



wwPDB X-ray Structure Validation Summary Report ⓘ

Apr 18, 2026 – 05:25 PM UTC

PDB ID : 4N58 / pdb_00004n58
Title : Crystal Structure of Pectocin M2 at 1.86 Angstroms
Authors : Grinter, R.; Roszak, A.W.; Zeth, K.; Cogdell, C.J.; Walker, D.
Deposited on : 2013-10-09
Resolution : 1.86 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

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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

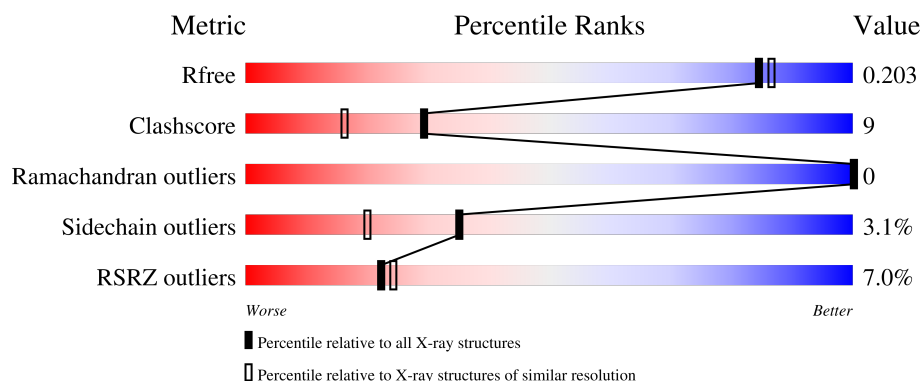
1 Overall quality at a glance

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	A	279	5% 81% 13% .. 5%
1	B	279	8% 78% 15% .. 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	GOL	A	312	-	X	-	-
4	GOL	A	322	-	-	X	-
4	GOL	B	314	-	-	X	-
5	MPD	A	325	-	X	-	-

2 Entry composition [i](#)

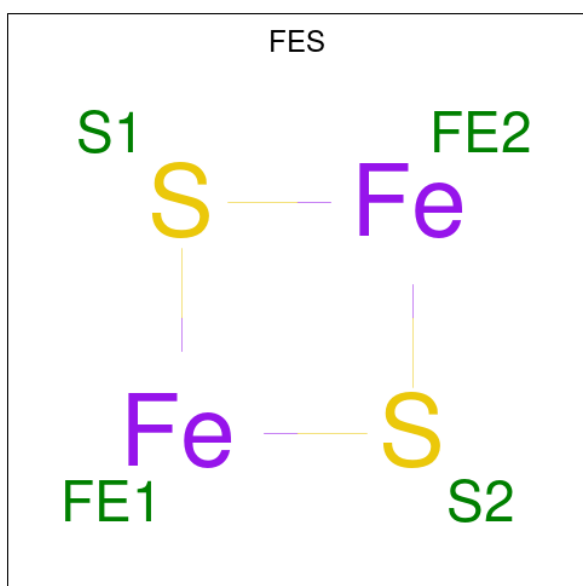
There are 7 unique types of molecules in this entry. The entry contains 5000 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Pectocin M2.

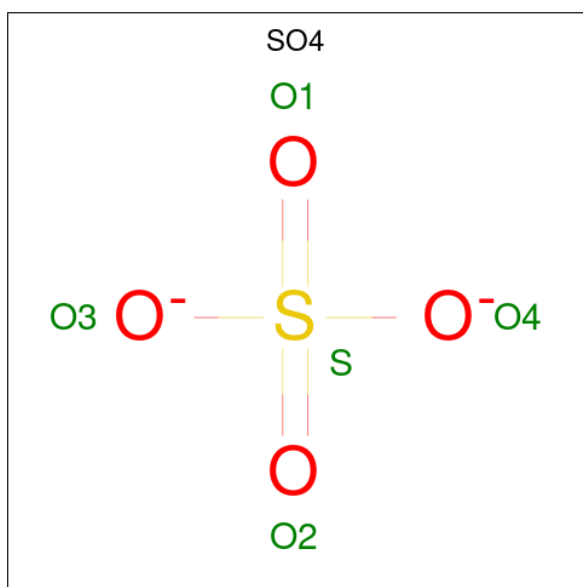
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	266	Total	C	N	O	S	0	5	0
			2088	1313	344	423	8			
1	B	266	Total	C	N	O	S	0	4	0
			2085	1309	343	424	9			

- Molecule 2 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe_2S_2).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	Fe	S	0	0
			4	2	2		
2	B	1	Total	Fe	S	0	0
			4	2	2		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O_4S).



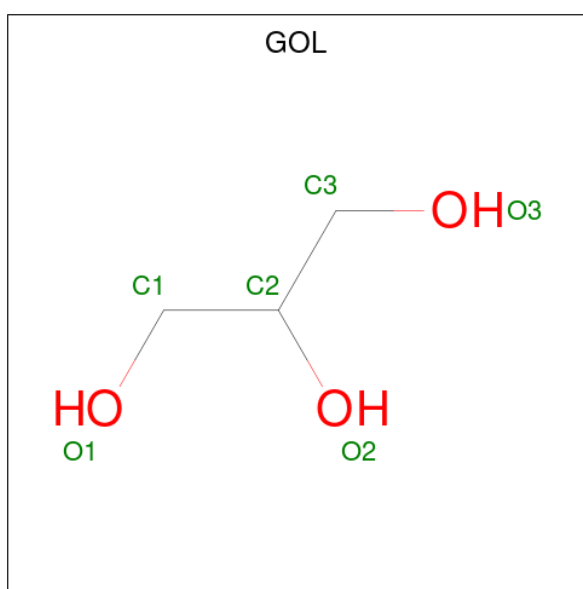
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

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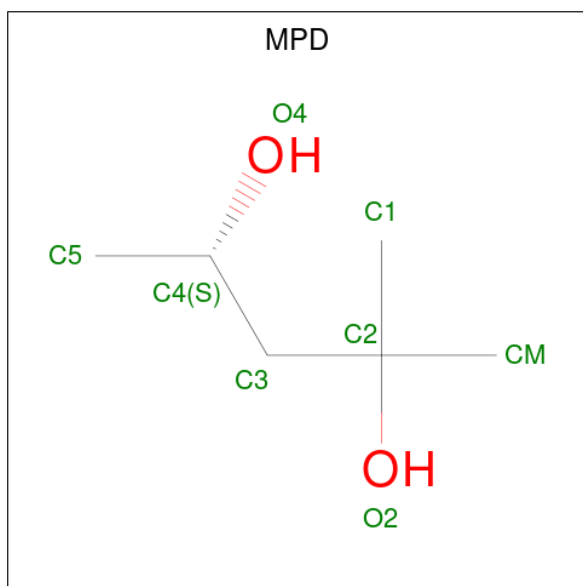
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total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

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

- Molecule 5 is (4S)-2-METHYL-2,4-PENTANEDIOL (CCD ID: MPD) (formula: C₆H₁₄O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			8	6	2		
5	A	1	Total	C	O	0	0
			8	6	2		
5	A	1	Total	C	O	0	0
			8	6	2		
5	A	1	Total	C	O	0	0
			8	6	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			8	6	2		
5	A	1	Total	C	O	0	0
			8	6	2		
5	B	1	Total	C	O	0	0
			8	6	2		
5	B	1	Total	C	O	0	0
			8	6	2		
5	B	1	Total	C	O	0	0
			8	6	2		
5	B	1	Total	C	O	0	0
			8	6	2		

- Molecule 6 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Cl	0	0
			1	1		

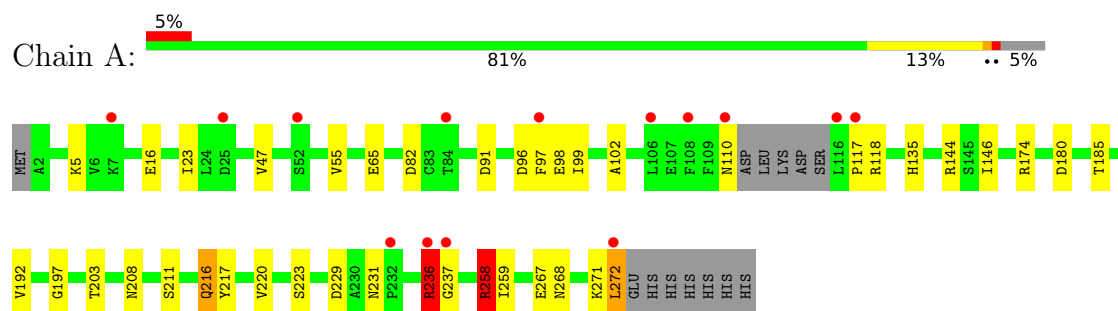
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	230	Total	O	0	0
			230	230		
7	B	208	Total	O	0	0
			208	208		

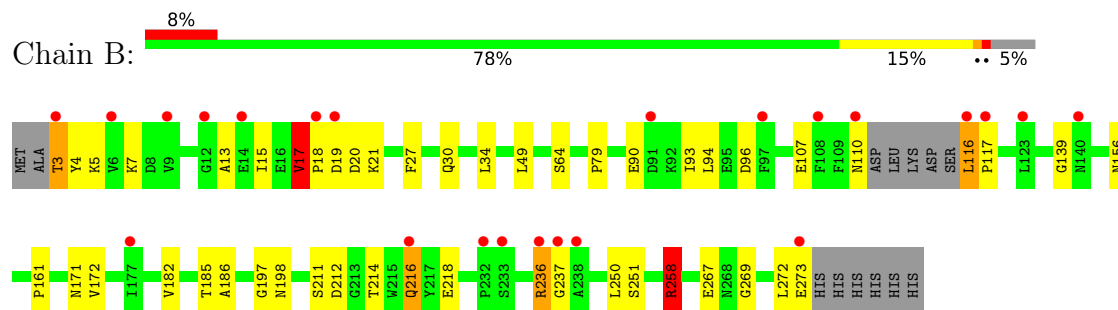
3 Residue-property plots [i](#)

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: Pectocin M2



• Molecule 1: Pectocin M2



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	117.45Å 117.45Å 128.45Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	58.79 – 1.86 58.79 – 1.86	Depositor EDS
% Data completeness (in resolution range)	100.0 (58.79-1.86) 100.0 (58.79-1.86)	Depositor EDS
R_{merge}	0.03	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.72 (at 1.86Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.168 , 0.197 0.176 , 0.203	Depositor DCC
R_{free} test set	4307 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	36.5	Xtriage
Anisotropy	0.458	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 56.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.023 for -h,-k,l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	5000	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 66.00 % 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 6.4851e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: FES, MPD, GOL, CL, SO4

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	1.47	10/2138 (0.5%)	1.28	9/2893 (0.3%)
1	B	1.50	14/2129 (0.7%)	1.27	12/2881 (0.4%)
All	All	1.49	24/4267 (0.6%)	1.28	21/5774 (0.4%)

The worst 5 of 24 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	267	GLU	CA-C	-8.89	1.41	1.52
1	B	34	LEU	CA-C	8.72	1.60	1.52
1	B	17	VAL	CA-CB	8.56	1.61	1.54
1	A	55	VAL	C-O	7.12	1.31	1.23
1	A	267	GLU	CA-C	-6.60	1.44	1.52

The worst 5 of 21 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	258	ARG	CG-CD-NE	-7.52	95.46	112.00
1	A	258	ARG	CG-CD-NE	-7.32	95.90	112.00
1	B	214	THR	CA-C-O	-6.99	113.51	121.19
1	A	216	GLN	CA-CB-CG	-6.87	100.36	114.10
1	B	237	GLY	N-CA-C	6.58	119.50	111.93

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	2088	0	2020	30	0
1	B	2085	0	2002	32	0
2	A	4	0	0	0	0
2	B	4	0	0	0	0
3	A	40	0	0	3	0
3	B	50	0	0	1	0
4	A	90	0	120	15	0
4	B	120	0	160	14	0
5	A	48	0	84	13	0
5	B	32	0	55	6	0
6	A	1	0	0	0	0
7	A	230	0	0	3	0
7	B	208	0	0	3	0
All	All	5000	0	4441	77	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 9.

The worst 5 of 77 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:236[A]:ARG:HH11	1:B:236[A]:ARG:HG3	1.15	1.10
1:B:171:ASN:HA	4:B:314:GOL:H12	1.44	0.98
5:A:328:MPD:H52	7:A:510:HOH:O	1.65	0.95
1:B:212:ASP:HB3	5:B:335:MPD:H4	1.54	0.90
1:B:236[A]:ARG:HH11	1:B:236[A]:ARG:CG	1.91	0.83

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	267/279 (96%)	263 (98%)	4 (2%)	0	100	100
1	B	266/279 (95%)	260 (98%)	6 (2%)	0	100	100
All	All	533/558 (96%)	523 (98%)	10 (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	231/239 (97%)	227 (98%)	4 (2%)	53	42
1	B	230/239 (96%)	217 (94%)	13 (6%)	18	6
All	All	461/478 (96%)	444 (96%)	17 (4%)	35	15

5 of 17 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	236[B]	ARG
1	B	273	GLU
1	B	30	GLN
1	B	107	GLU
1	B	110	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	216	GLN
1	A	225	ASN
1	B	268	ASN

5.3.3 RNA ⓘ

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 66 ligands modelled in this entry, 1 is monoatomic - leaving 65 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	SO4	A	303	-	4,4,4	0.28	0	6,6,6	0.43	0
4	GOL	A	316	-	5,5,5	0.43	0	5,5,5	0.60	0
4	GOL	A	321	-	5,5,5	0.46	0	5,5,5	0.17	0
5	MPD	A	328	-	7,7,7	1.32	1 (14%)	9,10,10	2.01	3 (33%)
3	SO4	B	306	-	4,4,4	0.18	0	6,6,6	0.59	0
3	SO4	B	302	-	4,4,4	0.26	0	6,6,6	1.44	1 (16%)
4	GOL	B	325	-	5,5,5	0.33	0	5,5,5	0.95	0
4	GOL	B	320	-	5,5,5	0.43	0	5,5,5	0.40	0
3	SO4	B	308	-	4,4,4	0.43	0	6,6,6	0.36	0
4	GOL	A	315	-	5,5,5	0.57	0	5,5,5	0.66	0
4	GOL	A	323	-	5,5,5	0.25	0	5,5,5	0.90	0
4	GOL	B	318	-	5,5,5	0.24	0	5,5,5	0.67	0
5	MPD	B	332	-	7,7,7	0.65	0	9,10,10	1.16	0
5	MPD	B	335	-	7,7,7	1.73	1 (14%)	9,10,10	1.18	1 (11%)
4	GOL	B	312	-	5,5,5	1.08	0	5,5,5	1.45	1 (20%)
2	FES	B	301	1	0,4,4	-	-	-	-	-
3	SO4	A	302	-	4,4,4	0.43	0	6,6,6	0.66	0
4	GOL	A	317	-	5,5,5	0.51	0	5,5,5	1.49	0
3	SO4	B	304	-	4,4,4	0.58	0	6,6,6	0.84	0
4	GOL	B	315	-	5,5,5	0.19	0	5,5,5	0.54	0
4	GOL	B	319	-	5,5,5	0.33	0	5,5,5	0.73	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	MPD	A	325	-	7,7,7	1.48	1 (14%)	9,10,10	3.20	5 (55%)
5	MPD	A	326	-	7,7,7	0.49	0	9,10,10	0.87	0
3	SO4	A	306	-	4,4,4	0.70	0	6,6,6	0.51	0
3	SO4	B	303	-	4,4,4	0.45	0	6,6,6	0.39	0
3	SO4	B	307	-	4,4,4	0.63	0	6,6,6	2.00	3 (50%)
4	GOL	A	311	-	5,5,5	0.47	0	5,5,5	0.57	0
4	GOL	A	313	-	5,5,5	1.11	0	5,5,5	2.61	1 (20%)
3	SO4	A	308	-	4,4,4	0.34	0	6,6,6	1.02	0
4	GOL	B	321	-	5,5,5	0.38	0	5,5,5	1.72	2 (40%)
4	GOL	B	324	-	5,5,5	0.29	0	5,5,5	0.48	0
4	GOL	B	327	-	5,5,5	0.96	0	5,5,5	1.14	1 (20%)
4	GOL	A	318	-	5,5,5	0.44	0	5,5,5	1.28	1 (20%)
3	SO4	A	309	-	4,4,4	0.51	0	6,6,6	0.52	0
4	GOL	A	312	-	5,5,5	1.09	1 (20%)	5,5,5	1.88	2 (40%)
3	SO4	B	311	-	4,4,4	0.47	0	6,6,6	0.36	0
4	GOL	B	316	-	5,5,5	0.54	0	5,5,5	0.61	0
4	GOL	B	331	-	5,5,5	0.74	0	5,5,5	1.29	1 (20%)
4	GOL	B	330	-	5,5,5	0.33	0	5,5,5	1.20	0
5	MPD	A	327	-	7,7,7	0.79	0	9,10,10	1.45	1 (11%)
4	GOL	B	329	-	5,5,5	0.29	0	5,5,5	1.00	0
4	GOL	B	314	-	5,5,5	0.74	0	5,5,5	1.46	1 (20%)
4	GOL	B	313	-	5,5,5	0.66	0	5,5,5	1.09	0
4	GOL	A	314	-	5,5,5	1.18	1 (20%)	5,5,5	3.80	3 (60%)
4	GOL	A	319	-	5,5,5	0.17	0	5,5,5	0.43	0
3	SO4	A	307	-	4,4,4	0.61	0	6,6,6	0.91	0
5	MPD	B	333	-	7,7,7	0.54	0	9,10,10	0.56	0
4	GOL	A	320	-	5,5,5	0.31	0	5,5,5	1.34	0
3	SO4	B	305	-	4,4,4	0.32	0	6,6,6	0.80	0
3	SO4	B	309	-	4,4,4	0.51	0	6,6,6	0.27	0
4	GOL	A	324	-	5,5,5	0.40	0	5,5,5	0.56	0
4	GOL	A	310	-	5,5,5	0.67	0	5,5,5	1.46	1 (20%)
4	GOL	B	323	-	5,5,5	0.80	0	5,5,5	1.57	2 (40%)
5	MPD	B	334	-	7,7,7	1.03	0	9,10,10	1.21	1 (11%)
3	SO4	A	304	-	4,4,4	0.42	0	6,6,6	0.49	0
3	SO4	B	310	-	4,4,4	0.59	0	6,6,6	0.58	0
4	GOL	B	317	-	5,5,5	0.38	0	5,5,5	1.23	1 (20%)
5	MPD	A	330	-	7,7,7	0.92	0	9,10,10	0.80	0
4	GOL	B	328	-	5,5,5	0.68	0	5,5,5	1.08	0
5	MPD	A	329	-	7,7,7	0.69	0	9,10,10	1.16	1 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	GOL	B	326	-	5,5,5	1.14	0	5,5,5	0.89	0
4	GOL	B	322	-	5,5,5	0.61	0	5,5,5	0.99	0
2	FES	A	301	1	0,4,4	-	-	-		
3	SO4	A	305	-	4,4,4	0.64	0	6,6,6	1.16	0
4	GOL	A	322	-	5,5,5	0.56	0	5,5,5	1.13	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	GOL	A	321	-	-	4/4/4/4	-
4	GOL	A	316	-	-	3/4/4/4	-
5	MPD	A	328	-	-	4/5/5/5	-
4	GOL	B	325	-	-	2/4/4/4	-
4	GOL	B	320	-	-	2/4/4/4	-
4	GOL	A	315	-	-	0/4/4/4	-
4	GOL	A	323	-	-	2/4/4/4	-
4	GOL	B	318	-	-	0/4/4/4	-
5	MPD	B	332	-	-	3/5/5/5	-
5	MPD	B	335	-	-	2/5/5/5	-
4	GOL	B	312	-	-	4/4/4/4	-
4	GOL	A	317	-	-	2/4/4/4	-
4	GOL	B	319	-	-	3/4/4/4	-
5	MPD	A	325	-	-	4/5/5/5	-
4	GOL	B	315	-	-	3/4/4/4	-
5	MPD	A	326	-	-	0/5/5/5	-
2	FES	B	301	1	-	-	0/1/1/1
4	GOL	B	321	-	-	2/4/4/4	-
4	GOL	A	311	-	-	2/4/4/4	-
4	GOL	B	324	-	-	3/4/4/4	-
4	GOL	A	313	-	-	2/4/4/4	-
4	GOL	B	327	-	-	4/4/4/4	-
4	GOL	A	318	-	-	4/4/4/4	-
4	GOL	A	312	-	-	4/4/4/4	-
4	GOL	B	316	-	-	2/4/4/4	-
4	GOL	B	331	-	-	3/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	B	330	-	-	4/4/4/4	-
5	MPD	A	327	-	-	4/5/5/5	-
4	GOL	B	329	-	-	2/4/4/4	-
4	GOL	B	314	-	-	2/4/4/4	-
4	GOL	B	313	-	-	2/4/4/4	-
4	GOL	A	314	-	-	0/4/4/4	-
4	GOL	A	319	-	-	3/4/4/4	-
5	MPD	B	333	-	-	2/5/5/5	-
4	GOL	A	320	-	-	2/4/4/4	-
4	GOL	A	324	-	-	2/4/4/4	-
4	GOL	A	310	-	-	3/4/4/4	-
4	GOL	B	323	-	-	2/4/4/4	-
5	MPD	B	334	-	-	1/5/5/5	-
4	GOL	B	317	-	-	0/4/4/4	-
5	MPD	A	330	-	-	4/5/5/5	-
4	GOL	B	328	-	-	2/4/4/4	-
5	MPD	A	329	-	-	5/5/5/5	-
4	GOL	B	326	-	-	4/4/4/4	-
4	GOL	B	322	-	-	4/4/4/4	-
2	FES	A	301	1	-	-	0/1/1/1
4	GOL	A	322	-	-	4/4/4/4	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	335	MPD	O2-C2	-3.64	1.35	1.44
5	A	325	MPD	C3-C2	2.70	1.62	1.54
5	A	328	MPD	O4-C4	2.04	1.51	1.43
4	A	314	GOL	O2-C2	2.01	1.49	1.43
4	A	312	GOL	O1-C1	2.00	1.50	1.42

The worst 5 of 33 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	325	MPD	CM-C2-C1	7.47	127.36	110.63
4	A	314	GOL	O2-C2-C3	5.76	133.04	109.18
4	A	313	GOL	C3-C2-C1	-5.50	91.64	111.80
4	A	314	GOL	O1-C1-C2	5.26	134.07	110.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	328	MPD	O4-C4-C3	4.47	129.13	111.35

There are no chirality outliers.

5 of 116 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	310	GOL	O1-C1-C2-C3
4	A	311	GOL	C1-C2-C3-O3
4	A	311	GOL	O2-C2-C3-O3
4	A	312	GOL	O1-C1-C2-C3
4	A	312	GOL	O2-C2-C3-O3

There are no ring outliers.

25 monomers are involved in 48 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	303	SO4	1	0
5	A	328	MPD	1	0
4	A	315	GOL	3	0
5	B	335	MPD	4	0
4	B	312	GOL	1	0
3	A	302	SO4	1	0
5	A	325	MPD	4	0
5	A	326	MPD	2	0
3	B	303	SO4	1	0
4	A	311	GOL	1	0
4	A	313	GOL	1	0
4	A	318	GOL	1	0
4	A	312	GOL	2	0
4	B	330	GOL	1	0
5	A	327	MPD	4	0
4	B	329	GOL	3	0
4	B	314	GOL	6	0
4	A	319	GOL	1	0
5	B	333	MPD	2	0
4	B	323	GOL	1	0
3	A	304	SO4	1	0
4	B	317	GOL	2	0
4	B	328	GOL	1	0
5	A	329	MPD	2	0
4	A	322	GOL	6	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	266/279 (95%)	0.47	14 (5%)	32 35	23, 40, 61, 105	5 (1%)
1	B	266/279 (95%)	0.54	23 (8%)	16 16	20, 43, 76, 85	4 (1%)
All	All	532/558 (95%)	0.50	37 (6%)	22 24	20, 41, 74, 105	9 (1%)

The worst 5 of 37 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	3	THR	5.8
1	A	236[A]	ARG	5.4
1	A	116	LEU	5.0
1	A	272	LEU	5.0
1	A	97	PHE	4.8

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
4	GOL	B	316	6/6	0.47	0.18	113,128,131,133	0
3	SO4	B	311	5/5	0.70	0.14	82,88,89,93	5
5	MPD	A	330	8/8	0.74	0.31	67,79,84,95	0
5	MPD	A	329	8/8	0.77	0.34	62,67,76,83	0
4	GOL	B	320	6/6	0.78	0.17	84,105,106,106	0
5	MPD	B	335	8/8	0.78	0.26	39,55,62,71	8
4	GOL	A	317	6/6	0.81	0.29	52,52,56,62	6
4	GOL	A	324	6/6	0.81	0.18	94,101,107,109	6
4	GOL	B	329	6/6	0.81	0.27	39,65,71,71	6
4	GOL	B	326	6/6	0.82	0.27	71,85,86,91	0
4	GOL	A	321	6/6	0.82	0.17	100,106,111,113	0
5	MPD	A	325	8/8	0.82	0.23	36,50,80,86	0
4	GOL	B	321	6/6	0.83	0.15	73,79,91,93	0
5	MPD	B	334	8/8	0.83	0.32	54,89,91,92	0
4	GOL	B	324	6/6	0.83	0.21	55,90,92,95	6
4	GOL	A	316	6/6	0.84	0.17	68,95,101,104	0
4	GOL	B	330	6/6	0.84	0.23	50,72,75,76	6
4	GOL	B	331	6/6	0.84	0.19	47,57,63,63	0
4	GOL	A	322	6/6	0.84	0.21	37,57,66,71	6
5	MPD	A	327	8/8	0.84	0.34	29,65,74,74	8
4	GOL	A	314	6/6	0.84	0.25	31,37,50,60	6
4	GOL	B	313	6/6	0.84	0.20	51,63,74,83	0
5	MPD	B	333	8/8	0.84	0.28	77,83,91,92	0
4	GOL	B	315	6/6	0.84	0.20	76,83,91,98	0
4	GOL	B	327	6/6	0.84	0.21	39,54,58,62	6
5	MPD	A	328	8/8	0.85	0.26	39,55,61,62	8
4	GOL	B	318	6/6	0.85	0.18	81,102,106,113	0
4	GOL	B	323	6/6	0.85	0.22	37,41,59,72	6
4	GOL	B	328	6/6	0.86	0.24	52,53,67,68	6
4	GOL	B	322	6/6	0.86	0.20	57,77,80,80	6
4	GOL	B	325	6/6	0.87	0.15	69,71,72,73	6
4	GOL	A	315	6/6	0.87	0.15	57,75,80,80	6
3	SO4	B	309	5/5	0.87	0.15	76,82,91,96	5
3	SO4	B	310	5/5	0.87	0.18	53,63,77,82	5
4	GOL	A	318	6/6	0.87	0.19	53,73,78,85	6
3	SO4	A	302	5/5	0.88	0.12	59,64,74,79	5
4	GOL	B	317	6/6	0.88	0.16	51,69,75,77	6
4	GOL	A	311	6/6	0.88	0.15	46,56,66,70	6
4	GOL	B	319	6/6	0.88	0.16	65,76,86,94	0
6	CL	A	331	1/1	0.88	0.14	84,84,84,84	0
4	GOL	B	314	6/6	0.89	0.15	33,45,49,54	6
4	GOL	A	320	6/6	0.89	0.22	47,76,77,79	6
4	GOL	A	323	6/6	0.89	0.17	90,91,97,98	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	MPD	B	332	8/8	0.90	0.18	39,46,51,60	5
3	SO4	B	305	5/5	0.91	0.13	46,65,72,81	5
3	SO4	B	308	5/5	0.91	0.12	61,81,95,99	5
4	GOL	A	312	6/6	0.91	0.20	31,39,50,53	6
4	GOL	A	313	6/6	0.91	0.19	33,37,45,59	6
5	MPD	A	326	8/8	0.91	0.22	51,68,74,87	3
3	SO4	A	305	5/5	0.91	0.12	36,56,58,71	5
3	SO4	A	307	5/5	0.91	0.11	36,51,65,66	5
4	GOL	A	310	6/6	0.92	0.17	45,52,58,64	6
3	SO4	B	304	5/5	0.92	0.11	39,57,67,71	5
3	SO4	A	306	5/5	0.92	0.10	50,55,63,76	5
3	SO4	A	308	5/5	0.92	0.13	42,66,71,77	5
3	SO4	A	309	5/5	0.93	0.13	74,87,94,103	0
3	SO4	B	303	5/5	0.93	0.07	54,62,67,81	5
4	GOL	B	312	6/6	0.93	0.16	49,64,71,76	0
3	SO4	B	307	5/5	0.94	0.09	50,54,59,66	0
4	GOL	A	319	6/6	0.94	0.12	50,62,70,74	0
3	SO4	A	304	5/5	0.95	0.06	58,70,74,85	5
3	SO4	A	303	5/5	0.95	0.10	44,55,61,72	5
3	SO4	B	306	5/5	0.96	0.13	49,61,66,77	5
3	SO4	B	302	5/5	0.96	0.08	40,53,58,70	5
2	FES	A	301	4/4	1.00	0.02	29,30,31,31	0
2	FES	B	301	4/4	1.00	0.02	34,34,34,35	0

6.5 Other polymers [i](#)

There are no such residues in this entry.