



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 09:29 am GMT

PDB ID : 2CY6
Title : Crystal structure of ConM in complex with trehalose and maltose
Authors : Cavada, B.S.; Azevedo Jr., W.F.; Delatorre, P.; Rocha, B.A.M.; Souza, E.P.; Gadelha, C.A.A.
Deposited on : 2005-07-05
Resolution : 2.00 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

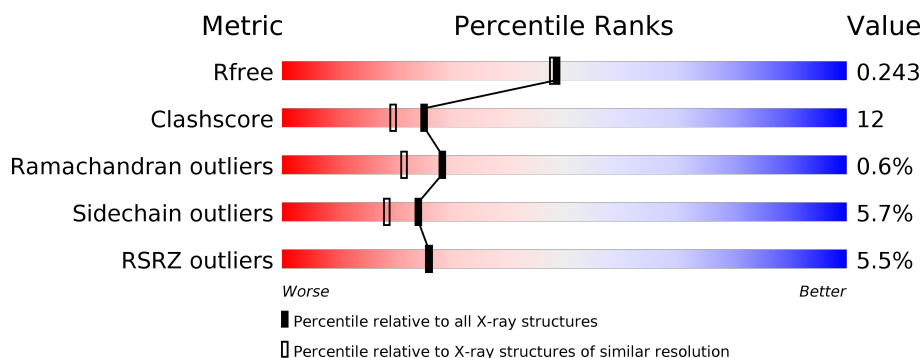
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	100719	6609 (2.00-2.00)
Clashscore	112137	7775 (2.00-2.00)
Ramachandran outliers	110173	7679 (2.00-2.00)
Sidechain outliers	110143	7678 (2.00-2.00)
RSRZ outliers	101464	6696 (2.00-2.00)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	237	<div> <div>5%</div> <div> <div></div> <div>74%</div> <div>21%</div> <div>5%</div> </div> </div>
1	D	237	<div> <div>6%</div> <div> <div></div> <div>78%</div> <div>16%</div> <div>5%</div> </div> </div>

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	TRE	A	5676	-	-	X	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3852 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 Lectin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	237	Total	C	N	O	S	0	0	0
			1800	1134	302	363	1			
1	D	237	Total	C	N	O	S	0	0	0
			1800	1134	302	363	1			

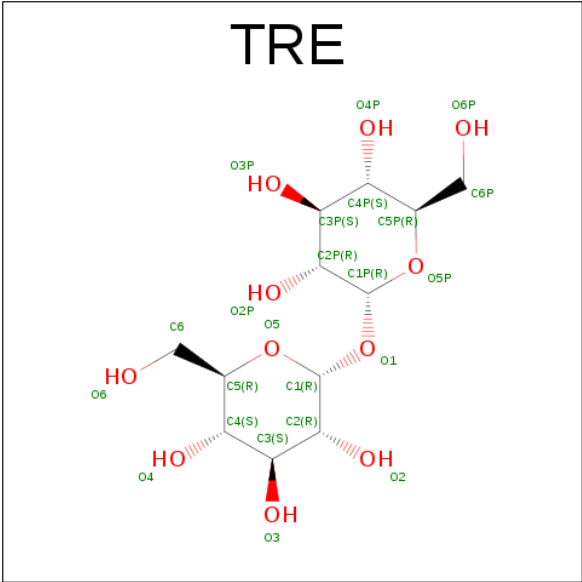
- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Ca	0	0
			1	1		
2	D	1	Total	Ca	0	0
			1	1		

- Molecule 3 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mn	0	0
			1	1		
3	D	1	Total	Mn	0	0
			1	1		

- Molecule 4 is TREHALOSE (three-letter code: TRE) (formula: C₁₂H₂₂O₁₁).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	D	1	Total	C	O	0	0
			23	12	11		
4	A	1	Total	C	O	0	0
			23	12	11		

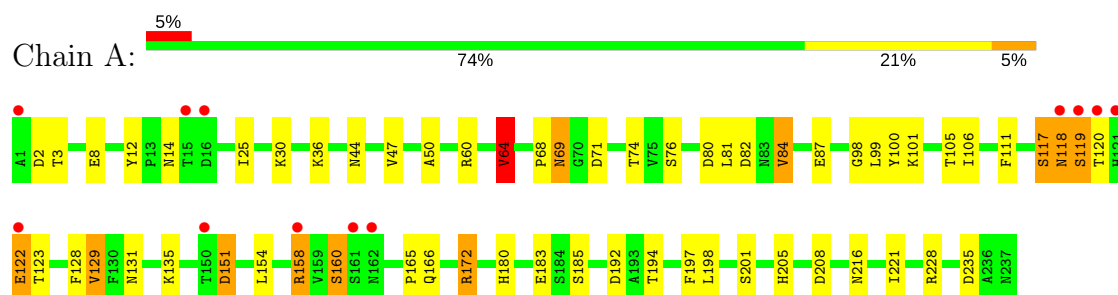
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	104	Total	O	0	0
			104	104		
5	D	98	Total	O	0	0
			98	98		

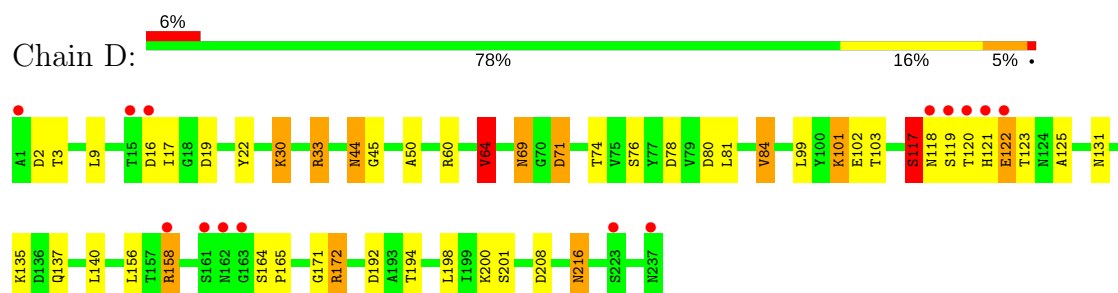
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: Lectin



• Molecule 1: Lectin



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2	Depositor
Cell constants a, b, c, α , β , γ	67.03Å 97.33Å 71.09Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	9.96 – 2.00 9.96 – 2.00	Depositor EDS
% Data completeness (in resolution range)	97.9 (9.96-2.00) 97.9 (9.96-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.75 (at 2.01Å)	Xtriage
Refinement program	REFMAC 5.2.0003	Depositor
R, R_{free}	0.185 , 0.239 0.194 , 0.243	Depositor DCC
R_{free} test set	1580 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	22.2	Xtriage
Anisotropy	0.059	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.52 , 53.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3852	wwPDB-VP
Average B, all atoms (Å ²)	23.0	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 55.12 % 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 3.3128e-05. 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CA, MN, TRE

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	1.11	2/1841 (0.1%)	1.09	9/2507 (0.4%)
1	D	1.08	2/1841 (0.1%)	1.10	9/2507 (0.4%)
All	All	1.09	4/3682 (0.1%)	1.10	18/5014 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	D	0	2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	129	VAL	CB-CG1	-6.75	1.38	1.52
1	A	172	ARG	CD-NE	-6.33	1.35	1.46
1	D	22	TYR	CD1-CE1	-5.70	1.30	1.39
1	D	172	ARG	CD-NE	-5.02	1.38	1.46

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	172	ARG	NE-CZ-NH2	-12.75	113.92	120.30
1	D	172	ARG	NE-CZ-NH2	-8.93	115.83	120.30
1	A	64	VAL	CB-CA-C	-7.94	96.32	111.40
1	D	30	LYS	CD-CE-NZ	7.06	127.93	111.70
1	D	71	ASP	CB-CG-OD2	6.98	124.58	118.30
1	D	64	VAL	CB-CA-C	-6.73	98.62	111.40
1	D	80	ASP	CB-CG-OD2	6.67	124.30	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	80	ASP	CB-CG-OD2	6.62	124.26	118.30
1	A	172	ARG	NE-CZ-NH1	6.54	123.57	120.30
1	A	158	ARG	NE-CZ-NH1	6.46	123.53	120.30
1	D	19	ASP	CB-CG-OD2	6.34	124.00	118.30
1	D	78	ASP	CB-CG-OD2	6.20	123.88	118.30
1	A	64	VAL	CG1-CB-CG2	6.07	120.62	110.90
1	A	82	ASP	CB-CG-OD2	5.87	123.58	118.30
1	D	16	ASP	CB-CG-OD2	5.63	123.37	118.30
1	A	235	ASP	CB-CG-OD2	5.46	123.21	118.30
1	A	151	ASP	CB-CG-OD2	5.17	122.95	118.30
1	D	33	ARG	NE-CZ-NH2	-5.17	117.72	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	117	SER	Peptide
1	D	122	GLU	Peptide

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	1800	0	1744	50	0
1	D	1800	0	1744	42	0
2	A	1	0	0	0	0
2	D	1	0	0	0	0
3	A	1	0	0	0	0
3	D	1	0	0	0	0
4	A	23	0	22	14	0
4	D	23	0	22	2	0
5	A	104	0	0	5	1
5	D	98	0	0	9	0
All	All	3852	0	3532	87	1

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 12.

All (87) 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:14:ASN:HD21	4:A:5676:TRE:C6	1.45	1.30
1:A:14:ASN:HD21	4:A:5676:TRE:HC61	1.23	1.01
1:A:14:ASN:ND2	4:A:5676:TRE:C6	2.27	0.98
1:D:71:ASP:HB2	5:D:4688:HOH:O	1.66	0.96
1:A:44:ASN:HD21	1:A:201:SER:H	1.05	0.93
1:A:14:ASN:ND2	4:A:5676:TRE:HC61	1.83	0.92
1:D:44:ASN:HD21	1:D:201:SER:H	1.18	0.90
1:A:120:THR:OG1	5:A:5779:HOH:O	2.03	0.75
1:A:166:GLN:HG3	5:A:5717:HOH:O	1.87	0.74
1:A:123:THR:H	1:D:131:ASN:HD22	1.33	0.73
1:D:120:THR:O	1:D:122:GLU:N	2.20	0.71
1:A:12:TYR:CD2	4:A:5676:TRE:O6	2.44	0.70
1:A:44:ASN:ND2	1:A:201:SER:H	1.87	0.69
1:A:208:ASP:OD2	4:A:5676:TRE:HC5	1.92	0.69
1:A:3:THR:H	1:A:216:ASN:ND2	1.90	0.69
1:D:3:THR:HG23	1:D:30:LYS:CE	2.25	0.67
1:D:3:THR:HG23	1:D:30:LYS:HE3	1.77	0.66
1:D:208:ASP:OD1	4:D:4676:TRE:H6P2	1.98	0.64
1:A:44:ASN:HD21	1:A:201:SER:N	1.87	0.63
1:A:98:GLY:HA3	4:A:5676:TRE:HC3	1.82	0.61
1:D:60:ARG:CZ	5:D:4690:HOH:O	2.49	0.61
1:A:99:LEU:HB2	4:A:5676:TRE:O3	2.01	0.61
1:A:172:ARG:HD2	1:A:221:ILE:HG13	1.83	0.61
1:D:3:THR:H	1:D:216:ASN:ND2	1.99	0.61
1:A:3:THR:H	1:A:216:ASN:HD22	1.50	0.59
1:D:64:VAL:HG13	1:D:74:THR:OG1	2.03	0.58
1:D:81:LEU:HA	1:D:84:VAL:HG13	1.86	0.57
1:A:172:ARG:HD2	1:A:221:ILE:CG1	2.35	0.57
1:A:2:ASP:CB	1:A:216:ASN:HD21	2.17	0.57
1:A:99:LEU:H	4:A:5676:TRE:C3	2.17	0.57
1:A:2:ASP:HB3	1:A:216:ASN:HD21	1.69	0.57
1:D:156:LEU:O	1:D:171:GLY:HA3	2.05	0.56
1:D:216:ASN:H	1:D:216:ASN:HD22	1.52	0.56
1:D:44:ASN:ND2	1:D:200:LYS:HA	2.21	0.56
1:A:69:ASN:ND2	1:A:71:ASP:H	2.02	0.55
1:D:71:ASP:CB	5:D:4688:HOH:O	2.37	0.55
1:A:81:LEU:HA	1:A:84:VAL:HG13	1.87	0.55
1:A:160:SER:HB3	1:A:166:GLN:NE2	2.22	0.54
1:A:99:LEU:H	4:A:5676:TRE:HC3	1.72	0.54
1:D:60:ARG:NE	5:D:4690:HOH:O	2.40	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:183:GLU:OE2	1:A:185:SER:CB	2.56	0.54
1:D:9:LEU:HD12	1:D:9:LEU:N	2.23	0.53
1:A:151:ASP:OD2	5:A:5700:HOH:O	2.19	0.53
1:D:60:ARG:NH2	5:D:4690:HOH:O	2.40	0.53
1:A:64:VAL:HG13	1:A:74:THR:OG1	2.08	0.52
1:A:60:ARG:HD3	1:A:76:SER:HB3	1.92	0.51
1:D:3:THR:HG23	1:D:30:LYS:HE2	1.92	0.51
1:D:101:LYS:HD2	1:D:165:PRO:O	2.09	0.51
1:D:172:ARG:NH2	5:D:4685:HOH:O	2.39	0.51
1:D:17:ILE:O	1:D:33:ARG:HG2	2.11	0.50
1:A:101:LYS:HD2	1:A:165:PRO:O	2.12	0.49
1:A:123:THR:O	1:D:131:ASN:ND2	2.43	0.49
1:D:60:ARG:HD3	1:D:76:SER:HB3	1.93	0.49
1:A:129:VAL:HG12	1:D:125:ALA:HB3	1.95	0.49
1:A:111:PHE:HB3	1:A:128:PHE:CZ	2.48	0.49
1:D:118:ASN:ND2	5:D:4760:HOH:O	2.45	0.49
1:D:216:ASN:H	1:D:216:ASN:ND2	2.10	0.49
1:D:2:ASP:HB3	1:D:216:ASN:HD21	1.78	0.49
1:A:117:SER:C	1:A:119:SER:H	2.17	0.48
1:A:14:ASN:ND2	4:A:5676:TRE:O6	2.42	0.48
1:D:69:ASN:ND2	1:D:71:ASP:H	2.13	0.47
1:D:50:ALA:O	1:D:194:THR:HA	2.15	0.47
1:D:2:ASP:CB	1:D:216:ASN:HD21	2.27	0.47
1:D:99:LEU:N	4:D:4676:TRE:H6P1	2.31	0.46
1:A:106:ILE:HB	1:A:154:LEU:HB3	1.99	0.45
1:D:137:GLN:HG2	1:D:140:LEU:HD12	1.99	0.45
1:D:44:ASN:HD21	1:D:201:SER:N	1.98	0.45
1:A:122:GLU:HG3	1:A:122:GLU:O	2.16	0.45
1:D:44:ASN:HD22	1:D:45:GLY:N	2.15	0.44
1:A:87:GLU:CG	1:A:180:HIS:HE2	2.31	0.44
1:D:117:SER:O	1:D:118:ASN:HB3	2.17	0.43
1:A:208:ASP:CG	4:A:5676:TRE:HC5	2.38	0.43
1:A:228:ARG:HB3	4:A:5676:TRE:HC61	2.01	0.43
1:D:192:ASP:HB2	5:D:4704:HOH:O	2.19	0.43
1:A:50:ALA:O	1:A:194:THR:HA	2.18	0.43
1:A:36:LYS:HE3	5:A:5732:HOH:O	2.18	0.42
1:A:131:ASN:HD22	1:D:123:THR:H	1.68	0.42
1:D:101:LYS:HG3	1:D:165:PRO:HG2	2.01	0.42
1:A:30:LYS:HA	5:A:5680:HOH:O	2.20	0.41
1:A:99:LEU:HD22	4:A:5676:TRE:O3	2.20	0.41
1:D:158:ARG:CZ	5:D:4764:HOH:O	2.68	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:105:THR:O	1:A:197:PHE:HA	2.19	0.41
1:A:100:TYR:HB3	1:A:205:HIS:O	2.20	0.41
1:A:8:GLU:O	1:A:25:ILE:HA	2.21	0.41
1:D:102:GLU:HG3	1:D:103:THR:O	2.21	0.40
1:A:47:VAL:O	1:A:68:PRO:HD3	2.21	0.40
1:A:123:THR:H	1:D:131:ASN:ND2	2.10	0.40

All (1) 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)
5:A:5705:HOH:O	5:A:5705:HOH:O[2_655]	2.15	0.05

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	235/237 (99%)	226 (96%)	8 (3%)	1 (0%)	38	33
1	D	235/237 (99%)	222 (94%)	11 (5%)	2 (1%)	20	12
All	All	470/474 (99%)	448 (95%)	19 (4%)	3 (1%)	28	21

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	118	ASN
1	D	121	HIS
1	D	119	SER

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	202/202 (100%)	190 (94%)	12 (6%)	23	17
1	D	202/202 (100%)	191 (95%)	11 (5%)	26	20
All	All	404/404 (100%)	381 (94%)	23 (6%)	24	18

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	64	VAL
1	A	69	ASN
1	A	84	VAL
1	A	117	SER
1	A	118	ASN
1	A	119	SER
1	A	122	GLU
1	A	135	LYS
1	A	158	ARG
1	A	160	SER
1	A	192	ASP
1	A	198	LEU
1	D	44	ASN
1	D	64	VAL
1	D	69	ASN
1	D	84	VAL
1	D	101	LYS
1	D	117	SER
1	D	135	LYS
1	D	158	ARG
1	D	164	SER
1	D	198	LEU
1	D	216	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	GLN
1	A	44	ASN
1	A	69	ASN
1	A	118	ASN
1	A	121	HIS
1	A	131	ASN
1	A	216	ASN
1	A	237	ASN
1	D	43	GLN
1	D	44	ASN
1	D	69	ASN
1	D	131	ASN
1	D	166	GLN
1	D	216	ASN
1	D	237	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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
4	TRE	A	5676	-	24,24,24	1.50	4 (16%)	35,35,35	3.23	17 (48%)
4	TRE	D	4676	-	24,24,24	0.68	0	35,35,35	1.53	8 (22%)

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	TRE	A	5676	-	-	0/8/48/48	0/2/2/2
4	TRE	D	4676	-	-	0/8/48/48	0/2/2/2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	5676	TRE	O1-C1P	2.17	1.47	1.41
4	A	5676	TRE	C3-C2	2.40	1.58	1.52
4	A	5676	TRE	O5P-C1P	2.69	1.48	1.41
4	A	5676	TRE	O4-C4	3.14	1.50	1.43

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	5676	TRE	O3-C3-C4	-8.07	92.80	110.36
4	A	5676	TRE	C1P-C2P-C3P	-4.16	102.24	109.98
4	A	5676	TRE	O6-C6-C5	-3.91	98.19	111.34
4	A	5676	TRE	O5-C1-C2	-3.33	103.86	110.30
4	A	5676	TRE	C3-C4-C5	-3.00	104.94	110.22
4	D	4676	TRE	C6P-C5P-C4P	-2.49	107.16	113.00
4	A	5676	TRE	C4P-C3P-C2P	-2.43	106.55	110.84
4	D	4676	TRE	C4-C3-C2	-2.23	106.91	110.84
4	D	4676	TRE	O5-C5-C6	2.01	111.23	106.41
4	A	5676	TRE	O4-C4-C3	2.18	115.10	110.36
4	A	5676	TRE	O2P-C2P-C3P	2.26	115.28	110.36
4	D	4676	TRE	C3-C4-C5	2.34	114.33	110.22
4	D	4676	TRE	O3P-C3P-C4P	2.36	115.49	110.36
4	A	5676	TRE	C1-C2-C3	2.47	114.56	109.98
4	D	4676	TRE	C1P-O5P-C5P	2.49	118.41	113.72
4	A	5676	TRE	C1P-O1-C1	2.60	119.06	114.37
4	D	4676	TRE	C3P-C4P-C5P	2.91	115.34	110.22
4	A	5676	TRE	O2P-C2P-C1P	3.04	116.39	110.03
4	A	5676	TRE	O5P-C5P-C6P	3.77	115.45	106.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	4676	TRE	O5P-C5P-C4P	3.88	116.80	109.66
4	A	5676	TRE	C1-O5-C5	3.97	121.20	113.72
4	A	5676	TRE	O5-C5-C4	4.55	118.05	109.66
4	A	5676	TRE	O1-C1-C2	4.74	118.80	108.11
4	A	5676	TRE	C1P-O5P-C5P	5.35	123.80	113.72
4	A	5676	TRE	C4-C3-C2	9.37	127.37	110.84

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	5676	TRE	14	0
4	D	4676	TRE	2	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	237/237 (100%)	-0.03	12 (5%)	29 29	12, 20, 36, 58	0
1	D	237/237 (100%)	0.04	14 (5%)	23 23	12, 22, 40, 59	0
All	All	474/474 (100%)	0.00	26 (5%)	26 26	12, 21, 38, 59	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	120	THR	4.9
1	D	118	ASN	4.8
1	D	162	ASN	4.7
1	A	118	ASN	4.7
1	D	1	ALA	4.0
1	D	122	GLU	4.0
1	D	16	ASP	3.9
1	A	122	GLU	3.8
1	A	150	THR	3.8
1	D	121	HIS	3.7
1	A	119	SER	3.5
1	A	162	ASN	3.5
1	A	1	ALA	3.4
1	D	161	SER	3.3
1	D	119	SER	3.3
1	A	16	ASP	3.1
1	A	120	THR	3.1
1	A	121	HIS	3.0
1	D	163	GLY	2.9
1	A	161	SER	2.6
1	D	237	ASN	2.4
1	D	223	SER	2.2
1	D	158	ARG	2.2
1	A	158	ARG	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	15	THR	2.0
1	D	15	THR	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
4	TRE	A	5676	23/23	0.57	0.45	9.73	23,48,60,63	0
4	TRE	D	4676	23/23	0.81	0.25	0.86	22,33,39,41	0
2	CA	D	1003	1/1	0.97	0.05	-2.78	19,19,19,19	0
3	MN	D	240	1/1	0.97	0.05	-3.01	22,22,22,22	0
2	CA	A	1003	1/1	0.99	0.04	-3.34	18,18,18,18	0
3	MN	A	240	1/1	0.99	0.03	-3.89	20,20,20,20	0

6.5 Other polymers [i](#)

There are no such residues in this entry.