



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:52 AM GMT

PDB ID : 2V4A
Title : CRYSTAL STRUCTURE OF THE SEMET-LABELED PROLYL-4 HYDROXYLASE (P4H) TYPE I FROM GREEN ALGAE CHLAMYDOMONAS REINHARDTII.
Authors : Koski, M.K.; Hieta, R.; Bollner, C.; Kivirikko, K.I.; Myllyharju, J.; Wierenga, R.K.
Deposited on : 2007-06-28
Resolution : 1.93 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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

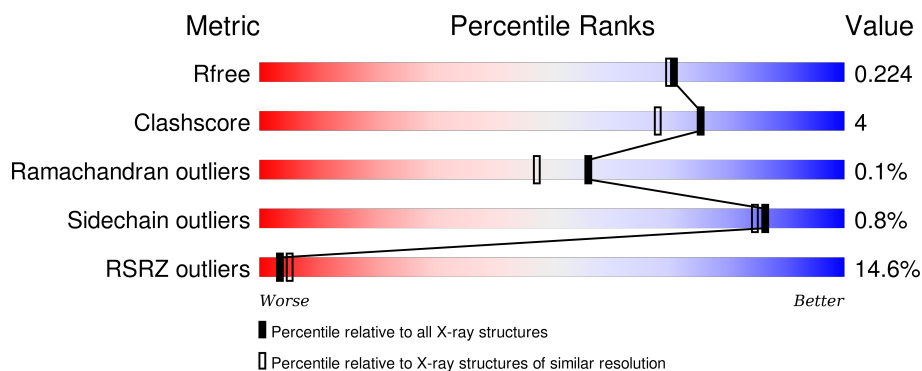
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.93 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}	91344	2910 (1.96-1.92)
Clashscore	102246	3095 (1.96-1.92)
Ramachandran outliers	100387	3062 (1.96-1.92)
Sidechain outliers	100360	3062 (1.96-1.92)
RSRZ outliers	91569	2915 (1.96-1.92)

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	233	<div> <div>7%</div> <div>73%</div> <div>22%</div> </div>
1	B	233	<div> <div>7%</div> <div>72%</div> <div>25%</div> </div>
1	C	233	<div> <div>14%</div> <div>75%</div> <div>5%</div> <div>20%</div> </div>
1	D	233	<div> <div>17%</div> <div>68%</div> <div>14%</div> <div>17%</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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	GOL	A	1251	-	-	-	X
3	GOL	B	1251	-	-	-	X
5	DMS	B	1252	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6127 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 PROLYL-4 HYDROXYLASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	182	Total	C	N	O	S	Se	0	0	0
			1423	906	243	265	3	6			
1	B	174	Total	C	N	O	S	Se	0	0	0
			1359	868	232	250	3	6			
1	C	187	Total	C	N	O	S	Se	0	0	0
			1461	931	250	271	3	6			
1	D	194	Total	C	N	O	S	Se	0	0	0
			1530	976	259	286	3	6			

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Cl	0	0
			1	1		
2	A	1	Total	Cl	0	0
			1	1		
2	D	1	Total	Cl	0	0
			1	1		
2	C	1	Total	Cl	0	0
			1	1		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



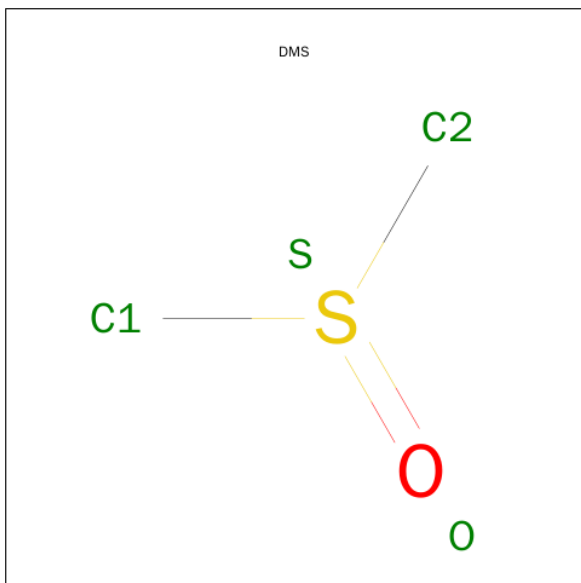
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

- Molecule 5 is DIMETHYL SULFOXIDE (three-letter code: DMS) (formula: C₂H₆OS).



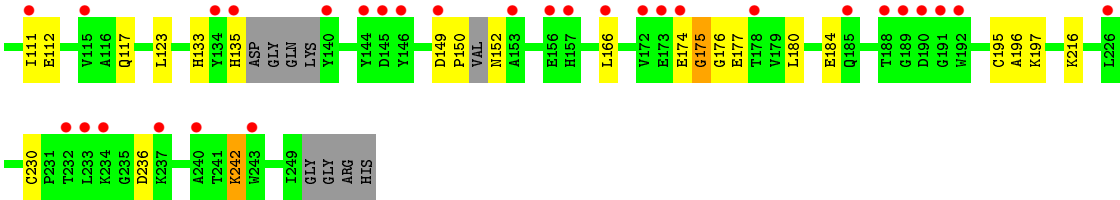
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	B	1	Total	C	O	S	0	0
			4	2	1	1		
5	B	1	Total	C	O	S	0	0
			4	2	1	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	99	Total	O	0	0
			99	99		
6	B	88	Total	O	0	0
			88	88		
6	C	67	Total	O	0	0
			67	67		
6	D	59	Total	O	0	0
			59	59		

- Molecule 1: PROLYL-4 HYDROXYLASE





4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	137.49 Å 137.49 Å 88.13 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.87 – 1.93 19.86 – 1.93	Depositor EDS
% Data completeness (in resolution range)	99.6 (19.87-1.93) 99.6 (19.86-1.93)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.88 (at 1.93 Å)	Xtriage
Refinement program	REFMAC 5.3.0028	Depositor
R, R_{free}	0.188 , 0.223 0.191 , 0.224	Depositor DCC
R_{free} test set	3176 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	27.4	Xtriage
Anisotropy	0.230	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 56.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Outliers	1 of 63507 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6127	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.30% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: GOL, DMS, SO4, CL

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	0.83	0/1451	0.75	2/1951 (0.1%)
1	B	0.94	2/1386 (0.1%)	0.78	1/1864 (0.1%)
1	C	0.80	2/1490 (0.1%)	0.68	1/2004 (0.0%)
1	D	0.87	5/1566 (0.3%)	0.76	3/2110 (0.1%)
All	All	0.86	9/5893 (0.2%)	0.74	7/7929 (0.1%)

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	32	GLY	C-O	19.94	1.55	1.23
1	C	194	GLU	CD-OE2	10.82	1.37	1.25
1	B	32	GLY	CA-C	10.28	1.68	1.51
1	C	194	GLU	CD-OE1	10.26	1.36	1.25
1	D	236	ASP	CG-OD2	9.36	1.46	1.25
1	D	174	GLU	C-N	9.01	1.49	1.33
1	D	72	PRO	C-N	7.93	1.52	1.34
1	D	197	LYS	C-N	5.87	1.47	1.34
1	D	135	HIS	C-O	5.73	1.34	1.23

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	32	GLY	CA-C-O	-11.46	99.98	120.60
1	D	236	ASP	CB-CG-OD1	-10.24	109.08	118.30
1	D	135	HIS	CA-CB-CG	-6.04	103.32	113.60
1	C	34	LEU	CA-CB-CG	5.59	128.15	115.30
1	D	236	ASP	CB-CG-OD2	5.33	123.10	118.30
1	A	185	GLN	CB-CA-C	-5.17	100.05	110.40
1	A	208	ASP	CB-CG-OD1	5.14	122.92	118.30

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	1423	0	1402	7	0
1	B	1359	0	1345	3	0
1	C	1461	0	1446	7	0
1	D	1530	0	1473	27	1
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	6	0	8	0	0
3	B	6	0	8	0	0
3	C	6	0	8	2	0
3	D	6	0	8	0	0
4	A	5	0	0	0	0
5	B	8	0	12	0	0
6	A	99	0	0	2	0
6	B	88	0	0	1	0
6	C	67	0	0	0	0
6	D	59	0	0	4	0
All	All	6127	0	5710	44	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 4.

All (44) 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:D:74:MSE:HE3	1:D:96:THR:HB	1.47	0.96
1:D:39:ARG:HH11	1:D:39:ARG:CG	1.85	0.89
1:C:50:ARG:HH11	3:C:1251:GOL:H11	1.39	0.87
1:D:195:CYS:HG	1:D:230:CYS:HG	0.89	0.83
1:D:39:ARG:HH11	1:D:39:ARG:HG2	1.48	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:156:GLU:HG2	1:A:157:HIS:H	1.60	0.67
1:D:39:ARG:HH11	1:D:39:ARG:HG3	1.61	0.64
1:D:175:GLY:HA2	1:D:177:GLU:OE1	1.99	0.63
1:A:156:GLU:HG2	1:A:157:HIS:N	2.13	0.61
1:C:48:SER:HB3	1:C:184:GLU:HB2	1.89	0.55
1:C:99:TRP:CZ3	1:C:130:GLN:HB2	2.42	0.54
1:D:177:GLU:CG	1:D:196:ALA:HB2	2.38	0.53
1:D:39:ARG:NH1	1:D:39:ARG:CG	2.54	0.53
1:D:149:ASP:HB3	1:D:152:ASN:HD22	1.74	0.53
1:C:38:TRP:CZ2	1:C:40:GLY:HA3	2.43	0.53
1:D:242:LYS:HE3	6:D:2024:HOH:O	2.11	0.51
1:D:216:LYS:NZ	6:D:2048:HOH:O	2.42	0.51
1:D:105:ASP:HB2	1:D:108:ILE:HG12	1.94	0.50
1:D:176:GLY:HA2	1:D:230:CYS:O	2.11	0.50
1:B:48:SER:HB3	1:B:184:GLU:HB3	1.94	0.50
1:C:50:ARG:NH1	3:C:1251:GOL:H11	2.19	0.49
1:D:74:MSE:HE3	1:D:96:THR:CB	2.31	0.48
1:D:39:ARG:NH1	1:D:39:ARG:HG2	2.21	0.47
1:D:62:GLU:HB3	1:D:111:ILE:HD11	1.95	0.47
1:B:55:LYS:NZ	6:B:2011:HOH:O	2.47	0.47
1:A:184:GLU:OE1	6:A:2067:HOH:O	2.20	0.47
1:D:74:MSE:HE1	1:D:133:HIS:HB3	1.97	0.47
1:D:38:TRP:CZ2	1:D:40:GLY:HA3	2.51	0.46
1:A:227:HIS:HD2	6:A:2089:HOH:O	1.99	0.45
1:D:74:MSE:CE	6:D:2028:HOH:O	2.64	0.45
1:C:186:LYS:HG2	1:C:199:GLY:HA2	1.99	0.44
1:A:55:LYS:HE3	1:A:55:LYS:HB3	1.78	0.44
1:A:157:HIS:O	1:A:249:ILE:HG12	2.19	0.43
1:D:49:PRO:HG3	1:D:180:LEU:HD22	2.00	0.43
1:D:74:MSE:HE2	6:D:2028:HOH:O	2.18	0.43
1:D:33:GLU:C	1:D:117:GLN:HE22	2.21	0.43
1:D:149:ASP:O	1:D:152:ASN:HB2	2.20	0.42
1:D:28:MSE:HE3	1:D:123:LEU:HD22	2.01	0.42
1:D:39:ARG:NH1	1:D:39:ARG:HG3	2.27	0.41
1:C:38:TRP:HZ2	1:C:55:LYS:O	2.03	0.41
1:D:48:SER:HB3	1:D:184:GLU:HB2	2.03	0.41
1:A:165:MSE:HB3	1:A:242:LYS:HB3	2.03	0.41
1:B:162:VAL:HB	1:B:246:VAL:HG12	2.03	0.41
1:D:112:GLU:HG2	1:D:242:LYS:CE	2.52	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)
1:D:150:PRO:CB	1:D:150:PRO:CB[8_554]	2.03	0.17

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	174/233 (75%)	171 (98%)	3 (2%)	0	100	100
1	B	166/233 (71%)	166 (100%)	0	0	100	100
1	C	179/233 (77%)	174 (97%)	5 (3%)	0	100	100
1	D	184/233 (79%)	178 (97%)	5 (3%)	1 (0%)	34	20
All	All	703/932 (75%)	689 (98%)	13 (2%)	1 (0%)	56	47

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	175	GLY

5.3.2 Protein sidechains [i](#)

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	154/192 (80%)	152 (99%)	2 (1%)	76	71
1	B	147/192 (77%)	147 (100%)	0	100	100
1	C	158/192 (82%)	158 (100%)	0	100	100
1	D	164/192 (85%)	161 (98%)	3 (2%)	66	58

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	623/768 (81%)	618 (99%)	5 (1%)	86	85

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

Mol	Chain	Res	Type
1	A	94	THR
1	A	185	GLN
1	D	39	ARG
1	D	166	LEU
1	D	242	LYS

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

Mol	Chain	Res	Type
1	A	185	GLN
1	B	185	GLN
1	D	117	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 11 ligands modelled in this entry, 4 are monoatomic - leaving 7 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	1251	-	5,5,5	0.30	0	5,5,5	0.89	0
4	SO4	A	1252	-	4,4,4	0.37	0	6,6,6	0.34	0
3	GOL	B	1251	-	5,5,5	0.38	0	5,5,5	0.27	0
5	DMS	B	1252	-	3,3,3	2.61	1 (33%)	3,3,3	0.72	0
5	DMS	B	1253	-	3,3,3	2.62	1 (33%)	3,3,3	0.51	0
3	GOL	C	1251	-	5,5,5	0.33	0	5,5,5	0.38	0
3	GOL	D	1251	-	5,5,5	0.40	0	5,5,5	0.43	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	GOL	A	1251	-	-	0/4/4/4	0/0/0/0
4	SO4	A	1252	-	-	0/0/0/0	0/0/0/0
3	GOL	B	1251	-	-	0/4/4/4	0/0/0/0
5	DMS	B	1252	-	-	0/0/0/0	0/0/0/0
5	DMS	B	1253	-	-	0/0/0/0	0/0/0/0
3	GOL	C	1251	-	-	0/4/4/4	0/0/0/0
3	GOL	D	1251	-	-	0/4/4/4	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	1252	DMS	O-S	4.35	1.80	1.50
5	B	1253	DMS	O-S	4.39	1.80	1.50

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	1251	GOL	2	0

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	176/233 (75%)	0.65	17 (9%) 10 15	20, 26, 31, 43	0
1	B	168/233 (72%)	0.68	16 (9%) 10 16	21, 26, 31, 35	0
1	C	181/233 (77%)	1.04	32 (17%) 2 3	20, 26, 31, 37	0
1	D	188/233 (80%)	1.11	39 (20%) 1 1	11, 26, 31, 34	0
All	All	713/932 (76%)	0.88	104 (14%) 3 5	11, 26, 31, 43	0

All (104) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	157	HIS	7.2
1	C	94	THR	6.6
1	D	153	ALA	6.0
1	A	36	GLU	6.0
1	D	140	TYR	5.8
1	D	173	GLU	5.5
1	A	108	ILE	5.1
1	C	192	TRP	5.0
1	C	76	LYS	4.8
1	A	190	ASP	4.6
1	A	94	THR	4.5
1	D	233	LEU	4.3
1	D	190	ASP	4.2
1	B	29	VAL	4.2
1	C	188	THR	4.2
1	C	36	GLU	4.2
1	A	156	GLU	4.1
1	C	226	LEU	4.0
1	D	157	HIS	3.9
1	B	135	HIS	3.8
1	D	156	GLU	3.8

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Mol	Chain	Res	Type	RSRZ
1	D	240	ALA	3.7
1	A	226	LEU	3.7
1	B	159	GLY	3.7
1	D	149	ASP	3.7
1	B	190	ASP	3.7
1	D	174	GLU	3.6
1	D	234	LYS	3.5
1	C	157	HIS	3.5
1	C	153	ALA	3.5
1	C	194	GLU	3.5
1	C	234	LYS	3.5
1	C	34	LEU	3.5
1	C	35	LYS	3.5
1	D	189	GLY	3.4
1	C	154	GLY	3.3
1	B	173	GLU	3.3
1	A	107	VAL	3.3
1	D	188	THR	3.2
1	B	115	VAL	3.2
1	D	135	HIS	3.2
1	D	29	VAL	3.1
1	D	191	GLY	3.1
1	B	240	ALA	3.1
1	C	99	TRP	3.1
1	D	144	TYR	3.1
1	A	66	ILE	3.1
1	C	174	GLU	3.0
1	C	178	THR	3.0
1	C	198	ARG	3.0
1	D	134	TYR	3.0
1	B	39	ARG	3.0
1	D	33	GLU	3.0
1	C	135	HIS	2.9
1	C	173	GLU	2.9
1	B	66	ILE	2.9
1	D	115	VAL	2.9
1	C	233	LEU	2.8
1	C	29	VAL	2.8
1	D	99	TRP	2.8
1	B	249	ILE	2.7
1	D	72	PRO	2.7
1	B	129	LEU	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	37	GLU	2.7
1	A	111	ILE	2.6
1	A	169	LEU	2.6
1	D	192	TRP	2.6
1	C	185	GLN	2.6
1	B	169	LEU	2.5
1	A	32	GLY	2.5
1	C	166	LEU	2.5
1	D	146	TYR	2.4
1	B	111	ILE	2.4
1	D	172	VAL	2.4
1	B	99	TRP	2.4
1	A	129	LEU	2.4
1	D	237	LYS	2.4
1	A	166	LEU	2.3
1	C	152	ASN	2.3
1	D	178	THR	2.3
1	C	199	GLY	2.3
1	C	235	GLY	2.3
1	D	166	LEU	2.3
1	C	155	PRO	2.3
1	B	234	LYS	2.2
1	D	226	LEU	2.2
1	B	96	THR	2.2
1	D	111	ILE	2.2
1	D	32	GLY	2.2
1	D	104	GLU	2.2
1	D	185	GLN	2.2
1	D	145	ASP	2.2
1	C	156	GLU	2.1
1	D	61	GLU	2.1
1	A	189	GLY	2.1
1	D	243	TRP	2.1
1	D	232	THR	2.1
1	C	210	LEU	2.1
1	A	240	ALA	2.1
1	C	248	PRO	2.1
1	D	97	GLY	2.1
1	D	107	VAL	2.0
1	A	202	VAL	2.0
1	C	33	GLU	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(\AA^2)	Q<0.9
3	GOL	A	1251	6/6	0.79	0.20	4.88	49,50,51,52	0
3	GOL	B	1251	6/6	0.80	0.20	3.09	55,56,57,57	0
5	DMS	B	1252	4/4	0.84	0.28	2.74	57,58,58,59	0
3	GOL	C	1251	6/6	0.88	0.15	1.32	45,47,48,48	0
2	CL	B	1250	1/1	0.99	0.15	0.15	28,28,28,28	0
4	SO4	A	1252	5/5	0.98	0.14	-0.11	31,32,35,36	0
5	DMS	B	1253	4/4	0.92	0.17	-0.28	68,69,69,69	0
3	GOL	D	1251	6/6	0.88	0.12	-0.48	54,56,57,58	0
2	CL	A	1250	1/1	0.99	0.11	-0.68	28,28,28,28	0
2	CL	C	1250	1/1	0.98	0.13	-1.12	35,35,35,35	0
2	CL	D	1250	1/1	0.98	0.09	-1.45	45,45,45,45	0

6.5 Other polymers [i](#)

There are no such residues in this entry.