



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 10:28 pm BST

PDB ID : 3RBY
Title : Crystal structure of uncharacterized protein YLR301w from *Saccharomyces cerevisiae*
Authors : Kim, K.-H.; Kim, E.E.
Deposited on : 2011-03-30
Resolution : 2.30 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

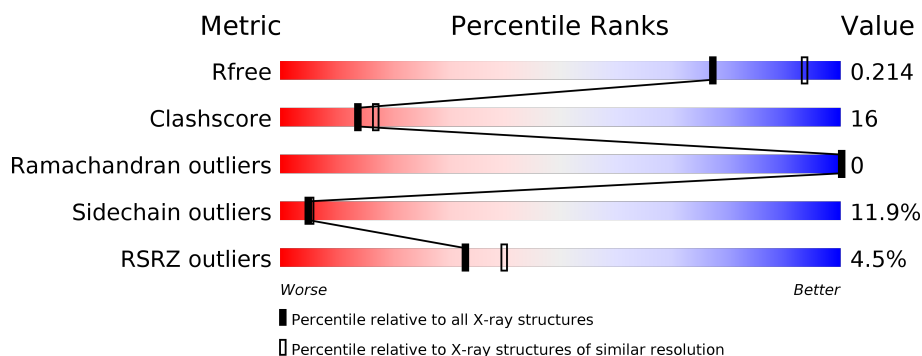
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (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 on 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	245	<div> <div>4%</div> <div>74%</div> <div>20%</div> <div>6%</div> </div>
1	B	245	<div> <div>5%</div> <div>69%</div> <div>27%</div> <div>.</div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4127 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.

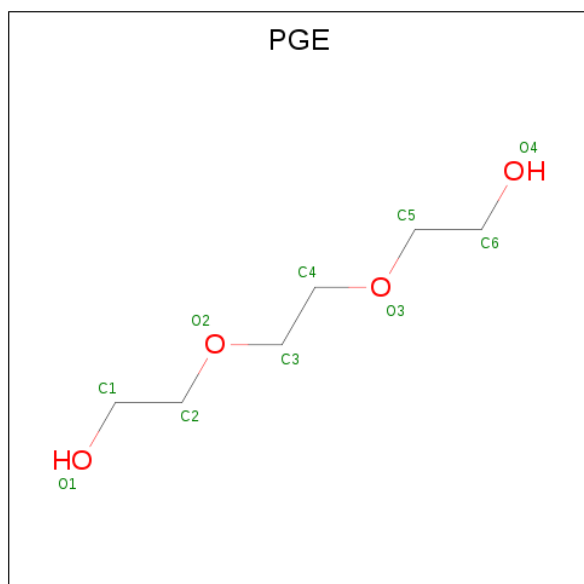
- Molecule 1 is a protein called Uncharacterized protein YLR301W.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	245	Total	C	N	O	S	0	0	0
			1958	1245	331	379	3			
1	B	245	Total	C	N	O	S	0	0	0
			1958	1245	331	379	3			

There are 4 discrepancies between the modelled and reference sequences:

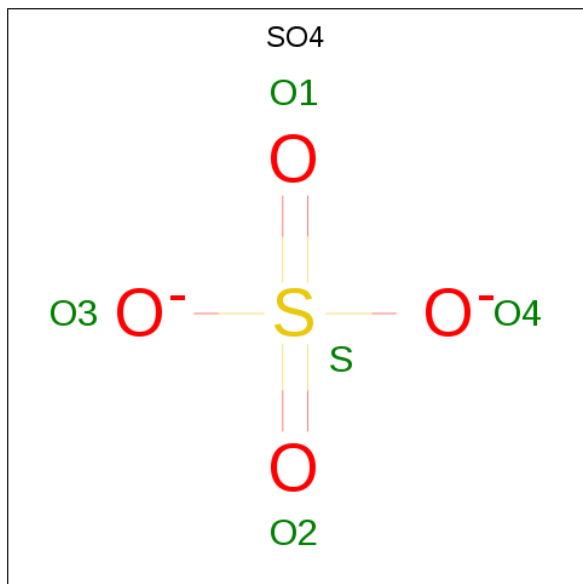
Chain	Residue	Modelled	Actual	Comment	Reference
A	0	HIS	-	EXPRESSION TAG	UNP Q05905
A	128	MET	LEU	CONFLICT	UNP Q05905
B	0	HIS	-	EXPRESSION TAG	UNP Q05905
B	128	MET	LEU	CONFLICT	UNP Q05905

- Molecule 2 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: $C_6H_{14}O_4$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 10 6 4	0	0
2	A	1	Total C O 10 6 4	0	0
2	A	1	Total C O 10 6 4	0	0
2	B	1	Total C O 10 6 4	0	0
2	B	1	Total C O 10 6 4	0	0
2	B	1	Total C O 10 6 4	0	0

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

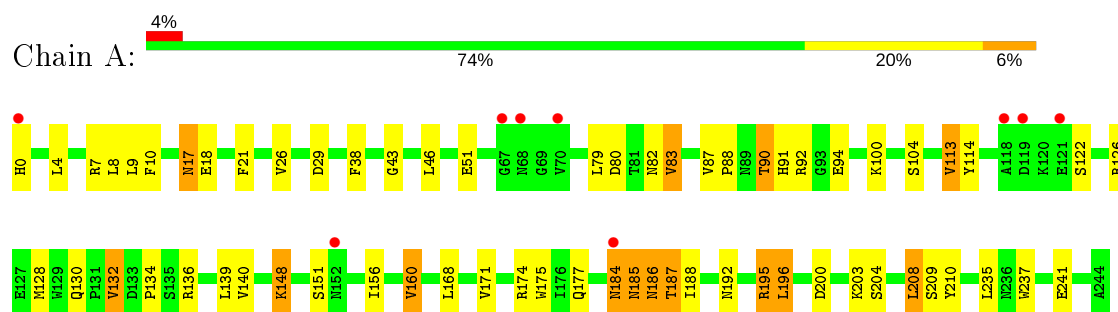
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	87	Total O 87 87	0	0
4	B	59	Total O 59 59	0	0

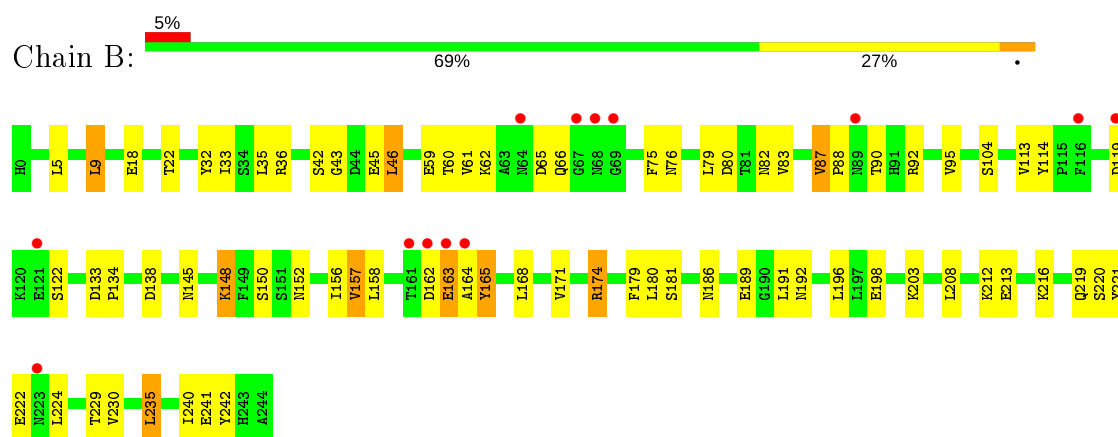
3 Residue-property plots [i](#)

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: Uncharacterized protein YLR301W



- Molecule 1: Uncharacterized protein YLR301W



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	122.32Å 122.32Å 174.00Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	40.24 – 2.30 40.24 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.0 (40.24-2.30) 98.0 (40.24-2.30)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.16 (at 2.29Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7_650)	Depositor
R, R_{free}	0.204 , 0.247 0.207 , 0.214	Depositor DCC
R_{free} test set	1707 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	45.7	Xtriage
Anisotropy	0.170	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 37.3	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	4127	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.13% 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: PGE, SO4

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	0.44	0/2009	0.60	0/2731
1	B	0.46	0/2009	0.66	1/2731 (0.0%)
All	All	0.45	0/4018	0.63	1/5462 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	165	TYR	N-CA-CB	-11.58	89.75	110.60

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	1958	0	1883	69	0
1	B	1958	0	1883	55	0
2	A	30	0	42	7	0
2	B	30	0	42	3	0
3	A	5	0	0	0	0
4	A	87	0	0	5	0
4	B	59	0	0	1	0
All	All	4127	0	3850	124	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 16.

All (124) 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:185:ASN:ND2	1:A:187:THR:H	1.57	1.01
1:A:7:ARG:NH1	1:A:10:PHE:HB2	1.76	1.00
1:A:91:HIS:HD2	1:A:114:TYR:H	1.02	0.98
1:B:90:THR:HG22	1:B:92:ARG:HH12	1.31	0.95
1:B:82:ASN:HB3	1:B:90:THR:HG21	1.47	0.93
1:B:90:THR:HG22	1:B:92:ARG:NH1	1.84	0.93
1:A:185:ASN:HD22	1:A:186:ASN:N	1.67	0.93
1:A:126:ARG:HD2	1:A:128:MET:HE1	1.50	0.93
1:A:82:ASN:HB3	1:A:90:THR:HG21	1.56	0.85
1:B:162:ASP:OD1	1:B:164:ALA:CB	2.26	0.84
1:A:91:HIS:CD2	1:A:114:TYR:H	1.93	0.83
1:A:209:SER:H	2:A:703:PGE:H22	1.41	0.83
1:B:90:THR:CG2	1:B:92:ARG:HH12	1.94	0.80
1:A:114:TYR:CE2	1:A:122:SER:HB3	2.17	0.79
1:B:162:ASP:OD1	1:B:164:ALA:HB3	1.84	0.78
1:A:91:HIS:HD2	1:A:114:TYR:N	1.83	0.75
1:B:164:ALA:HB3	1:B:165:TYR:CD2	2.23	0.73
1:A:185:ASN:HD22	1:A:187:THR:H	1.35	0.73
1:B:43:GLY:HA2	1:B:46:LEU:HD22	1.69	0.72
1:A:186:ASN:ND2	1:A:186:ASN:H	1.88	0.71
1:B:192:ASN:ND2	1:B:213:GLU:H	1.90	0.69
1:A:7:ARG:HH12	1:A:10:PHE:HB2	1.53	0.69
1:A:185:ASN:HD21	1:A:187:THR:HG23	1.57	0.69
1:A:177:GLN:HG2	1:A:195:ARG:HG2	1.74	0.67
1:A:186:ASN:HD22	1:A:186:ASN:H	1.39	0.67
1:B:61:VAL:HG12	1:B:75:PHE:CD2	2.31	0.65
1:A:208:LEU:C	1:A:208:LEU:HD23	2.17	0.65
1:B:208:LEU:HB2	2:B:703:PGE:H32	1.79	0.65
1:B:22:THR:OG1	1:B:36:ARG:HG2	1.97	0.65
1:A:148:LYS:H	1:A:148:LYS:CD	2.10	0.64
1:B:157:VAL:HG13	1:B:241:GLU:HB3	1.77	0.64
1:B:61:VAL:CG1	1:B:75:PHE:CD2	2.83	0.62
1:B:82:ASN:CB	1:B:90:THR:HG21	2.27	0.61
1:A:17:ASN:ND2	1:A:18:GLU:H	1.98	0.61
1:B:95:VAL:HG11	2:B:702:PGE:H22	1.82	0.61
1:A:184:ASN:C	1:A:184:ASN:HD22	2.03	0.60
1:A:185:ASN:HD22	1:A:185:ASN:C	2.03	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:174:ARG:HH21	1:A:200:ASP:HA	1.66	0.60
2:A:702:PGE:H42	4:A:332:HOH:O	2.01	0.59
1:B:192:ASN:HD21	1:B:213:GLU:H	1.49	0.58
1:A:90:THR:HG22	1:A:92:ARG:HH21	1.67	0.58
1:A:186:ASN:ND2	1:A:186:ASN:N	2.51	0.58
1:A:209:SER:N	2:A:703:PGE:H22	2.14	0.58
1:A:185:ASN:ND2	1:A:187:THR:N	2.40	0.58
1:A:90:THR:CG2	1:A:92:ARG:HH21	2.17	0.57
1:B:162:ASP:OD1	1:B:164:ALA:N	2.37	0.57
1:A:185:ASN:ND2	1:A:187:THR:HG23	2.19	0.57
2:A:701:PGE:H12	4:A:264:HOH:O	2.03	0.57
1:B:174:ARG:HG2	1:B:198:GLU:O	2.05	0.56
1:A:87:VAL:O	1:A:90:THR:HB	2.05	0.56
1:A:156:ILE:HB	1:A:171:VAL:HB	1.88	0.56
1:A:91:HIS:CG	1:A:113:VAL:HG22	2.41	0.56
1:B:162:ASP:C	1:B:164:ALA:H	2.08	0.55
1:B:221:TYR:HA	1:B:224:LEU:HD11	1.89	0.55
1:A:91:HIS:CD2	1:A:113:VAL:HG22	2.42	0.55
1:B:114:TYR:CE2	1:B:122:SER:HB3	2.45	0.52
1:B:219:GLN:NE2	4:B:295:HOH:O	2.32	0.51
1:A:185:ASN:HD21	1:A:187:THR:H	1.51	0.51
1:B:32:TYR:CE2	1:B:61:VAL:HG13	2.46	0.51
1:B:61:VAL:HG11	1:B:75:PHE:CE2	2.45	0.51
1:B:213:GLU:HB3	1:B:216:LYS:HD2	1.93	0.51
1:B:22:THR:OG1	1:B:36:ARG:CG	2.59	0.50
1:A:208:LEU:HD23	1:A:209:SER:N	2.26	0.50
1:A:38:PHE:CE2	1:A:51:GLU:HB2	2.46	0.50
1:A:91:HIS:HE1	1:A:94:GLU:O	1.95	0.49
1:A:192:ASN:HA	1:A:210:TYR:O	2.13	0.49
1:A:100:LYS:HD2	4:A:283:HOH:O	2.13	0.48
1:A:136:ARG:HG3	1:A:140:VAL:HG11	1.94	0.48
1:A:26:VAL:HG21	2:A:701:PGE:H22	1.95	0.48
1:B:162:ASP:OD1	1:B:164:ALA:HB2	2.10	0.48
1:A:126:ARG:CD	1:A:128:MET:HE1	2.34	0.48
1:A:148:LYS:H	1:A:148:LYS:HD2	1.79	0.47
1:A:29:ASP:HB2	1:A:241:GLU:OE1	2.14	0.47
1:B:158:LEU:HD22	1:B:230:VAL:HG11	1.96	0.47
1:A:139:LEU:CD1	1:A:175:TRP:CH2	2.98	0.46
1:A:18:GLU:HG3	1:B:104:SER:HB3	1.96	0.46
1:B:224:LEU:HB2	1:B:242:TYR:CE1	2.50	0.46
1:A:80:ASP:HB3	1:A:83:VAL:HG13	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:43:GLY:HA2	1:A:46:LEU:HD12	1.97	0.46
1:A:9:LEU:C	1:A:9:LEU:HD12	2.37	0.46
1:A:21:PHE:CD1	1:B:145:ASN:HB2	2.51	0.46
1:A:136:ARG:NH1	1:B:138:ASP:OD1	2.49	0.45
1:B:163:GLU:H	1:B:163:GLU:HG2	1.42	0.45
1:A:174:ARG:NH1	4:A:313:HOH:O	2.50	0.45
1:A:186:ASN:HD22	1:A:186:ASN:N	2.04	0.45
1:A:8:LEU:HD11	1:A:130:GLN:HB2	1.98	0.44
1:A:160:VAL:HG23	1:A:237:TRP:CZ2	2.53	0.44
2:B:702:PGE:H6	2:B:702:PGE:H42	1.88	0.44
1:B:156:ILE:HB	1:B:171:VAL:HB	1.99	0.44
1:B:162:ASP:C	1:B:164:ALA:N	2.72	0.44
1:B:164:ALA:CB	1:B:165:TYR:CD2	2.98	0.44
1:B:133:ASP:HA	1:B:134:PRO:HD3	1.91	0.43
1:B:42:SER:OG	1:B:45:GLU:HG2	2.18	0.43
1:A:148:LYS:N	1:A:148:LYS:HD2	2.33	0.43
1:B:164:ALA:HB3	1:B:165:TYR:HD2	1.79	0.43
1:B:220:SER:C	1:B:221:TYR:CG	2.92	0.43
1:A:208:LEU:C	1:A:208:LEU:CD2	2.86	0.43
1:A:10:PHE:HZ	2:A:702:PGE:H5	1.84	0.43
1:A:82:ASN:HB3	1:A:90:THR:CG2	2.40	0.43
1:B:157:VAL:HG22	1:B:240:ILE:CG1	2.49	0.43
1:A:139:LEU:CD1	1:A:175:TRP:CZ2	3.02	0.42
1:A:196:LEU:HD22	1:A:204:SER:HB3	2.02	0.42
1:A:160:VAL:HG23	1:A:237:TRP:CE2	2.54	0.42
1:B:9:LEU:HD12	1:B:9:LEU:C	2.40	0.42
2:A:702:PGE:H62	4:A:331:HOH:O	2.20	0.42
1:B:33:ILE:HD11	1:B:35:LEU:HD21	2.02	0.42
1:B:46:LEU:HA	1:B:46:LEU:HD12	1.76	0.42
1:B:60:THR:HB	1:B:76:ASN:O	2.19	0.42
1:B:62:LYS:NZ	1:B:62:LYS:HB3	2.34	0.42
1:A:91:HIS:HB2	1:A:113:VAL:HG22	2.02	0.41
1:B:87:VAL:HA	1:B:88:PRO:HD3	1.82	0.41
1:B:162:ASP:HB3	1:B:235:LEU:HD21	2.03	0.41
1:A:185:ASN:HD21	1:A:187:THR:CG2	2.30	0.41
1:B:165:TYR:N	1:B:165:TYR:CD2	2.88	0.41
1:A:148:LYS:N	1:A:148:LYS:CD	2.81	0.41
1:A:87:VAL:HA	1:A:88:PRO:HD3	1.82	0.41
1:B:80:ASP:OD1	1:B:82:ASN:HB2	2.21	0.41
1:A:7:ARG:NH1	1:A:18:GLU:OE1	2.54	0.41
1:B:148:LYS:H	1:B:148:LYS:HG2	1.58	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:42:SER:H	1:B:45:GLU:CG	2.34	0.41
1:A:160:VAL:HG23	1:A:237:TRP:CH2	2.55	0.41
1:B:164:ALA:HB3	1:B:165:TYR:CE2	2.53	0.40
1:A:132:VAL:O	1:A:134:PRO:HD3	2.21	0.40
1:A:104:SER:HB3	1:B:18:GLU:HG3	2.03	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	243/245 (99%)	233 (96%)	10 (4%)	0	100	100
1	B	243/245 (99%)	235 (97%)	8 (3%)	0	100	100
All	All	486/490 (99%)	468 (96%)	18 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	218/218 (100%)	196 (90%)	22 (10%)	7	9
1	B	218/218 (100%)	188 (86%)	30 (14%)	3	3
All	All	436/436 (100%)	384 (88%)	52 (12%)	5	5

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	0	HIS
1	A	4	LEU
1	A	17	ASN
1	A	79	LEU
1	A	83	VAL
1	A	90	THR
1	A	113	VAL
1	A	132	VAL
1	A	148	LYS
1	A	151	SER
1	A	160	VAL
1	A	168	LEU
1	A	184	ASN
1	A	185	ASN
1	A	186	ASN
1	A	187	THR
1	A	188	ILE
1	A	195	ARG
1	A	196	LEU
1	A	203	LYS
1	A	208	LEU
1	A	235	LEU
1	B	5	LEU
1	B	9	LEU
1	B	46	LEU
1	B	59	GLU
1	B	65	ASP
1	B	66	GLN
1	B	79	LEU
1	B	83	VAL
1	B	87	VAL
1	B	113	VAL
1	B	119	ASP
1	B	148	LYS
1	B	150	SER
1	B	152	ASN
1	B	157	VAL
1	B	163	GLU
1	B	168	LEU
1	B	174	ARG
1	B	179	PHE
1	B	180	LEU

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Mol	Chain	Res	Type
1	B	181	SER
1	B	186	ASN
1	B	189	GLU
1	B	191	LEU
1	B	196	LEU
1	B	203	LYS
1	B	212	LYS
1	B	222	GLU
1	B	229	THR
1	B	235	LEU

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	11	GLN
1	A	17	ASN
1	A	76	ASN
1	A	91	HIS
1	A	96	ASN
1	A	146	ASN
1	A	184	ASN
1	A	185	ASN
1	A	186	ASN
1	B	11	GLN
1	B	96	ASN
1	B	146	ASN
1	B	152	ASN
1	B	192	ASN
1	B	236	ASN

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

7 ligands are 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	$\# Z > 2$
2	PGE	A	701	-	9,9,9	0.81	0	8,8,8	0.65	0
2	PGE	B	701	-	9,9,9	0.86	0	8,8,8	0.69	0
2	PGE	B	702	-	9,9,9	0.97	0	8,8,8	0.51	0
3	SO4	A	600	-	4,4,4	0.18	0	6,6,6	0.27	0
2	PGE	A	703	-	9,9,9	1.00	0	8,8,8	0.49	0
2	PGE	B	703	-	9,9,9	0.79	0	8,8,8	0.67	0
2	PGE	A	702	-	9,9,9	0.65	0	8,8,8	0.84	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
2	PGE	A	701	-	-	4/7/7/7	-
2	PGE	B	701	-	-	1/7/7/7	-
2	PGE	B	702	-	-	4/7/7/7	-
2	PGE	A	703	-	-	2/7/7/7	-
2	PGE	B	703	-	-	5/7/7/7	-
2	PGE	A	702	-	-	4/7/7/7	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	703	PGE	O2-C3-C4-O3
2	A	703	PGE	O3-C5-C6-O4
2	B	702	PGE	C6-C5-O3-C4
2	B	701	PGE	O3-C5-C6-O4
2	A	702	PGE	O1-C1-C2-O2
2	A	701	PGE	O1-C1-C2-O2
2	B	703	PGE	O3-C5-C6-O4
2	B	702	PGE	O3-C5-C6-O4
2	A	701	PGE	C3-C4-O3-C5
2	B	703	PGE	C6-C5-O3-C4
2	A	701	PGE	C1-C2-O2-C3
2	B	703	PGE	C4-C3-O2-C2
2	A	703	PGE	C1-C2-O2-C3
2	A	702	PGE	C3-C4-O3-C5
2	B	703	PGE	C1-C2-O2-C3
2	A	702	PGE	C4-C3-O2-C2
2	A	701	PGE	O2-C3-C4-O3
2	B	702	PGE	C3-C4-O3-C5
2	B	702	PGE	O1-C1-C2-O2
2	A	702	PGE	O2-C3-C4-O3

There are no ring outliers.

5 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	701	PGE	2	0
2	B	702	PGE	2	0
2	A	703	PGE	2	0
2	B	703	PGE	1	0
2	A	702	PGE	3	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	245/245 (100%)	-0.11	9 (3%) 41 48	33, 44, 68, 95	0
1	B	245/245 (100%)	-0.14	13 (5%) 26 33	33, 48, 73, 90	0
All	All	490/490 (100%)	-0.12	22 (4%) 33 40	33, 46, 72, 95	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	119	ASP	5.2
1	B	161	THR	5.0
1	A	118	ALA	4.1
1	B	67	GLY	3.2
1	A	67	GLY	3.2
1	B	69	GLY	3.0
1	B	116	PHE	3.0
1	A	0	HIS	2.7
1	A	152	ASN	2.6
1	A	184	ASN	2.6
1	B	163	GLU	2.5
1	B	164	ALA	2.4
1	B	119	ASP	2.4
1	A	70	VAL	2.4
1	B	64	ASN	2.3
1	B	89	ASN	2.2
1	A	121	GLU	2.2
1	B	68	ASN	2.2
1	B	121	GLU	2.1
1	A	68	ASN	2.1
1	B	223	ASN	2.0
1	B	162	ASP	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. 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(\AA^2)	Q<0.9
2	PGE	B	702	10/10	0.84	0.14	51,55,64,65	0
2	PGE	A	703	10/10	0.88	0.16	45,46,55,56	0
2	PGE	B	703	10/10	0.88	0.17	44,48,52,54	0
2	PGE	A	701	10/10	0.91	0.14	42,47,63,65	0
2	PGE	A	702	10/10	0.93	0.10	44,52,60,66	0
2	PGE	B	701	10/10	0.94	0.12	47,53,71,79	0
3	SO4	A	600	5/5	0.97	0.16	49,50,62,62	0

6.5 Other polymers [i](#)

There are no such residues in this entry.