



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 04:29 AM UTC

PDB ID : 3L57 / pdb_00003l57
Title : Crystal Structure of the Plasmid pCU1 TraI Relaxase Domain
Authors : Redinbo, M.R.; Nash, R.P.
Deposited on : 2009-12-21
Resolution : 2.29 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

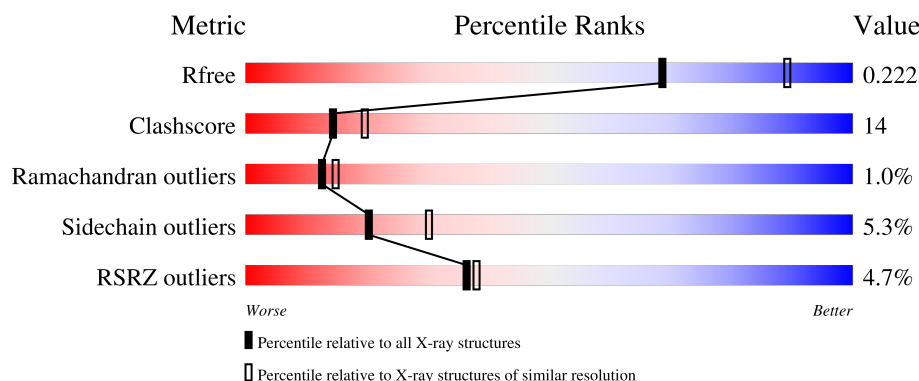
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.29 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	299	3% 60% 12% • 25%
1	B	299	4% 52% 13% • 32%

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	EDO	A	316	-	-	X	-
3	EDO	B	304	-	-	X	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 3684 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.

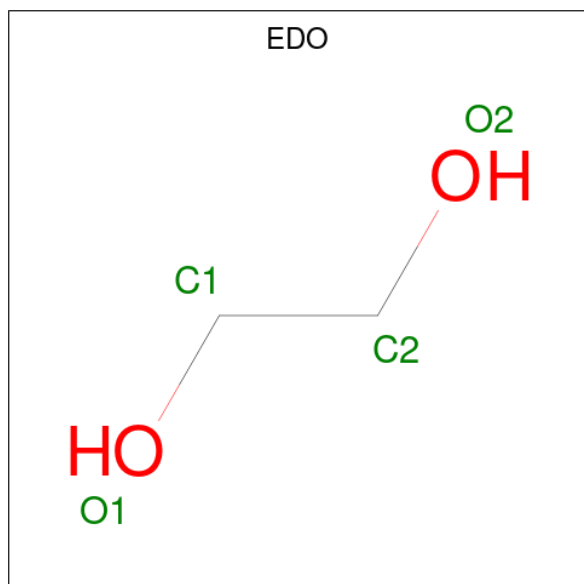
- Molecule 1 is a protein called Mobilization protein TraI.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	223	Total	C	N	O	S	0	8	0
			1762	1095	315	343	9			
1	B	203	Total	C	N	O	S	0	7	0
			1574	983	284	299	8			

- Molecule 2 is MANGANESE (III) ION (CCD ID: MN3) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mn	0	0
			1	1		
2	B	1	Total	Mn	0	0
			1	1		

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



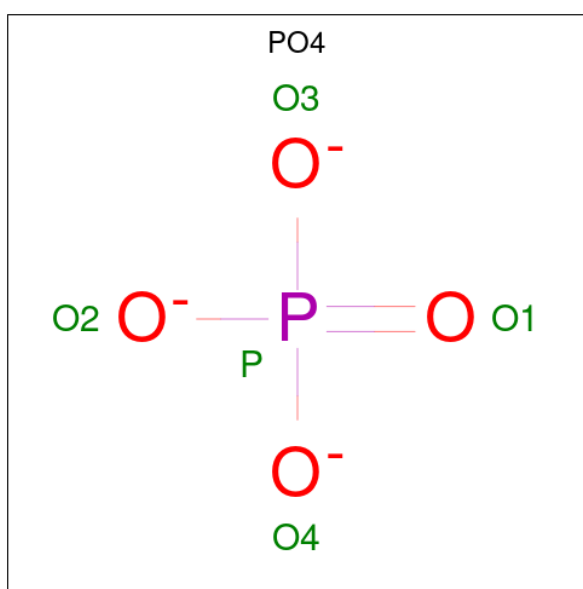
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0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

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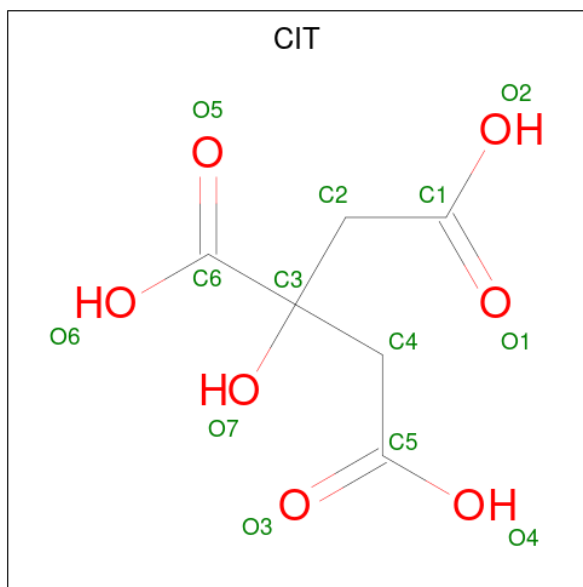
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	P	0	0
			5	4	1		

- Molecule 5 is CITRIC ACID (CCD ID: CIT) (formula: $\text{C}_6\text{H}_8\text{O}_7$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			13	6	7		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	156	Total	O	0	2
			158	158		
6	B	63	Total	O	0	3
			66	66		

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	50.28Å 58.52Å 87.12Å 90.00° 83.88° 90.00°	Depositor
Resolution (Å)	48.49 – 2.29 48.49 – 2.29	Depositor EDS
% Data completeness (in resolution range)	97.0 (48.49-2.29) 96.9 (48.49-2.29)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.91 (at 2.29Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.173 , 0.226 0.173 , 0.222	Depositor DCC
R_{free} test set	1105 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	36.2	Xtriage
Anisotropy	0.327	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 72.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3684	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.94% 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: PO4, EDO, CIT, MN3

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	0.33	0/1810	0.69	2/2444 (0.1%)
1	B	0.30	0/1618	0.65	1/2181 (0.0%)
All	All	0.32	0/3428	0.67	3/4625 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	156	ASP	CA-C-N	5.41	125.72	119.93
1	A	156	ASP	C-N-CA	5.41	125.72	119.93
1	B	150	GLU	N-CA-C	-5.12	106.16	112.72

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	1762	0	1696	45	0
1	B	1574	0	1511	49	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	76	0	114	13	0
3	B	28	0	42	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	5	0	0	0	0
5	A	13	0	5	0	0
6	A	158	0	0	6	0
6	B	66	0	0	2	0
All	All	3684	0	3368	96	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 14.

All (96) 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:A:9[A]:ARG:HH11	1:A:9[A]:ARG:HG3	1.05	1.18
1:A:153:ARG:HH11	1:A:153:ARG:HG3	1.18	1.07
1:A:9[A]:ARG:HG3	1:A:9[A]:ARG:NH1	1.83	0.90
1:B:137:GLN:HE22	1:B:172:GLU:H	1.17	0.87
1:A:153:ARG:HH11	1:A:153:ARG:CG	1.88	0.85
1:B:8[B]:THR:H	1:B:11:ASN:HD21	1.23	0.82
1:B:8[A]:THR:H	1:B:11:ASN:HD21	1.24	0.81
1:B:188:LYS:HB3	3:B:302:EDO:H11	1.66	0.78
1:A:153:ARG:HG3	1:A:153:ARG:NH1	1.89	0.77
1:B:8[B]:THR:H	1:B:11:ASN:ND2	1.82	0.77
1:B:8[A]:THR:H	1:B:11:ASN:ND2	1.83	0.76
1:B:59:LEU:HA	1:B:166[B]:MET:HE1	1.67	0.76
1:A:95[A]:MET:HG2	1:A:157:PRO:HG2	1.68	0.76
1:A:190:HIS:H	3:A:316:EDO:H21	1.50	0.74
1:B:213:ASN:C	1:B:213:ASN:HD22	1.95	0.74
1:A:8:THR:O	1:A:12[A]:VAL:HG22	1.90	0.72
1:A:190:HIS:N	3:A:316:EDO:H21	2.05	0.72
1:A:128:THR:HG23	1:A:135:VAL:HG13	1.72	0.70
1:A:48:GLY:HA2	3:A:302:EDO:H21	1.74	0.70
1:B:9:ARG:HG3	1:B:81:LEU:HD11	1.77	0.66
1:B:11:ASN:HD22	1:B:11:ASN:C	2.01	0.65
1:B:8[A]:THR:HG23	1:B:10:GLN:H	1.61	0.65
1:B:59:LEU:HA	1:B:166[B]:MET:CE	2.28	0.64
1:A:9[A]:ARG:HH11	1:A:9[A]:ARG:CG	1.95	0.64
1:B:110:GLU:OE2	1:B:148:ARG:HD3	1.98	0.64
1:B:59:LEU:HD22	1:B:81:LEU:HD23	1.80	0.63
1:A:115:ALA:HB1	3:A:309:EDO:H11	1.80	0.62
1:A:211:ARG:NH2	1:A:223:HIS:H	1.95	0.62
1:B:11:ASN:C	1:B:13:THR:H	2.07	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:4:ILE:HG12	1:A:85[A]:LEU:HD22	1.81	0.61
1:B:11:ASN:HD22	1:B:12:VAL:N	1.97	0.61
1:A:35:SER:O	1:A:146[B]:THR:HG22	2.01	0.61
3:B:304:EDO:H21	6:B:328:HOH:O	2.00	0.61
1:A:201:LEU:HD13	3:B:304:EDO:O1	2.02	0.59
1:B:4:ILE:HG12	1:B:85:LEU:HD22	1.84	0.59
1:A:190:HIS:HB2	3:A:316:EDO:H12	1.84	0.59
1:A:210:LEU:HB2	3:A:317:EDO:H11	1.86	0.57
1:B:59:LEU:HD23	1:B:81:LEU:HB3	1.86	0.56
1:B:1:MET:O	1:B:87:PHE:HA	2.05	0.56
1:B:59:LEU:CD2	1:B:81:LEU:HB3	2.36	0.55
1:A:10:GLN:O	3:A:301:EDO:H12	2.07	0.55
1:B:137:GLN:NE2	1:B:172:GLU:H	1.96	0.55
1:A:33:PHE:N	6:A:471:HOH:O	2.40	0.55
1:B:213:ASN:C	1:B:213:ASN:ND2	2.65	0.55
1:B:54:ARG:HA	1:B:54:ARG:HE	1.72	0.54
1:B:7:ILE:HG13	1:B:164[A]:PHE:CZ	2.43	0.54
1:B:98:LEU:HD22	1:B:150:GLU:HB3	1.90	0.54
1:A:187:ASN:HA	3:A:316:EDO:O2	2.08	0.54
1:B:7:ILE:HG21	1:B:164[A]:PHE:HZ	1.73	0.53
1:A:9[A]:ARG:NH1	1:A:9[A]:ARG:CG	2.62	0.53
1:A:16:VAL:HG12	6:A:410:HOH:O	2.09	0.52
1:A:16:VAL:HG11	1:A:162:HIS:CG	2.45	0.52
1:B:8[A]:THR:HG23	1:B:10:GLN:N	2.23	0.52
1:B:189:MET:HE2	3:B:304:EDO:H12	1.91	0.52
1:B:11:ASN:ND2	1:B:11:ASN:C	2.68	0.52
1:B:187:ASN:HA	3:B:304:EDO:O2	2.09	0.52
1:B:162:HIS:HB3	1:B:164[B]:PHE:CE2	2.45	0.51
1:B:2:LEU:HD13	1:B:188:LYS:HG3	1.93	0.50
1:A:13:THR:HG23	3:A:301:EDO:H21	1.94	0.50
1:A:217:ASN:OD1	3:A:318:EDO:H12	2.13	0.49
1:B:162:HIS:HB3	1:B:164[B]:PHE:HE2	1.79	0.48
1:B:64:ASP:OD1	1:B:64:ASP:C	2.56	0.48
1:A:1:MET:HE2	6:A:354:HOH:O	2.13	0.47
1:B:46:LEU:HD11	1:B:58:LEU:HD11	1.95	0.47
1:A:197:GLN:HA	3:B:304:EDO:H12	1.96	0.47
1:B:131:GLN:O	1:B:131:GLN:HG2	2.14	0.47
1:A:9[A]:ARG:HA	1:A:12[A]:VAL:CG2	2.45	0.46
1:B:36:TRP:CZ2	1:B:50:VAL:HG13	2.50	0.46
1:B:54:ARG:HH21	1:B:57:GLU:HB3	1.80	0.46
1:A:33:PHE:HB2	6:A:471:HOH:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:166[B]:MET:HE3	1:B:166[B]:MET:HA	1.99	0.45
1:B:11:ASN:C	1:B:13:THR:N	2.74	0.45
1:A:99:ILE:CD1	1:A:157:PRO:HG3	2.47	0.45
1:B:8[B]:THR:HG22	1:B:9:ARG:N	2.31	0.45
1:A:13:THR:HG23	3:A:301:EDO:C2	2.46	0.45
1:A:152:SER:C	1:A:153:ARG:HD2	2.42	0.45
1:B:173:ASP:OD1	3:B:306:EDO:H11	2.17	0.44
1:A:79:GLU:O	1:A:79:GLU:HG3	2.17	0.44
1:A:1:MET:HE3	1:A:1:MET:HB2	1.91	0.43
1:B:91:LYS:HG3	6:B:323:HOH:O	2.18	0.43
1:B:166[B]:MET:HB2	1:B:168:MET:HG2	2.01	0.42
1:A:37:GLN:HE22	1:A:114:ALA:HA	1.85	0.42
1:A:1:MET:O	1:A:87:PHE:HA	2.19	0.41
1:A:190:HIS:H	3:A:316:EDO:C2	2.26	0.41
1:A:198:GLU:HG2	6:A:447:HOH:O	2.19	0.41
1:B:84:ASP:OD1	1:B:84:ASP:C	2.63	0.41
1:A:99:ILE:HD11	1:A:157:PRO:HG3	2.03	0.41
3:A:313:EDO:H11	6:A:326:HOH:O	2.21	0.41
1:B:208:TYR:CD2	1:B:221:MET:HE1	2.56	0.40
1:A:146[A]:THR:HA	1:A:162:HIS:O	2.22	0.40
1:B:54:ARG:HH21	1:B:57:GLU:CB	2.34	0.40
1:B:71:ARG:O	1:B:72:HIS:C	2.63	0.40
1:A:199:LEU:HG	1:A:219:PHE:HZ	1.87	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	227/299 (76%)	219 (96%)	8 (4%)	0	100	100
1	B	203/299 (68%)	194 (96%)	5 (2%)	4 (2%)	6	5

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	430/598 (72%)	413 (96%)	13 (3%)	4 (1%)	12	17

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	14	SER
1	B	12	VAL
1	B	151	THR
1	B	78	LYS

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	182/251 (72%)	170 (93%)	12 (7%)	15	21
1	B	156/251 (62%)	149 (96%)	7 (4%)	24	37
All	All	338/502 (67%)	319 (94%)	19 (6%)	20	28

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9[A]	ARG
1	A	9[C]	ARG
1	A	12[A]	VAL
1	A	12[B]	VAL
1	A	59	LEU
1	A	65	THR
1	A	99	ILE
1	A	128	THR
1	A	135	VAL
1	A	153	ARG
1	A	201	LEU
1	A	216	ASN
1	B	2	LEU
1	B	11	ASN

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Mol	Chain	Res	Type
1	B	54	ARG
1	B	59	LEU
1	B	106	ILE
1	B	188	LYS
1	B	213	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	GLN
1	A	37	GLN
1	A	70	GLN
1	A	96	GLN
1	A	170	GLN
1	A	216	ASN
1	B	11	ASN
1	B	37	GLN
1	B	137	GLN
1	B	170	GLN
1	B	213	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 30 ligands modelled in this entry, 2 are monoatomic - leaving 28 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	EDO	A	313	-	3,3,3	0.45	0	2,2,2	0.72	0
3	EDO	A	305	-	3,3,3	0.45	0	2,2,2	0.60	0
3	EDO	A	308	-	3,3,3	0.44	0	2,2,2	0.71	0
3	EDO	B	307	-	3,3,3	0.44	0	2,2,2	0.70	0
3	EDO	B	302	-	3,3,3	0.44	0	2,2,2	0.75	0
5	CIT	A	321	2	12,12,12	2.94	5 (41%)	17,17,17	1.45	1 (5%)
3	EDO	A	318	-	3,3,3	0.46	0	2,2,2	0.73	0
3	EDO	B	301	-	3,3,3	0.44	0	2,2,2	0.71	0
3	EDO	A	314	-	3,3,3	0.44	0	2,2,2	0.64	0
3	EDO	A	315	-	3,3,3	0.44	0	2,2,2	0.67	0
3	EDO	A	316	-	3,3,3	0.46	0	2,2,2	0.71	0
3	EDO	B	304	-	3,3,3	0.46	0	2,2,2	0.84	0
3	EDO	A	304	-	3,3,3	0.45	0	2,2,2	0.72	0
3	EDO	B	303	-	3,3,3	0.42	0	2,2,2	0.73	0
3	EDO	A	306	-	3,3,3	0.44	0	2,2,2	0.59	0
3	EDO	A	311	-	3,3,3	0.44	0	2,2,2	0.69	0
4	PO4	A	320	-	4,4,4	0.95	0	6,6,6	0.49	0
3	EDO	A	309	-	3,3,3	0.44	0	2,2,2	0.77	0
3	EDO	A	301	-	3,3,3	0.45	0	2,2,2	0.67	0
3	EDO	A	302	-	3,3,3	0.44	0	2,2,2	0.79	0
3	EDO	A	312	-	3,3,3	0.44	0	2,2,2	0.73	0
3	EDO	A	310	-	3,3,3	0.44	0	2,2,2	0.71	0
3	EDO	B	306	-	3,3,3	0.44	0	2,2,2	0.69	0
3	EDO	A	303	-	3,3,3	0.44	0	2,2,2	0.68	0
3	EDO	B	305	-	3,3,3	0.44	0	2,2,2	0.77	0
3	EDO	A	307	-	3,3,3	0.45	0	2,2,2	0.67	0
3	EDO	A	319	-	3,3,3	0.44	0	2,2,2	0.76	0
3	EDO	A	317	-	3,3,3	0.45	0	2,2,2	0.67	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	EDO	A	313	-	-	0/1/1/1	-
3	EDO	A	305	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EDO	A	308	-	-	1/1/1/1	-
3	EDO	B	307	-	-	1/1/1/1	-
3	EDO	B	302	-	-	1/1/1/1	-
5	CIT	A	321	2	-	3/16/16/16	-
3	EDO	A	318	-	-	1/1/1/1	-
3	EDO	B	301	-	-	1/1/1/1	-
3	EDO	A	314	-	-	1/1/1/1	-
3	EDO	A	315	-	-	1/1/1/1	-
3	EDO	A	316	-	-	1/1/1/1	-
3	EDO	B	304	-	-	1/1/1/1	-
3	EDO	A	304	-	-	1/1/1/1	-
3	EDO	B	303	-	-	1/1/1/1	-
3	EDO	A	306	-	-	1/1/1/1	-
3	EDO	A	311	-	-	1/1/1/1	-
3	EDO	A	309	-	-	1/1/1/1	-
3	EDO	A	301	-	-	1/1/1/1	-
3	EDO	A	302	-	-	1/1/1/1	-
3	EDO	A	312	-	-	1/1/1/1	-
3	EDO	A	310	-	-	1/1/1/1	-
3	EDO	B	306	-	-	1/1/1/1	-
3	EDO	A	303	-	-	1/1/1/1	-
3	EDO	B	305	-	-	0/1/1/1	-
3	EDO	A	307	-	-	1/1/1/1	-
3	EDO	A	319	-	-	1/1/1/1	-
3	EDO	A	317	-	-	1/1/1/1	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	321	CIT	O5-C6	5.81	1.40	1.22
5	A	321	CIT	O3-C5	5.66	1.40	1.22
5	A	321	CIT	O4-C5	-3.60	1.18	1.30
5	A	321	CIT	O6-C6	-3.25	1.18	1.30
5	A	321	CIT	O2-C1	2.96	1.40	1.30

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	321	CIT	O6-C6-C3	3.57	119.99	113.14

There are no chirality outliers.

All (27) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	321	CIT	C1-C2-C3-C4
5	A	321	CIT	C1-C2-C3-C6
3	A	302	EDO	O1-C1-C2-O2
3	A	307	EDO	O1-C1-C2-O2
3	A	316	EDO	O1-C1-C2-O2
3	A	317	EDO	O1-C1-C2-O2
5	A	321	CIT	C1-C2-C3-O7
3	A	308	EDO	O1-C1-C2-O2
3	A	310	EDO	O1-C1-C2-O2
3	B	304	EDO	O1-C1-C2-O2
3	B	306	EDO	O1-C1-C2-O2
3	B	303	EDO	O1-C1-C2-O2
3	A	319	EDO	O1-C1-C2-O2
3	A	301	EDO	O1-C1-C2-O2
3	A	314	EDO	O1-C1-C2-O2
3	A	303	EDO	O1-C1-C2-O2
3	A	306	EDO	O1-C1-C2-O2
3	A	311	EDO	O1-C1-C2-O2
3	A	312	EDO	O1-C1-C2-O2
3	A	304	EDO	O1-C1-C2-O2
3	A	305	EDO	O1-C1-C2-O2
3	A	315	EDO	O1-C1-C2-O2
3	A	318	EDO	O1-C1-C2-O2
3	B	301	EDO	O1-C1-C2-O2
3	B	307	EDO	O1-C1-C2-O2
3	A	309	EDO	O1-C1-C2-O2
3	B	302	EDO	O1-C1-C2-O2

There are no ring outliers.

10 monomers are involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	313	EDO	1	0
3	B	302	EDO	1	0
3	A	318	EDO	1	0
3	A	316	EDO	5	0
3	B	304	EDO	5	0
3	A	309	EDO	1	0
3	A	301	EDO	3	0
3	A	302	EDO	1	0
3	B	306	EDO	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	317	EDO	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	A	223/299 (74%)	-0.19	8 (3%) 46 48	15, 36, 80, 93	9 (4%)
1	B	203/299 (67%)	0.30	12 (5%) 28 30	18, 48, 94, 102	14 (6%)
All	All	426/598 (71%)	0.04	20 (4%) 36 38	15, 41, 88, 102	23 (5%)

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	72	HIS	5.3
1	B	16	VAL	3.2
1	A	152	SER	3.2
1	B	15	VAL	3.1
1	B	17	GLY	3.0
1	B	153	ARG	2.8
1	B	152	SER	2.8
1	B	65	THR	2.7
1	A	72	HIS	2.7
1	B	223	HIS	2.6
1	B	33	PHE	2.5
1	A	225	SER	2.5
1	A	224	PHE	2.4
1	A	131	GLN	2.2
1	B	136	THR	2.1
1	A	130	ARG	2.1
1	A	28	SER	2.1
1	B	63	ILE	2.0
1	B	13	THR	2.0
1	A	75	ASP	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 ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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	EDO	A	314	4/4	0.59	0.13	77,82,85,86	0
3	EDO	A	308	4/4	0.70	0.12	75,75,75,76	0
3	EDO	A	319	4/4	0.70	0.18	80,82,83,85	0
3	EDO	B	307	4/4	0.75	0.23	85,89,95,98	0
3	EDO	A	315	4/4	0.76	0.17	90,92,92,92	0
3	EDO	A	307	4/4	0.77	0.15	62,69,75,75	0
3	EDO	A	306	4/4	0.77	0.19	72,75,75,76	0
3	EDO	A	313	4/4	0.80	0.17	67,71,73,77	0
3	EDO	B	305	4/4	0.83	0.24	56,61,65,70	0
3	EDO	A	311	4/4	0.83	0.23	48,49,49,49	4
3	EDO	A	302	4/4	0.84	0.15	52,60,64,64	0
3	EDO	B	301	4/4	0.84	0.11	64,66,66,68	0
3	EDO	B	306	4/4	0.86	0.11	74,76,76,79	0
3	EDO	A	309	4/4	0.86	0.23	63,63,65,65	0
3	EDO	A	304	4/4	0.87	0.14	57,58,64,64	0
4	PO4	A	320	5/5	0.87	0.09	79,80,83,84	5
3	EDO	A	312	4/4	0.88	0.22	61,62,64,66	0
3	EDO	A	305	4/4	0.88	0.24	50,51,52,53	0
3	EDO	A	316	4/4	0.88	0.23	51,54,62,69	0
3	EDO	A	317	4/4	0.89	0.13	60,61,61,65	0
3	EDO	A	318	4/4	0.89	0.22	57,62,65,68	0
3	EDO	A	301	4/4	0.90	0.25	49,51,51,52	0
3	EDO	B	303	4/4	0.91	0.15	47,55,59,65	0
3	EDO	B	302	4/4	0.92	0.12	47,54,59,68	0
3	EDO	A	310	4/4	0.92	0.17	54,61,64,68	0
5	CIT	A	321	13/13	0.93	0.10	44,51,65,66	4
3	EDO	B	304	4/4	0.94	0.14	44,48,50,51	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	EDO	A	303	4/4	0.95	0.12	66,70,74,77	0
2	MN3	B	300	1/1	0.99	0.03	47,47,47,47	0
2	MN3	A	300	1/1	1.00	0.02	28,28,28,28	0

6.5 Other polymers [i](#)

There are no such residues in this entry.