



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 11:50 AM GMT

PDB ID : 1KB0
Title : Crystal Structure of Quinohemoprotein Alcohol Dehydrogenase from Comamonas testosteroni
Authors : Rozeboom, H.J.; Oubrie, A.
Deposited on : 2001-11-05
Resolution : 1.44 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

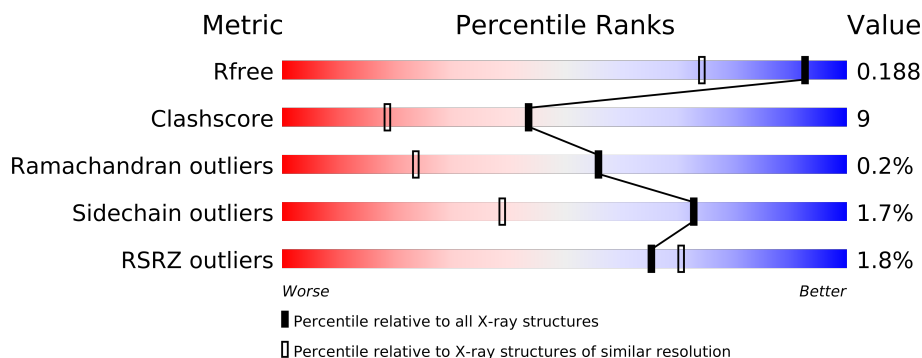
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance


The reported resolution of this entry is 1.44 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}	66092	1819 (1.48-1.40)
Clashscore	79885	2050 (1.48-1.40)
Ramachandran outliers	78287	2002 (1.48-1.40)
Sidechain outliers	78261	2001 (1.48-1.40)
RSRZ outliers	66119	1819 (1.48-1.40)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	677	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
4	TFB	A	1810	-	X
6	GOL	A	1820	-	X
6	GOL	A	1823	-	X
6	GOL	A	1824	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 6286 atoms, of which 0 are hydrogen and 0 are deuterium.

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 quinohemoprotein alcohol dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	670	Total	C	N	O	S	0	6	0
			5172	3290	903	960	19			

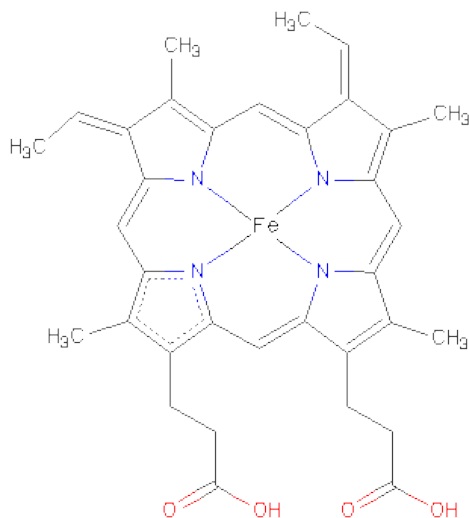
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	181	ASN	LYS	see remark 999	UNP Q46444
A	182	GLY	ARG	see remark 999	UNP Q46444
A	566	GLY	ALA	see remark 999	UNP Q46444
A	512	TRO	TRP	MODIFIED RESIDUE	UNP Q46444

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

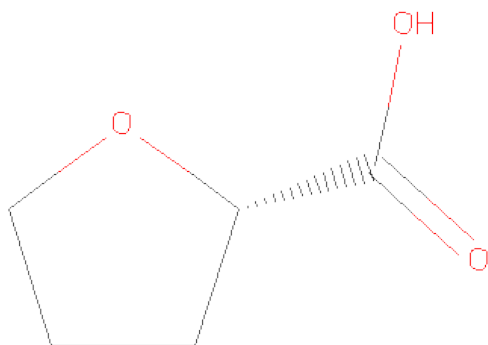
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Ca	0	0
			1	1		

- Molecule 3 is HEME C (three-letter code: HEC) (formula: C₃₄H₃₄FeN₄O₄).



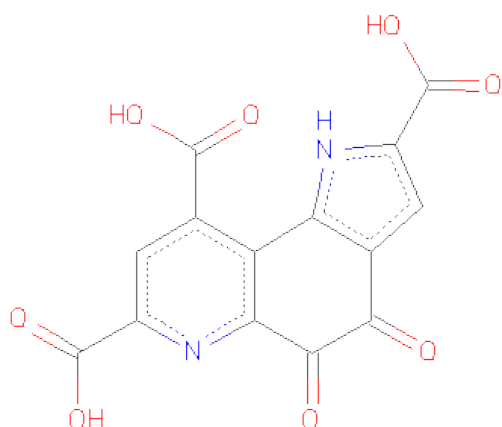
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 4 is TETRAHYDROFURAN-2-CARBOXYLICACID (three-letter code: TFB) (formula: $C_5H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			8	5	3		

- Molecule 5 is PYRROLOQUINOLINE QUINONE (three-letter code: PQQ) (formula: $C_{14}H_6N_2O_8$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			24	14	2	8		

- Molecule 6 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 7 is water.

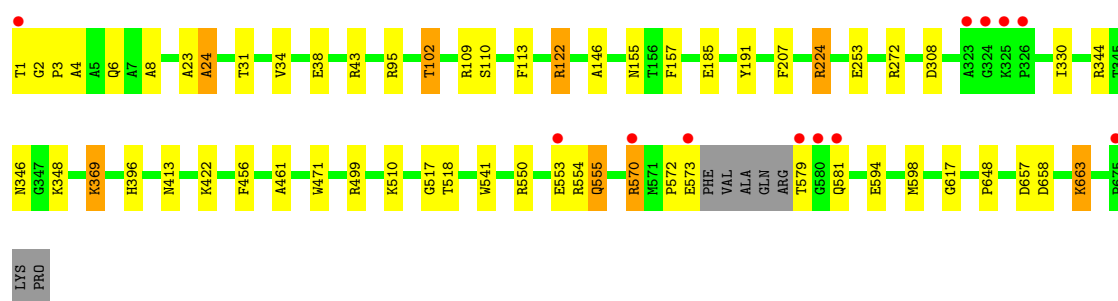
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	1016	Total 1020	O 1020	0	5

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: quinoheprotein alcohol dehydrogenase

Chain A: 



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	97.92Å 74.33Å 92.33Å 90.00° 105.67° 90.00°	Depositor
Resolution (Å)	29.63 – 1.44 29.63 – 1.44	Depositor EDS
% Data completeness (in resolution range)	96.3 (29.63-1.44) 96.5 (29.63-1.44)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.60 (at 1.44Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.160 , 0.188 0.162 , 0.188	Depositor DCC
R_{free} test set	5671 reflections (5.38%)	DCC
Wilson B-factor (Å ²)	16.4	Xtriage
Anisotropy	0.025	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 50.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 111119 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	6286	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.57% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, TRO, CA, TFB, PQQ, HEC

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	0.82	0/5307	0.96	4/7227 (0.1%)

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	A	0	1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	272	ARG	NE-CZ-NH1	-6.35	117.12	120.30
1	A	272	ARG	NE-CZ-NH2	5.95	123.27	120.30
1	A	157	PHE	CB-CG-CD1	-5.41	117.02	120.80
1	A	122	ARG	NE-CZ-NH2	-5.16	117.72	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	122	ARG	Sidechain

5.2 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 hydrogens

added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5172	0	4979	86	0
2	A	1	0	0	0	0
3	A	43	0	30	0	0
4	A	8	0	7	3	0
5	A	24	0	3	0	0
6	A	18	0	23	14	0
7	A	1020	0	0	40	0
All	All	6286	0	5042	90	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 9.

All (90) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:594:GLU:HG3	7:A:2679:HOH:O	1.70	0.92
1:A:253:GLU:HG3	7:A:2235:HOH:O	1.72	0.89
1:A:663:LYS:NZ	1:A:663:LYS:HB3	1.93	0.83
1:A:550:ARG:HH11	1:A:579:THR:HG23	1.45	0.80
1:A:541:TRP:HD1	6:A:1823:GOL:H2	1.46	0.80
6:A:1823:GOL:H11	7:A:2575:HOH:O	1.81	0.79
1:A:553:GLU:HG2	1:A:579:THR:HG22	1.63	0.79
1:A:541:TRP:CD1	6:A:1823:GOL:H2	2.17	0.78
1:A:663:LYS:HE3	7:A:2788:HOH:O	1.85	0.77
1:A:224:ARG:HG3	7:A:2487:HOH:O	1.86	0.75
1:A:581:GLN:HG3	7:A:2334:HOH:O	1.88	0.72
1:A:570:ARG:H	1:A:570:ARG:NE	1.88	0.72
1:A:553:GLU:HG2	1:A:579:THR:CG2	2.20	0.72
1:A:43:ARG:NH2	7:A:2677:HOH:O	2.23	0.71
1:A:554:ARG:HD3	7:A:2561:HOH:O	1.92	0.69
1:A:658:ASP:HB2	7:A:2538:HOH:O	1.93	0.68
1:A:102:THR:HG22	7:A:2255:HOH:O	1.94	0.67
1:A:517:GLY:HA2	1:A:554:ARG:HE	1.59	0.67
1:A:663:LYS:HZ3	1:A:663:LYS:HB3	1.58	0.67
1:A:663:LYS:NZ	1:A:663:LYS:CB	2.58	0.67
1:A:570:ARG:H	1:A:570:ARG:CD	2.09	0.66
1:A:517:GLY:O	1:A:554:ARG:NE	2.28	0.66
1:A:598[B]:MET:HE1	7:A:2797:HOH:O	1.95	0.66
1:A:155:ASN:HB3	7:A:2589:HOH:O	1.95	0.65

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:553:GLU:C	7:A:2798:HOH:O	2.34	0.65
1:A:23[B]:ALA:HA	7:A:2548:HOH:O	1.96	0.64
1:A:369:LYS:HE2	7:A:2540:HOH:O	1.98	0.63
1:A:510:LYS:HD3	6:A:1820:GOL:H31	1.79	0.63
1:A:330:ILE:HG13	1:A:344:ARG:HB3	1.82	0.61
1:A:572:PRO:O	1:A:573:GLU:HB3	2.02	0.60
1:A:43:ARG:NH2	7:A:2317:HOH:O	2.26	0.60
1:A:23[B]:ALA:O	1:A:43:ARG:CZ	2.50	0.59
1:A:422:LYS:HD3	7:A:2796:HOH:O	2.02	0.59
1:A:517:GLY:CA	1:A:554:ARG:HE	2.14	0.59
1:A:550:ARG:NH1	1:A:579:THR:HG23	2.17	0.59
1:A:554:ARG:N	7:A:2798:HOH:O	2.34	0.59
1:A:663:LYS:HZ2	1:A:663:LYS:HB3	1.67	0.57
1:A:663:LYS:CE	7:A:2788:HOH:O	2.48	0.56
1:A:663:LYS:HZ2	1:A:663:LYS:CB	2.20	0.55
1:A:346:ASN:ND2	1:A:348:LYS:HE3	2.21	0.55
1:A:102:THR:HG21	7:A:2592:HOH:O	2.09	0.53
1:A:3:PRO:HA	1:A:6:GLN:HE21	1.74	0.53
4:A:1810:TFB:HD1	7:A:2840:HOH:O	2.09	0.53
1:A:8:ALA:HB2	1:A:146:ALA:HB2	1.90	0.52
1:A:517:GLY:O	1:A:554:ARG:CG	2.58	0.51
1:A:23[B]:ALA:O	1:A:43:ARG:NE	2.44	0.50
1:A:308:ASP:OD2	4:A:1810:TFB:OXT	2.28	0.50
1:A:555:GLN:OE1	6:A:1823:GOL:C3	2.60	0.49
1:A:573:GLU:HG2	7:A:2837:HOH:O	2.12	0.49
1:A:2:GLY:N	1:A:3:PRO:CD	2.76	0.49
1:A:43:ARG:CZ	7:A:2677:HOH:O	2.60	0.49
1:A:517:GLY:C	1:A:554:ARG:HE	2.15	0.49
1:A:617:GLY:O	6:A:1823:GOL:O2	2.30	0.49
1:A:517:GLY:HA2	1:A:554:ARG:NE	2.24	0.48
1:A:617:GLY:C	6:A:1823:GOL:O2	2.51	0.48
1:A:31:THR:OG1	7:A:2680:HOH:O	2.20	0.48
1:A:555:GLN:CD	6:A:1823:GOL:H32	2.34	0.48
1:A:657[B]:ASP:HA	7:A:2015:HOH:O	2.14	0.47
1:A:598[A]:MET:HE3	7:A:2732:HOH:O	2.12	0.47
1:A:550:ARG:HD2	6:A:1823:GOL:H31	1.97	0.47
1:A:555:GLN:OE1	6:A:1823:GOL:H32	2.14	0.47
1:A:422:LYS:CD	7:A:2796:HOH:O	2.60	0.46
1:A:4:ALA:HB2	7:A:2392:HOH:O	2.13	0.46
1:A:517:GLY:O	1:A:554:ARG:HG3	2.15	0.46
1:A:553:GLU:CB	7:A:2798:HOH:O	2.64	0.46
1:A:413:ASN:HB2	1:A:456:PHE:CE2	2.50	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:A:1810:TFB:CD	7:A:2840:HOH:O	2.64	0.45
1:A:369:LYS:HG3	7:A:2643:HOH:O	2.15	0.45
6:A:1823:GOL:C1	7:A:2575:HOH:O	2.53	0.45
1:A:570:ARG:CD	1:A:570:ARG:N	2.77	0.44
1:A:43:ARG:NH1	7:A:2677:HOH:O	2.50	0.44
1:A:570:ARG:H	1:A:570:ARG:HD2	1.80	0.43
1:A:110:SER:HA	7:A:1941:HOH:O	2.19	0.43
1:A:113:PHE:HB3	1:A:648:PRO:HG3	2.01	0.43
1:A:396:HIS:CD2	1:A:396:HIS:H	2.35	0.43
1:A:541:TRP:HB3	6:A:1823:GOL:C1	2.48	0.43
1:A:38:GLU:HG2	7:A:2680:HOH:O	2.19	0.43
1:A:1:THR:HB	1:A:2:GLY:H	1.45	0.43
1:A:663:LYS:CD	7:A:2788:HOH:O	2.66	0.42
1:A:553:GLU:HB3	7:A:2798:HOH:O	2.18	0.42
1:A:461:ALA:HB2	1:A:471:TRP:CE2	2.54	0.42
1:A:579:THR:N	7:A:2800:HOH:O	2.52	0.42
1:A:1:THR:HG22	1:A:2:GLY:N	2.35	0.41
1:A:499:ARG:HH12	6:A:1820:GOL:H2	1.86	0.41
1:A:23[B]:ALA:CA	1:A:24[B]:ALA:HB2	2.51	0.40
1:A:191:TYR:HB3	1:A:207:PHE:CD2	2.57	0.40
1:A:541:TRP:HB3	6:A:1823:GOL:H12	2.04	0.40
1:A:23[B]:ALA:N	1:A:24[B]:ALA:HB2	2.36	0.40
1:A:109:ARG:NH2	7:A:2519:HOH:O	2.37	0.40
1:A:95:ARG:HD2	7:A:2192:HOH:O	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

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	671/677 (99%)	644 (96%)	25 (4%)	2 (0%)	50 19

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	24[A]	ALA
1	A	24[B]	ALA

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	521/523 (100%)	512 (98%)	9 (2%)	73 39

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	34	VAL
1	A	102	THR
1	A	185	GLU
1	A	224	ARG
1	A	369	LYS
1	A	518	THR
1	A	555	GLN
1	A	570	ARG
1	A	663	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	GLN
1	A	396	HIS

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

1 non-standard protein/DNA/RNA residue is modelled in this entry.

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	$\# Z > 2$
1	TRO	A	512	1	16,16,17	4.77	4 (25%)	19,22,24	1.82	3 (15%)

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
1	TRO	A	512	1	-	0/4/6/8	0/0/2/2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	512	TRO	O-C	18.42	1.24	1.11
1	A	512	TRO	CB-CG	2.72	1.55	1.51
1	A	512	TRO	CA-C	2.49	1.52	1.48
1	A	512	TRO	CG-CD2	2.18	1.44	1.41

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	512	TRO	C-CA-N	-4.42	109.42	113.83
1	A	512	TRO	OD1-CD1-CG	-4.25	124.09	131.21
1	A	512	TRO	CG-CD1-NE1	3.60	110.78	107.76

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 7 ligands modelled in this entry, 1 is monoatomic - leaving 6 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
5	PQQ	A	1800	2	26,26,26	3.37	12 (46%)	39,40,40	2.34	16 (41%)
4	TFB	A	1810	2	8,8,8	3.68	2 (25%)	10,10,10	1.32	1 (10%)
6	GOL	A	1820	-	5,5,5	1.39	1 (20%)	5,5,5	1.20	1 (20%)
6	GOL	A	1823	-	5,5,5	0.88	0	5,5,5	0.57	0
6	GOL	A	1824	-	5,5,5	0.63	0	5,5,5	0.47	0
3	HEC	A	802	1	50,50,50	2.01	8 (16%)	56,82,82	2.03	8 (14%)

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
5	PQQ	A	1800	2	-	0/10/28/28	0/0/3/3
4	TFB	A	1810	2	-	0/4/11/11	0/1/1/1
6	GOL	A	1820	-	-	0/4/4/4	0/0/0/0
6	GOL	A	1823	-	-	0/4/4/4	0/0/0/0
6	GOL	A	1824	-	-	0/4/4/4	0/0/0/0
3	HEC	A	802	1	-	0/10/54/54	0/0/8/8

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1810	TFB	CA-C	-10.01	1.38	1.51
3	A	802	HEC	C3C-CAC	8.74	1.53	1.35
5	A	1800	PQQ	C2-C2X	-7.45	1.42	1.51
3	A	802	HEC	C3B-CAB	7.05	1.50	1.35
5	A	1800	PQQ	C7-C7X	-6.21	1.42	1.50
5	A	1800	PQQ	O4-C4	5.90	1.36	1.23
5	A	1800	PQQ	C3A-C4	-5.07	1.38	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1800	PQQ	O5-C5	4.88	1.34	1.23
5	A	1800	PQQ	C9-C9A	4.82	1.46	1.40
5	A	1800	PQQ	C5-C4	-4.12	1.41	1.53
5	A	1800	PQQ	C3-C3A	3.53	1.45	1.39
5	A	1800	PQQ	C9A-C1A	-3.52	1.42	1.49
5	A	1800	PQQ	C9A-C6A	3.36	1.43	1.40
5	A	1800	PQQ	C2-N1	3.18	1.39	1.36
3	A	802	HEC	C4C-NC	3.17	1.41	1.37
3	A	802	HEC	C3C-C4C	-2.53	1.38	1.41
3	A	802	HEC	C4D-C3D	-2.50	1.39	1.43
5	A	1800	PQQ	C3-C2	-2.36	1.36	1.39
6	A	1820	GOL	O2-C2	-2.35	1.36	1.43
3	A	802	HEC	FE-ND	2.26	2.02	1.92
3	A	802	HEC	CMD-C2D	2.20	1.56	1.51
4	A	1810	TFB	OA-CD	2.16	1.49	1.42
3	A	802	HEC	C4A-NA	2.03	1.39	1.37

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1800	PQQ	C9-C9A-C1A	8.04	129.81	122.52
3	A	802	HEC	CBB-CAB-C3B	-6.42	110.37	128.44
3	A	802	HEC	C3C-C2C-C1C	-5.78	103.22	107.07
3	A	802	HEC	CBC-CAC-C3C	-5.64	112.55	128.44
3	A	802	HEC	C2C-C1C-NC	5.30	113.41	109.41
3	A	802	HEC	C4C-C3C-C2C	4.26	109.40	106.68
5	A	1800	PQQ	O5-C5-C6A	-3.95	118.12	122.24
5	A	1800	PQQ	O9B-C9X-C9	3.78	126.75	115.47
3	A	802	HEC	CBA-CAA-C2A	-3.68	105.48	112.35
4	A	1810	TFB	CB-CA-C	3.48	116.74	112.55
5	A	1800	PQQ	C9A-C6A-N6	3.14	128.06	123.37
5	A	1800	PQQ	C9-C9A-C6A	-3.03	113.05	118.53
3	A	802	HEC	CMC-C2C-C3C	2.99	129.02	125.72
5	A	1800	PQQ	C1A-C9A-C6A	-2.93	117.14	119.78
5	A	1800	PQQ	C8-C7-C7X	2.92	124.92	119.55
5	A	1800	PQQ	O9B-C9X-O9A	-2.88	116.81	123.35
5	A	1800	PQQ	C5-C6A-N6	-2.73	111.92	116.23
5	A	1800	PQQ	C7X-C7-N6	-2.66	112.63	116.51
5	A	1800	PQQ	C3A-C3-C2	-2.65	106.39	109.86
5	A	1800	PQQ	C9A-C9-C9X	2.61	125.44	121.77
6	A	1820	GOL	O2-C2-C1	2.57	119.93	108.22
5	A	1800	PQQ	O7B-C7X-C7	2.35	120.04	114.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1800	PQQ	O9A-C9X-C9	-2.33	115.96	121.82
3	A	802	HEC	C3A-C4A-NA	2.27	111.12	109.41
5	A	1800	PQQ	O2B-C2X-C2	2.19	118.81	113.39
5	A	1800	PQQ	C8-C9-C9A	-2.09	117.89	120.72

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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	670/677 (98%)	-0.27	12 (1%) 65 70	11, 17, 29, 50	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	579	THR	6.0
1	A	1	THR	5.8
1	A	675	PRO	3.2
1	A	580	GLY	3.1
1	A	324	GLY	3.0
1	A	553	GLU	2.7
1	A	573	GLU	2.7
1	A	581	GLN	2.6
1	A	326	PRO	2.4
1	A	570	ARG	2.4
1	A	323	ALA	2.3
1	A	325	LYS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
1	TRO	A	512	15/16	0.06	0.03	14,17,20,22	0

6.3 Carbohydrates ⓘ

There are no carbohydrates 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. 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
6	GOL	A	1820	6/6	0.15	13.63	17,24,29,32	0
6	GOL	A	1824	6/6	0.17	6.27	57,58,58,59	0
4	TFB	A	1810	8/8	0.16	6.14	27,31,32,33	0
6	GOL	A	1823	6/6	0.27	3.51	36,38,39,39	0
3	HEC	A	802	43/43	0.07	0.72	11,13,24,29	0
5	PQQ	A	1800	24/24	0.07	-0.53	10,11,13,14	0
2	CA	A	801	1/1	0.04	-1.85	21,21,21,21	0

6.5 Other polymers ⓘ

There are no such residues in this entry.