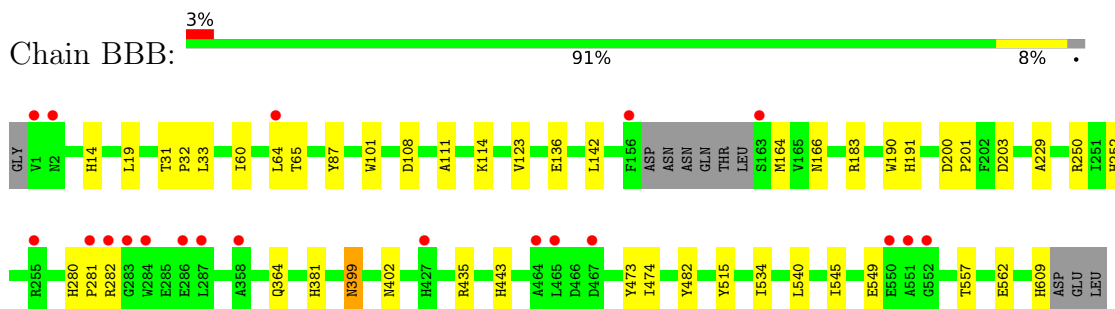
3 Residue-property plots [i](#)

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.

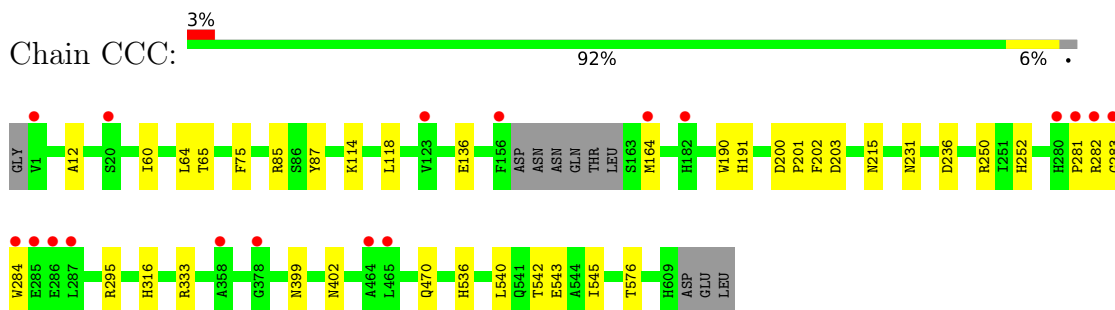
- Molecule 1: *Shewanella putrefaciens* PE-like toxin catalytically inactive mutant



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- Molecule 1: *Shewanella putrefaciens* PE-like toxin catalytically inactive mutant



4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α , β , γ	121.20Å 121.20Å 125.22Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	104.83 – 1.86 104.83 – 1.86	Depositor EDS
% Data completeness (in resolution range)	99.7 (104.83-1.86) 99.5 (104.83-1.86)	Depositor EDS
R_{merge}	0.28	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.10 (at 1.86Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.185 , 0.220 0.184 , 0.220	Depositor DCC
R_{free} test set	8479 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	31.0	Xtriage
Anisotropy	0.235	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 32.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.022 for -h,-k,l 0.059 for h,-h-k,-l 0.032 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	28572	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.52% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, MG

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	AAA	0.98	0/4750	1.27	2/6495 (0.0%)
1	BBB	0.98	0/4750	1.29	4/6495 (0.1%)
1	CCC	0.98	0/4750	1.27	1/6495 (0.0%)
All	All	0.98	0/14250	1.28	7/19485 (0.0%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AAA	548	PRO	CA-C-N	5.61	127.80	120.28
1	AAA	548	PRO	C-N-CA	5.61	127.80	120.28
1	BBB	399	ASN	CA-C-N	5.06	125.70	120.03
1	BBB	399	ASN	C-N-CA	5.06	125.70	120.03
1	CCC	333	ARG	CG-CD-NE	-5.04	100.91	112.00
1	BBB	19	LEU	CA-C-N	5.03	127.02	120.28
1	BBB	19	LEU	C-N-CA	5.03	127.02	120.28

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	AAA	4635	4446	4424	15	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	BBB	4635	4446	4424	25	1
1	CCC	4635	4446	4424	22	1
2	AAA	12	16	16	0	0
2	BBB	12	16	16	1	0
2	CCC	12	16	16	2	0
3	AAA	1	0	0	0	0
3	BBB	2	0	0	0	1
3	CCC	1	0	0	0	0
4	AAA	400	0	0	1	0
4	BBB	380	0	0	3	0
4	CCC	461	0	0	6	0
All	All	15186	13386	13320	63	2

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 (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:CCC:470:GLN:NE2	4:CCC:804:HOH:O	2.29	0.62
1:CCC:215:ASN:HB2	4:CCC:1183:HOH:O	1.98	0.62
1:BBB:399:ASN:H	1:BBB:402:ASN:HD22	1.47	0.62
1:CCC:399:ASN:H	1:CCC:402:ASN:HD22	1.50	0.59
1:AAA:470:GLN:NE2	4:AAA:801:HOH:O	2.22	0.56
1:BBB:482:TYR:O	2:BBB:701:GOL:O1	2.18	0.56
1:CCC:60:ILE:HD12	1:CCC:64:LEU:HD23	1.87	0.56
1:AAA:278:THR:HG21	1:AAA:282:ARG:HG2	1.88	0.56
1:CCC:236:ASP:OD2	4:CCC:801:HOH:O	2.18	0.56
1:AAA:399:ASN:H	1:AAA:402:ASN:HD22	1.55	0.55
1:CCC:136:GLU:H	1:CCC:252:HIS:HD2	1.55	0.55
1:BBB:114:LYS:HZ1	1:BBB:191:HIS:HD2	1.55	0.55
1:BBB:136:GLU:H	1:BBB:252:HIS:HD2	1.56	0.54
1:CCC:282:ARG:HG2	1:CCC:283:GLY:N	2.23	0.53
1:BBB:381:HIS:HD2	1:BBB:562:GLU:OE2	1.91	0.53
1:CCC:282:ARG:HD3	4:CCC:990:HOH:O	2.08	0.53
1:AAA:200:ASP:N	1:AAA:201:PRO:HD2	2.25	0.52
2:CCC:702:GOL:O1	4:CCC:802:HOH:O	2.19	0.51
1:BBB:31:THR:OG1	1:BBB:32:PRO:HD3	2.10	0.51
1:BBB:14:HIS:CE1	1:BBB:166:ASN:HD22	2.29	0.51
1:BBB:435:ARG:HD2	4:BBB:1122:HOH:O	2.11	0.50
1:CCC:87:TYR:CE1	1:CCC:250:ARG:HD3	2.46	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CCC:12:ALA:O	1:CCC:231:ASN:HB3	2.12	0.50
1:AAA:453:ILE:HG22	1:AAA:569:ALA:HB3	1.93	0.49
1:BBB:200:ASP:N	1:BBB:201:PRO:HD2	2.28	0.49
1:BBB:540:LEU:HD11	1:BBB:545:ILE:HG22	1.95	0.48
1:CCC:118:LEU:HD22	1:CCC:202:PHE:CE1	2.49	0.48
1:BBB:364:GLN:NE2	4:BBB:810:HOH:O	2.47	0.48
1:CCC:85:ARG:HD2	1:CCC:252:HIS:CE1	2.48	0.48
1:BBB:123:VAL:HG12	1:BBB:123:VAL:O	2.13	0.47
1:CCC:200:ASP:N	1:CCC:201:PRO:HD2	2.30	0.47
1:AAA:465:LEU:HD11	1:AAA:470:GLN:HG2	1.96	0.47
1:BBB:280:HIS:O	1:BBB:282:ARG:N	2.48	0.47
1:CCC:536:HIS:HE2	1:CCC:543:GLU:CD	2.24	0.46
1:BBB:33:LEU:HD13	1:BBB:229:ALA:HB2	1.98	0.46
1:CCC:542:THR:C	1:CCC:543:GLU:HG3	2.41	0.46
1:BBB:473:TYR:OH	1:BBB:549:GLU:HG3	2.15	0.46
1:AAA:87:TYR:CE1	1:AAA:250:ARG:HD3	2.51	0.45
1:BBB:443:HIS:HB2	1:BBB:474:ILE:HG22	1.99	0.45
1:BBB:435:ARG:NH1	4:BBB:802:HOH:O	2.34	0.45
1:CCC:282:ARG:CG	1:CCC:283:GLY:N	2.80	0.45
1:BBB:114:LYS:HE2	1:BBB:190:TRP:CE2	2.53	0.44
1:AAA:60:ILE:HD12	1:AAA:64:LEU:HD23	2.00	0.44
1:BBB:474:ILE:HG13	1:BBB:557:THR:HB	1.99	0.44
1:AAA:540:LEU:HD11	1:AAA:545:ILE:HG22	1.99	0.44
1:CCC:295:ARG:HD3	4:CCC:850:HOH:O	2.17	0.44
1:CCC:114:LYS:HE2	1:CCC:190:TRP:CE2	2.54	0.43
1:AAA:118:LEU:HD22	1:AAA:202:PHE:CE1	2.53	0.43
1:BBB:108:ASP:O	1:BBB:183:ARG:NE	2.36	0.43
1:AAA:46:SER:HA	1:AAA:99:LEU:O	2.19	0.43
1:BBB:515:TYR:CD2	1:BBB:534:ILE:HG22	2.53	0.43
1:CCC:540:LEU:HD11	1:CCC:545:ILE:HG22	2.00	0.43
1:AAA:199:LEU:C	1:AAA:201:PRO:HD2	2.44	0.42
1:BBB:111:ALA:HA	1:BBB:142:LEU:HD11	2.01	0.42
1:CCC:114:LYS:HZ1	1:CCC:191:HIS:HD2	1.67	0.42
1:BBB:87:TYR:CE1	1:BBB:250:ARG:HD3	2.55	0.42
1:AAA:101:TRP:HA	1:AAA:114:LYS:O	2.20	0.42
1:AAA:586:ASP:OD1	1:AAA:588:SER:HB3	2.20	0.42
1:BBB:60:ILE:HD12	1:BBB:64:LEU:HD23	2.02	0.42
1:BBB:101:TRP:HA	1:BBB:114:LYS:O	2.20	0.42
1:CCC:576:THR:HB	2:CCC:702:GOL:H32	2.01	0.42
1:CCC:65:THR:O	1:CCC:75:PHE:HA	2.21	0.41
1:AAA:111:ALA:HA	1:AAA:142:LEU:HD11	2.03	0.40

All (2) 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	Interatomic distance (Å)	Clash overlap (Å)
1:BBB:65:THR:HG23	1:BBB:609:HIS:HD1[3_644]	1.31	0.29
1:CCC:316:HIS:HE2	3:BBB:704:MG:MG[1_665]	1.58	0.02

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	AAA	599/613 (98%)	584 (98%)	15 (2%)	0	100	100
1	BBB	599/613 (98%)	582 (97%)	16 (3%)	1 (0%)	43	31
1	CCC	599/613 (98%)	581 (97%)	17 (3%)	1 (0%)	43	31
All	All	1797/1839 (98%)	1747 (97%)	48 (3%)	2 (0%)	48	35

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	BBB	281	PRO
1	CCC	281	PRO

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	AAA	496/505 (98%)	494 (100%)	2 (0%)	84	82

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	BBB	496/505 (98%)	494 (100%)	2 (0%)	84	82
1	CCC	496/505 (98%)	493 (99%)	3 (1%)	78	73
All	All	1488/1515 (98%)	1481 (100%)	7 (0%)	81	77

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AAA	203	ASP
1	AAA	284	TRP
1	BBB	164	MET
1	BBB	203	ASP
1	CCC	164	MET
1	CCC	203	ASP
1	CCC	284	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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](#)

Of 10 ligands modelled in this entry, 4 are monoatomic - leaving 6 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	$\# Z > 2$
2	GOL	AAA	702	-	5,5,5	0.10	0	5,5,5	0.29	0
2	GOL	BBB	702	-	5,5,5	0.09	0	5,5,5	0.26	0
2	GOL	CCC	701	-	5,5,5	0.08	0	5,5,5	0.32	0
2	GOL	CCC	702	-	5,5,5	0.07	0	5,5,5	0.33	0
2	GOL	AAA	701	-	5,5,5	0.09	0	5,5,5	0.27	0
2	GOL	BBB	701	-	5,5,5	0.09	0	5,5,5	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
2	GOL	AAA	702	-	-	2/4/4/4	-
2	GOL	BBB	702	-	-	2/4/4/4	-
2	GOL	CCC	701	-	-	2/4/4/4	-
2	GOL	CCC	702	-	-	2/4/4/4	-
2	GOL	AAA	701	-	-	0/4/4/4	-
2	GOL	BBB	701	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	BBB	701	GOL	O1-C1-C2-C3
2	CCC	701	GOL	O1-C1-C2-C3
2	CCC	702	GOL	O1-C1-C2-O2
2	CCC	702	GOL	O1-C1-C2-C3
2	BBB	701	GOL	O1-C1-C2-O2
2	CCC	701	GOL	O1-C1-C2-O2
2	AAA	702	GOL	O1-C1-C2-C3
2	BBB	702	GOL	O1-C1-C2-C3
2	BBB	702	GOL	O1-C1-C2-O2
2	AAA	702	GOL	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	CCC	702	GOL	2	0
2	BBB	701	GOL	1	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, 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	AAA	603/613 (98%)	-0.01	15 (2%) 58 62	21, 33, 54, 119	0
1	BBB	603/613 (98%)	-0.01	20 (3%) 49 53	23, 34, 55, 119	0
1	CCC	603/613 (98%)	-0.06	18 (2%) 52 56	20, 33, 52, 112	0
All	All	1809/1839 (98%)	-0.03	53 (2%) 53 57	20, 33, 54, 119	0

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	AAA	284	TRP	9.7
1	BBB	284	TRP	8.0
1	CCC	284	TRP	7.9
1	AAA	156	PHE	6.9
1	BBB	281	PRO	5.7
1	BBB	465	LEU	5.4
1	CCC	156	PHE	5.2
1	AAA	281	PRO	5.1
1	CCC	281	PRO	5.1
1	CCC	465	LEU	5.1
1	BBB	156	PHE	5.0
1	BBB	1	VAL	4.8
1	BBB	282	ARG	4.8
1	CCC	283	GLY	4.2
1	AAA	1	VAL	4.2
1	CCC	1	VAL	4.2
1	BBB	283	GLY	3.9
1	BBB	287	LEU	3.7
1	AAA	283	GLY	3.7
1	AAA	282	ARG	3.6
1	CCC	287	LEU	3.4
1	BBB	552	GLY	3.1
1	CCC	282	ARG	3.1

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Mol	Chain	Res	Type	RSRZ
1	BBB	286	GLU	3.1
1	CCC	286	GLU	2.9
1	CCC	358	ALA	2.9
1	AAA	285	GLU	2.8
1	BBB	464	ALA	2.8
1	AAA	551	ALA	2.8
1	CCC	464	ALA	2.7
1	AAA	164	MET	2.7
1	CCC	378	GLY	2.7
1	CCC	182	HIS	2.7
1	AAA	163	SER	2.6
1	BBB	427	HIS	2.5
1	BBB	2	ASN	2.5
1	CCC	123	VAL	2.5
1	AAA	155	ALA	2.4
1	BBB	255	ARG	2.3
1	BBB	358	ALA	2.3
1	BBB	467	ASP	2.3
1	BBB	163	SER	2.3
1	CCC	285	GLU	2.3
1	CCC	164	MET	2.3
1	CCC	20	SER	2.2
1	BBB	64	LEU	2.2
1	CCC	280	HIS	2.2
1	BBB	550	GLU	2.2
1	BBB	551	ALA	2.1
1	AAA	286	GLU	2.1
1	AAA	2	ASN	2.1
1	AAA	182	HIS	2.0
1	AAA	378	GLY	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 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	GOL	AAA	701	6/6	0.82	0.21	41,52,61,61	2
2	GOL	BBB	702	6/6	0.86	0.15	38,51,57,58	2
2	GOL	BBB	701	6/6	0.87	0.14	50,57,64,64	2
2	GOL	CCC	701	6/6	0.90	0.12	36,42,46,47	2
2	GOL	AAA	702	6/6	0.91	0.12	41,48,53,54	2
2	GOL	CCC	702	6/6	0.91	0.12	44,46,52,53	2
3	MG	AAA	703	1/1	0.98	0.04	29,29,29,29	0
3	MG	BBB	704	1/1	0.98	0.06	35,35,35,35	0
3	MG	CCC	703	1/1	0.99	0.02	27,27,27,27	0
3	MG	BBB	703	1/1	1.00	0.02	28,28,28,28	0

6.5 Other polymers [i](#)

There are no such residues in this entry.