



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 03:12 AM UTC

PDB ID : 3RK8 / pdb_00003rk8
Title : Crystal structure of the chloride inhibited dihydrodipicolinate synthase from *Acinetobacter baumannii* complexed with pyruvate at 1.8 Å resolution
Authors : Kaushik, S.; Singh, A.; Sinha, M.; Kaur, P.; Sharma, S.; Singh, T.P.
Deposited on : 2011-04-17
Resolution : 1.80 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

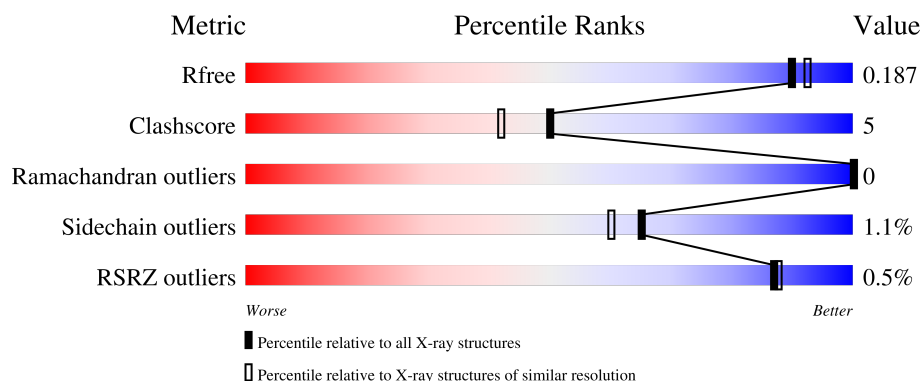
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	291	
1	B	291	

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
2	PYR	A	292	-	X	-	-
3	GOL	A	293	-	X	-	-
3	GOL	A	296	-	-	X	-
3	GOL	B	296	-	-	X	-

2 Entry composition [i](#)

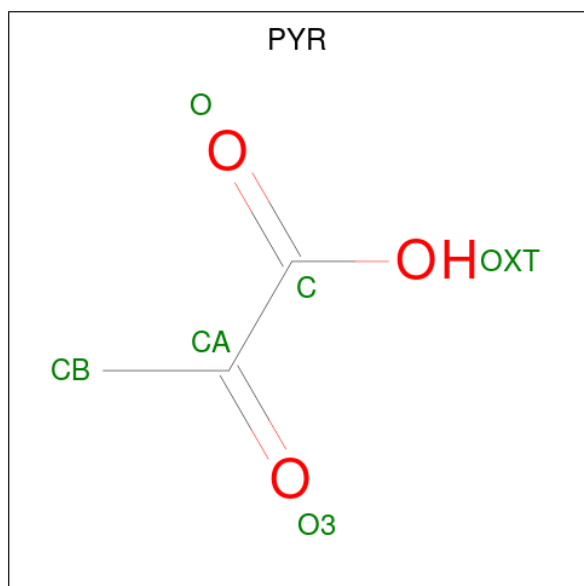
There are 6 unique types of molecules in this entry. The entry contains 5075 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Dihydrodipicolinate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	291	Total	C	N	O	S	0	0	0
			2195	1387	374	426	8			
1	B	291	Total	C	N	O	S	0	0	0
			2195	1387	374	426	8			

- Molecule 2 is PYRUVIC ACID (CCD ID: PYR) (formula: $C_3H_4O_3$).



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

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 4 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			7	4	3		
4	B	1	Total	C	O	0	0
			7	4	3		

- Molecule 5 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Cl	0	0
			1	1		
5	B	1	Total	Cl	0	0
			1	1		

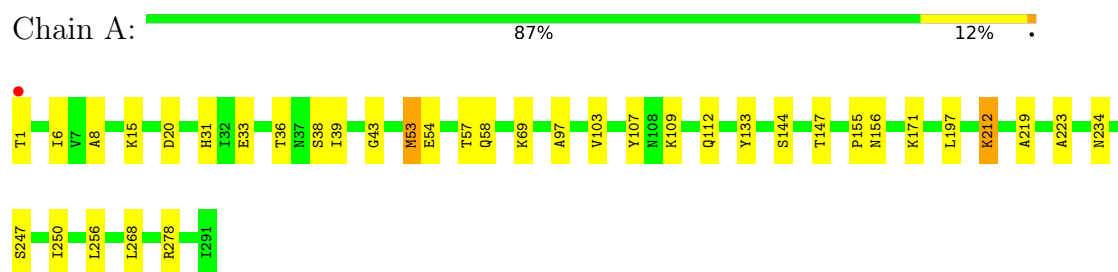
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	300	Total	O	0	0
			300	300		
6	B	303	Total	O	0	0
			303	303		

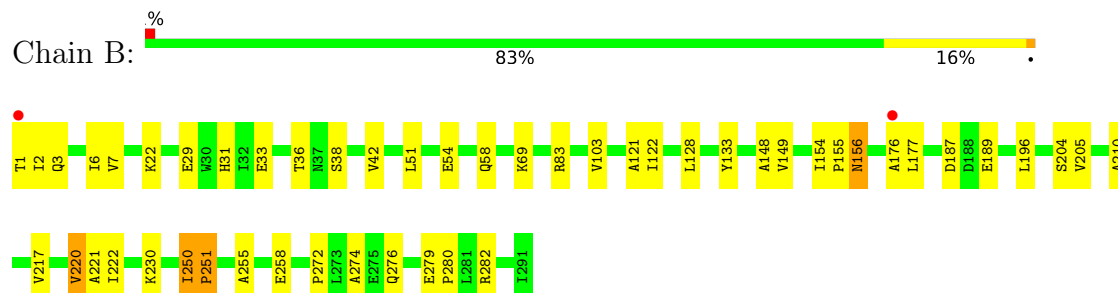
3 Residue-property plots [i](#)

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: Dihydrodipicolinate synthase



• Molecule 1: Dihydrodipicolinate synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	51.25Å 122.27Å 52.56Å 90.00° 116.31° 90.00°	Depositor
Resolution (Å)	47.12 – 1.80 47.12 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.7 (47.12-1.80) 99.7 (47.12-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.32 (at 1.79Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.148 , 0.188 0.149 , 0.187	Depositor DCC
R_{free} test set	2722 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	14.7	Xtriage
Anisotropy	0.757	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 44.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.018 for l,-k,h	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	5075	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.02% 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: PYR, CL, PEG, 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	1.53	14/2229 (0.6%)	1.24	2/3033 (0.1%)
1	B	1.56	20/2229 (0.9%)	1.28	8/3033 (0.3%)
All	All	1.55	34/4458 (0.8%)	1.26	10/6066 (0.2%)

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	220	VAL	C-N	9.79	1.46	1.33
1	A	268	LEU	C-O	-8.45	1.13	1.24
1	B	205	VAL	C-O	-7.58	1.15	1.24
1	B	250	ILE	CA-C	7.56	1.59	1.52
1	B	204	SER	C-O	-7.38	1.15	1.24
1	A	33	GLU	C-O	-7.03	1.16	1.24
1	A	39	ILE	C-O	-6.99	1.16	1.24
1	A	250	ILE	CA-C	6.36	1.59	1.52
1	B	121	ALA	CA-CB	6.32	1.63	1.53
1	B	210	ALA	CA-CB	6.25	1.60	1.53
1	B	220	VAL	C-O	-6.21	1.16	1.24
1	B	156	ASN	C-O	-6.13	1.16	1.24
1	B	217	VAL	CA-CB	6.09	1.62	1.54
1	B	230	LYS	CA-C	-5.95	1.45	1.52
1	A	43	GLY	C-O	-5.82	1.18	1.24
1	A	107	TYR	N-CA	5.79	1.53	1.46
1	A	223	ALA	N-CA	5.71	1.54	1.46
1	B	122	ILE	CA-CB	5.68	1.61	1.54
1	A	219	ALA	CA-CB	5.57	1.62	1.53
1	A	147	THR	CA-CB	5.46	1.61	1.53
1	B	149	VAL	N-CA	5.38	1.52	1.46
1	B	196	LEU	N-CA	5.26	1.52	1.46
1	B	272	PRO	CA-C	-5.24	1.46	1.52
1	A	8	ALA	CA-CB	5.23	1.59	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	251	PRO	C-O	5.20	1.30	1.24
1	A	155	PRO	CA-C	5.19	1.59	1.52
1	B	220	VAL	CA-C	5.18	1.59	1.52
1	A	247	SER	N-CA	5.14	1.52	1.46
1	A	97	ALA	CA-CB	5.12	1.62	1.53
1	B	148	ALA	CA-CB	5.11	1.61	1.53
1	B	177	LEU	N-CA	5.10	1.52	1.46
1	B	2	ILE	C-O	-5.09	1.18	1.24
1	A	256	LEU	C-O	-5.05	1.18	1.24
1	B	274	ALA	CA-CB	-5.04	1.44	1.53

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	128	LEU	CA-C-N	-6.64	113.12	119.76
1	B	128	LEU	C-N-CA	-6.64	113.12	119.76
1	B	255	ALA	N-CA-C	-6.61	104.00	111.07
1	B	176	ALA	CA-C-N	6.28	128.69	120.28
1	B	176	ALA	C-N-CA	6.28	128.69	120.28
1	B	276	GLN	N-CA-C	6.02	119.73	112.38
1	B	250	ILE	O-C-N	5.72	124.08	120.42
1	A	278	ARG	N-CA-C	5.55	117.41	111.36
1	B	177	LEU	N-CA-C	5.55	117.33	111.28
1	A	155	PRO	N-CA-C	5.44	121.09	113.53

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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	2195	0	2251	22	0
1	B	2195	0	2251	25	0
2	A	6	0	0	0	0
2	B	6	0	0	0	0
3	A	30	0	40	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	24	0	32	6	0
4	A	7	0	10	0	0
4	B	7	0	10	2	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	A	300	0	0	9	0
6	B	303	0	0	4	0
All	All	5075	0	4594	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 5.

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:B:189:GLU:OE2	3:B:296:GOL:H2	1.31	1.28
1:B:58:GLN:HG2	4:B:297:PEG:O4	1.69	0.91
1:B:187:ASP:OD2	3:B:296:GOL:H32	1.72	0.89
1:A:58:GLN:OE1	3:A:296:GOL:H2	1.80	0.80
1:B:189:GLU:OE2	3:B:296:GOL:C2	2.23	0.79
1:A:54:GLU:OE1	3:A:296:GOL:H31	1.83	0.79
1:A:20:ASP:OD1	3:A:295:GOL:H12	1.87	0.75
1:A:212:LYS:HD2	6:A:559:HOH:O	1.89	0.70
1:B:58:GLN:HB3	4:B:297:PEG:H31	1.78	0.65
1:B:29:GLU:O	1:B:33:GLU:HG3	1.96	0.65
1:A:20:ASP:OD1	3:A:295:GOL:C1	2.49	0.59
1:B:31:HIS:HD2	1:B:36:THR:OG1	1.87	0.57
1:B:187:ASP:CG	3:B:296:GOL:H32	2.28	0.57
1:A:31:HIS:HE1	6:A:398:HOH:O	1.89	0.55
1:A:15:LYS:O	3:A:293:GOL:H11	2.11	0.51
1:A:234:ASN:CG	6:A:492:HOH:O	2.54	0.50
1:A:6:ILE:HG12	1:A:38:SER:HB3	1.94	0.49
1:B:42:VAL:HG13	1:B:51:LEU:HD12	1.94	0.49
1:A:234:ASN:HB3	6:A:492:HOH:O	2.13	0.49
1:A:109:LYS:HE3	6:A:544:HOH:O	2.13	0.48
1:A:171:LYS:HE2	1:A:197:LEU:HD22	1.95	0.48
1:B:54:GLU:O	1:B:58:GLN:HG3	2.14	0.48
1:B:103:VAL:HA	1:B:133:TYR:HB3	1.96	0.47
1:A:31:HIS:HD2	1:A:36:THR:OG1	1.98	0.47
1:B:7:VAL:HG12	1:B:36:THR:HG21	1.96	0.47
1:B:279:GLU:N	1:B:280:PRO:HD2	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:112:GLN:HE22	1:A:144:SER:H	1.62	0.46
1:A:53:MET:HE1	6:A:324:HOH:O	2.16	0.46
1:B:1:THR:O	1:B:3:GLN:HG3	2.16	0.45
1:A:234:ASN:CB	6:A:492:HOH:O	2.64	0.45
1:B:222:ILE:HG23	3:B:295:GOL:H32	1.99	0.45
1:B:31:HIS:HE1	6:B:324:HOH:O	1.99	0.45
3:A:296:GOL:H12	6:A:436:HOH:O	2.17	0.45
1:B:83:ARG:NH1	6:B:307:HOH:O	2.50	0.45
1:B:220:VAL:HG23	1:B:221:ALA:N	2.32	0.44
1:B:250:ILE:HB	1:B:251:PRO:HD3	1.99	0.44
1:B:187:ASP:OD2	3:B:296:GOL:C3	2.56	0.43
1:A:103:VAL:HA	1:A:133:TYR:HB3	2.00	0.43
1:B:69:LYS:NZ	6:B:380:HOH:O	2.52	0.42
1:B:155:PRO:O	1:B:156:ASN:HB2	2.20	0.42
1:A:58:GLN:OE1	3:A:296:GOL:C2	2.60	0.42
1:B:154:ILE:HA	1:B:155:PRO:HD3	1.96	0.42
1:A:109:LYS:HD3	6:B:555:HOH:O	2.19	0.41
1:A:156:ASN:OD1	1:A:156:ASN:N	2.44	0.41
1:A:112:GLN:NE2	1:A:144:SER:H	2.17	0.41
1:A:1:THR:N	6:A:578:HOH:O	2.54	0.41
1:B:6:ILE:HG12	1:B:38:SER:HB3	2.02	0.41
1:B:258:GLU:CD	1:B:282:ARG:HD3	2.46	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	289/291 (99%)	286 (99%)	3 (1%)	0	100	100
1	B	289/291 (99%)	285 (99%)	4 (1%)	0	100	100
All	All	578/582 (99%)	571 (99%)	7 (1%)	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	234/234 (100%)	230 (98%)	4 (2%)	53	45
1	B	234/234 (100%)	233 (100%)	1 (0%)	84	83
All	All	468/468 (100%)	463 (99%)	5 (1%)	65	60

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	53	MET
1	A	57	THR
1	A	69	LYS
1	A	212	LYS
1	B	22	LYS

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	31	HIS
1	A	108	ASN
1	A	112	GLN
1	A	241	ASN
1	A	283	ASN
1	B	31	HIS
1	B	68	ASN
1	B	238	ASN
1	B	283	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 15 ligands modelled in this entry, 2 are monoatomic - leaving 13 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	A	294	-	5,5,5	0.49	0	5,5,5	1.37	1 (20%)
3	GOL	A	297	-	5,5,5	0.51	0	5,5,5	1.99	1 (20%)
2	PYR	B	292	-	5,5,5	3.38	3 (60%)	3,6,6	1.39	0
4	PEG	A	298	-	6,6,6	0.52	0	5,5,5	0.36	0
3	GOL	B	295	-	5,5,5	0.40	0	5,5,5	1.29	1 (20%)
3	GOL	B	294	-	5,5,5	0.54	0	5,5,5	0.43	0
4	PEG	B	297	-	6,6,6	0.64	0	5,5,5	0.54	0
3	GOL	A	295	-	5,5,5	0.86	0	5,5,5	1.57	1 (20%)
3	GOL	B	296	-	5,5,5	0.85	0	5,5,5	1.22	0
3	GOL	A	296	-	5,5,5	0.35	0	5,5,5	0.75	0
2	PYR	A	292	-	5,5,5	3.83	3 (60%)	3,6,6	1.70	1 (33%)
3	GOL	B	293	-	5,5,5	0.45	0	5,5,5	1.00	0
3	GOL	A	293	-	5,5,5	0.55	0	5,5,5	1.89	3 (60%)

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	A	294	-	-	2/4/4/4	-
3	GOL	A	297	-	-	4/4/4/4	-
2	PYR	B	292	-	-	1/4/4/4	-
4	PEG	A	298	-	-	3/4/4/4	-
3	GOL	B	295	-	-	2/4/4/4	-
3	GOL	B	294	-	-	2/4/4/4	-
4	PEG	B	297	-	-	2/4/4/4	-
3	GOL	A	295	-	-	4/4/4/4	-
3	GOL	B	296	-	-	3/4/4/4	-
3	GOL	A	296	-	-	3/4/4/4	-
2	PYR	A	292	-	-	2/4/4/4	-
3	GOL	B	293	-	-	4/4/4/4	-
3	GOL	A	293	-	-	3/4/4/4	-

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	292	PYR	O-C	5.86	1.37	1.22
2	B	292	PYR	O-C	4.74	1.34	1.22
2	A	292	PYR	O3-CA	4.59	1.33	1.23
2	B	292	PYR	O3-CA	4.17	1.32	1.23
2	A	292	PYR	CA-C	-3.73	1.42	1.54
2	B	292	PYR	CA-C	-3.73	1.42	1.54

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	297	GOL	C3-C2-C1	-4.31	95.99	111.80
3	A	295	GOL	O2-C2-C3	2.98	121.52	109.18
3	B	295	GOL	O1-C1-C2	-2.58	98.76	110.38
3	A	293	GOL	O3-C3-C2	2.49	121.57	110.38
3	A	293	GOL	O2-C2-C1	-2.33	99.55	109.18
3	A	293	GOL	O1-C1-C2	-2.20	100.45	110.38
2	A	292	PYR	O3-CA-CB	-2.18	114.88	119.77
3	A	294	GOL	O2-C2-C1	2.04	117.62	109.18

There are no chirality outliers.

All (35) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	292	PYR	OXT-C-CA-CB
2	B	292	PYR	OXT-C-CA-CB
3	A	295	GOL	O1-C1-C2-C3
3	A	295	GOL	C1-C2-C3-O3
3	A	296	GOL	O1-C1-C2-C3
3	A	297	GOL	O1-C1-C2-C3
3	A	297	GOL	C1-C2-C3-O3
3	A	297	GOL	O2-C2-C3-O3
3	B	293	GOL	O1-C1-C2-C3
3	B	294	GOL	O1-C1-C2-C3
3	B	296	GOL	C1-C2-C3-O3
3	A	295	GOL	O1-C1-C2-O2
3	B	293	GOL	O2-C2-C3-O3
3	B	296	GOL	O2-C2-C3-O3
4	A	298	PEG	O2-C3-C4-O4
4	B	297	PEG	O2-C3-C4-O4
3	A	293	GOL	C1-C2-C3-O3
3	A	294	GOL	C1-C2-C3-O3
3	A	296	GOL	C1-C2-C3-O3
3	B	293	GOL	C1-C2-C3-O3
3	B	295	GOL	C1-C2-C3-O3
3	B	296	GOL	O1-C1-C2-C3
3	A	293	GOL	O2-C2-C3-O3
3	A	295	GOL	O2-C2-C3-O3
3	A	297	GOL	O1-C1-C2-O2
3	B	294	GOL	O1-C1-C2-O2
4	A	298	PEG	O1-C1-C2-O2
3	A	296	GOL	O1-C1-C2-O2
3	B	293	GOL	O1-C1-C2-O2
4	B	297	PEG	O1-C1-C2-O2
3	A	294	GOL	O2-C2-C3-O3
3	B	295	GOL	O2-C2-C3-O3
4	A	298	PEG	C1-C2-O2-C3
2	A	292	PYR	O-C-CA-CB
3	A	293	GOL	O1-C1-C2-O2

There are no ring outliers.

6 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	295	GOL	1	0
4	B	297	PEG	2	0
3	A	295	GOL	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	296	GOL	5	0
3	A	296	GOL	4	0
3	A	293	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 [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	291/291 (100%)	-0.66	1 (0%) 90 90	7, 13, 26, 31	0
1	B	291/291 (100%)	-0.58	2 (0%) 84 85	8, 13, 26, 33	0
All	All	582/582 (100%)	-0.62	3 (0%) 87 88	7, 13, 26, 33	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1	THR	5.9
1	A	1	THR	4.3
1	B	176	ALA	3.0

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(Å ²)	Q<0.9
3	GOL	A	296	6/6	0.72	0.19	54,56,57,59	0
4	PEG	A	298	7/7	0.75	0.17	43,49,52,52	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	GOL	A	297	6/6	0.78	0.18	27,35,40,41	0
4	PEG	B	297	7/7	0.78	0.20	52,54,56,57	0
3	GOL	B	293	6/6	0.83	0.16	29,45,50,51	0
2	PYR	B	292	6/6	0.83	0.19	12,28,31,34	6
3	GOL	A	294	6/6	0.83	0.15	35,41,43,44	0
2	PYR	A	292	6/6	0.84	0.19	16,33,36,37	0
3	GOL	B	294	6/6	0.85	0.14	40,48,50,52	0
3	GOL	B	296	6/6	0.85	0.12	20,31,37,40	0
3	GOL	A	293	6/6	0.86	0.13	25,38,41,42	0
3	GOL	B	295	6/6	0.92	0.10	18,31,39,41	0
3	GOL	A	295	6/6	0.94	0.09	14,22,28,31	0
5	CL	B	298	1/1	0.98	0.05	24,24,24,24	0
5	CL	A	299	1/1	0.99	0.04	18,18,18,18	0

6.5 Other polymers [i](#)

There are no such residues in this entry.