Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:04 AM GMT

PDB ID	:	2ZFE
Title	:	Crystal structure of bacteriorhodopsin-xenon complex
Authors	:	Kouyama, T.
Deposited on	:	2007-12-31
Resolution	:	2.50 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:

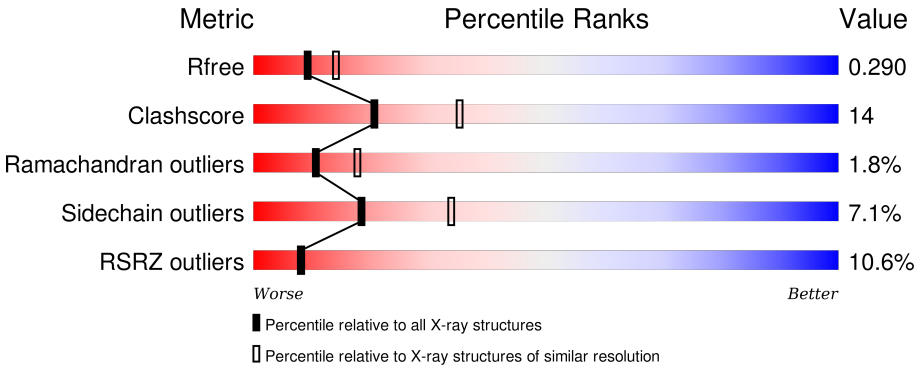
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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 <=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	262	9%, 60%, 23%, 13%

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
2	GLC	A	400	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	L3P	A	260	-	-	-	X
5	L2P	A	270	-	-	-	X
6	L1P	A	280	-	-	-	X
6	L1P	A	290	-	-	-	X
6	L1P	A	300	-	-	-	X
7	XE	A	638	-	-	X	-

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 2099 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.

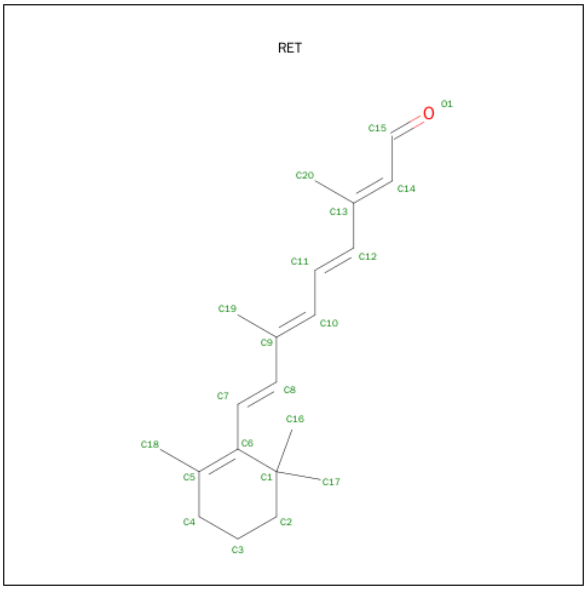
- Molecule 1 is a protein called Bacteriorhodopsin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	227	Total	C	N	O	S	0	0	0
			1756	1180	268	299	9			

- Molecule 2 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	3	Total	C	O	S	0	0
			37	18	18	1		

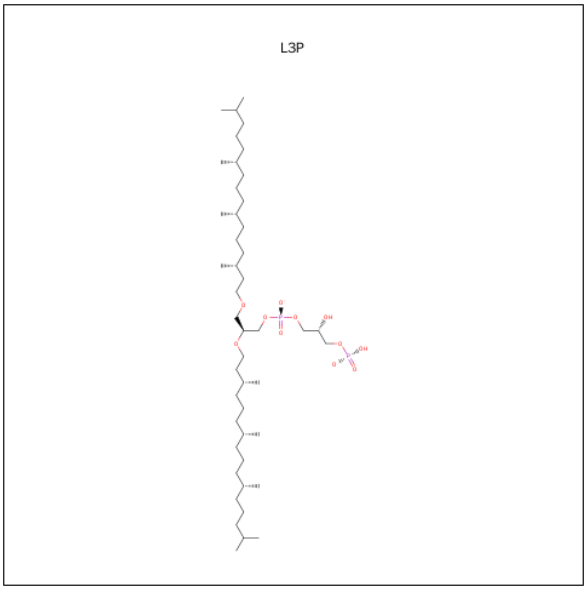
- Molecule 3 is RETINAL (three-letter code: RET) (formula: $C_{20}H_{28}O$).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	C	0	0
			20	20		

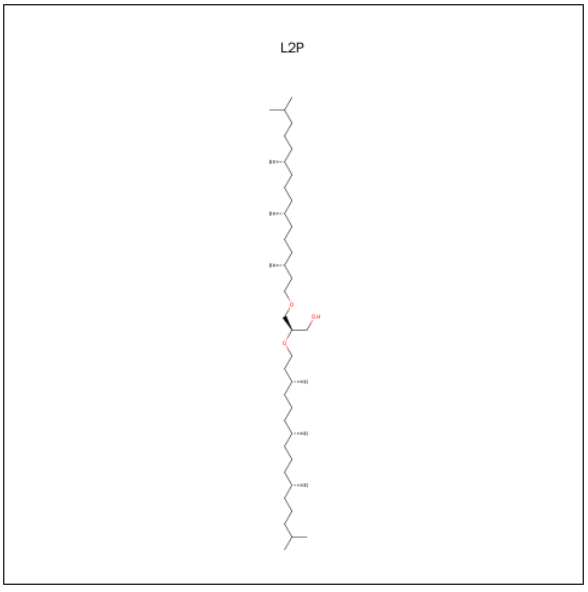
- Molecule 4 is 2,3-DI-O-PHYTANLY-3-SN-GLYCERO-1-PHOSPHORYL-3'-SN-GLYCER

OL-1'-PHOSPHATE (three-letter code: L3P) (formula: $C_{46}H_{94}O_{11}P_2$).



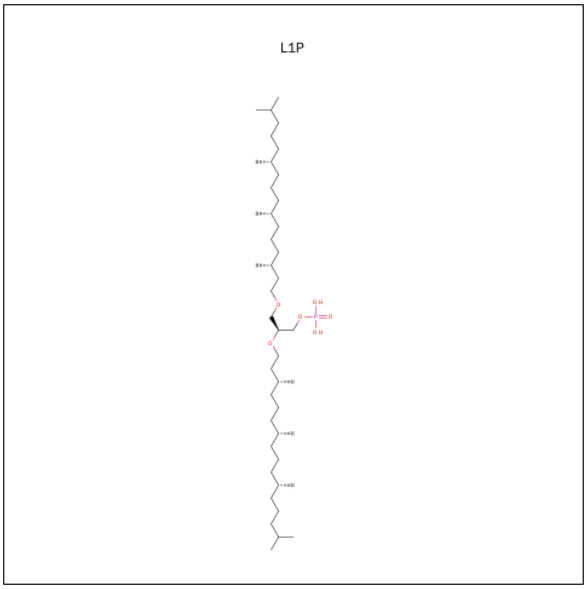
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	O	P		
4	A	1	59	46	11	2	0	0

- Molecule 5 is 2,3-DI-PHYTANYL-GLYCEROL (three-letter code: L2P) (formula: $C_{43}H_{88}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
5	A	1	46	43	3	0	0

- Molecule 6 is 3-PHOSPHORYL-[1,2-DI-PHYTANYL]GLYCEROL (three-letter code: L1P) (formula: $C_{43}H_{89}O_6P$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	O	P	0	0
			40	33	6	1		
6	A	1	Total	C	O	P	0	0
			50	43	6	1		
6	A	1	Total	C	O	P	0	0
			50	43	6	1		

- Molecule 7 is XENON (three-letter code: XE) (formula: Xe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	1	Total	Xe	0	0
			1	1		

- Molecule 8 is water.

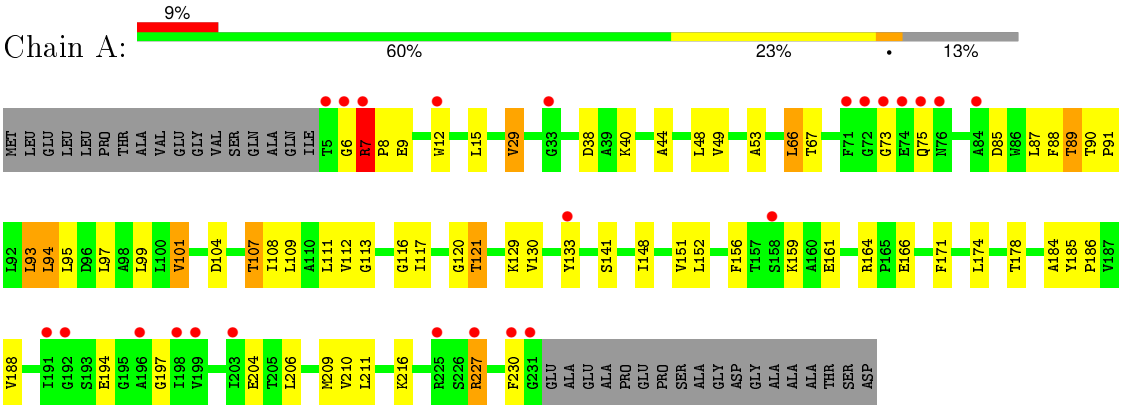
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	40	Total	O	0	0
			40	40		

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 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: Bacteriorhodopsin

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 6 2 2	Depositor
Cell constants a, b, c, α, β, γ	102.41Å 102.41Å 112.47Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	15.00 – 2.50 37.86 – 2.50	Depositor EDS
% Data completeness (in resolution range)	98.8 (15.00-2.50) 98.7 (37.86-2.50)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$< I/\sigma(I) >$ [1]	4.75 (at 2.51Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.250 , 0.279 0.258 , 0.290	Depositor DCC
R_{free} test set	627 reflections (5.07%)	DCC
Wilson B-factor (Å^2)	45.0	Xtriage
Anisotropy	0.554	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 58.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.35$	Xtriage
Outliers	0 of 12439 reflections	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	2099	wwPDB-VP
Average B, all atoms (Å^2)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.04% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: GLC, L1P, RET, XE, L2P, L3P, SGA, MAN

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	0.41	0/1804	0.56	0/2464

There are no bond length outliers.

There are no bond angle outliers.

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	1756	0	1813	62	0
2	A	37	0	28	1	0
3	A	20	0	27	4	0
4	A	59	0	93	5	0
5	A	46	0	87	7	0
6	A	140	0	238	0	0
7	A	1	0	0	3	0
8	A	40	0	0	2	0
All	All	2099	0	2286	63	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 (63) 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:87:LEU:HD11	5:A:270:L2P:H271	1.32	1.11
1:A:121:THR:HG22	1:A:141:SER:HB2	1.43	0.99
1:A:108:ILE:HG23	7:A:638:XE:XE	2.48	0.91
1:A:87:LEU:HD21	5:A:270:L2P:H262	1.55	0.88
1:A:29:VAL:O	1:A:29:VAL:HG12	1.80	0.81
1:A:174:LEU:O	1:A:178:THR:HG23	1.84	0.77
1:A:121:THR:CG2	1:A:141:SER:HB2	2.15	0.74
1:A:111:LEU:HD21	1:A:151:VAL:HG11	1.69	0.74
1:A:15:LEU:HB3	1:A:209:MET:HE2	1.70	0.72
1:A:117:ILE:O	1:A:121:THR:HB	1.90	0.71
1:A:44:ALA:HA	4:A:260:L3P:H111	1.76	0.67
1:A:6:GLY:HA2	1:A:9:GLU:OE1	1.95	0.67
1:A:94:LEU:HB3	7:A:638:XE:XE	2.74	0.65
1:A:107:THR:O	1:A:111:LEU:HD23	1.97	0.65
1:A:40:LYS:HZ1	4:A:260:L3P:H42	1.63	0.63
1:A:227:ARG:HH11	1:A:227:ARG:HB3	1.62	0.63
1:A:104:ASP:HB2	1:A:107:THR:HG22	1.81	0.62
1:A:101:VAL:HG22	1:A:159:LYS:HB3	1.82	0.61
1:A:116:GLY:HA3	5:A:270:L2P:H541	1.82	0.61
1:A:104:ASP:HB2	1:A:107:THR:CG2	2.32	0.60
1:A:49:VAL:HG21	1:A:93:LEU:HD13	1.83	0.60
1:A:164:ARG:NH2	1:A:230:PHE:O	2.35	0.60
1:A:85:ASP:O	1:A:89:THR:HB	2.03	0.58
1:A:90:THR:OG1	1:A:91:PRO:HD3	2.06	0.55
1:A:15:LEU:HD23	1:A:209:MET:CE	2.36	0.55
1:A:164:ARG:NH1	1:A:166:GLU:OE2	2.40	0.54
1:A:152:LEU:O	1:A:156:PHE:HB2	2.09	0.53
1:A:186:PRO:HB3	3:A:250:RET:H183	1.91	0.52
1:A:12:TRP:CE3	1:A:206:LEU:HD23	2.44	0.52
1:A:141:SER:HB3	3:A:250:RET:H41	1.91	0.51
1:A:53:ALA:HB2	1:A:216:LYS:HE2	1.93	0.51
1:A:129:LYS:HE2	2:A:410:MAN:H61	1.94	0.50
1:A:40:LYS:HZ2	4:A:260:L3P:P1	2.36	0.49
1:A:66:LEU:HD23	1:A:67:THR:N	2.28	0.49
1:A:87:LEU:CD1	5:A:270:L2P:H271	2.24	0.48
1:A:194:GLU:OE1	1:A:204:GLU:OE2	2.30	0.48
1:A:29:VAL:O	1:A:29:VAL:CG1	2.53	0.48
1:A:112:VAL:HG23	7:A:638:XE:XE	2.92	0.47
1:A:109:LEU:HD11	5:A:270:L2P:H592	1.97	0.47
1:A:166:GLU:HG3	8:A:630:HOH:O	2.14	0.46
1:A:120:GLY:HA3	5:A:270:L2P:H491	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:48:LEU:HG	4:A:260:L3P:H143	1.98	0.46
1:A:130:VAL:HB	1:A:133:TYR:CD2	2.52	0.45
1:A:15:LEU:HD23	1:A:209:MET:HE1	1.98	0.45
1:A:185:TYR:N	1:A:186:PRO:HD2	2.32	0.45
1:A:7:ARG:H	1:A:8:PRO:HD2	1.83	0.44
1:A:90:THR:N	1:A:91:PRO:CD	2.80	0.44
1:A:113:GLY:HA2	5:A:270:L2P:H543	2.00	0.44
3:A:250:RET:H7	3:A:250:RET:H181	1.84	0.43
1:A:107:THR:HG21	8:A:629:HOH:O	2.18	0.43
1:A:75:GLN:OE1	1:A:75:GLN:HA	2.18	0.43
1:A:88:PHE:C	1:A:91:PRO:HD2	2.40	0.42
1:A:95:LEU:O	1:A:99:LEU:HG	2.18	0.42
1:A:12:TRP:CD2	1:A:206:LEU:HD23	2.55	0.42
1:A:206:LEU:O	1:A:210:VAL:HG23	2.19	0.42
1:A:156:PHE:HB3	1:A:171:PHE:CZ	2.55	0.42
1:A:15:LEU:HD23	1:A:209:MET:HE2	2.01	0.41
1:A:184:ALA:O	1:A:188:VAL:HG23	2.21	0.41
1:A:7:ARG:N	1:A:8:PRO:HD2	2.36	0.41
1:A:141:SER:CB	3:A:250:RET:H41	2.51	0.41
1:A:111:LEU:HD12	1:A:148:ILE:HG23	2.03	0.41
1:A:53:ALA:HA	1:A:85:ASP:OD2	2.21	0.41
1:A:44:ALA:HA	4:A:260:L3P:H142	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	225/262 (86%)	209 (93%)	12 (5%)	4 (2%)	11 18

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	7	ARG
1	A	29	VAL
1	A	197	GLY
1	A	73	GLY

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	182/206 (88%)	169 (93%)	13 (7%)	18 34

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	ARG
1	A	38	ASP
1	A	66	LEU
1	A	89	THR
1	A	93	LEU
1	A	94	LEU
1	A	97	LEU
1	A	101	VAL
1	A	107	THR
1	A	121	THR
1	A	161	GLU
1	A	211	LEU
1	A	227	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	105	GLN

5.3.3 RNA

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)

3 carbohydrates 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	GLC	A	400	2,5	11,11,12	3.06	5 (45%)	14,15,17	1.24	2 (14%)
2	MAN	A	410	2	11,11,12	2.34	5 (45%)	14,15,17	3.82	7 (50%)
2	SGA	A	420	2	15,15,16	1.93	5 (33%)	18,22,24	3.93	9 (50%)

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	GLC	A	400	2,5	-	0/2/19/22	0/1/1/1
2	MAN	A	410	2	-	0/2/19/22	0/1/1/1
2	SGA	A	420	2	-	1/7/24/27	0/1/1/1

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	420	SGA	O3-C3	-2.41	1.41	1.46
2	A	400	GLC	C4-C5	2.02	1.57	1.53
2	A	410	MAN	O5-C5	2.10	1.48	1.43
2	A	420	SGA	O5-C5	2.45	1.48	1.43
2	A	410	MAN	O5-C1	2.54	1.48	1.43
2	A	420	SGA	C2-C3	2.56	1.58	1.52
2	A	410	MAN	C1-C2	2.67	1.58	1.52
2	A	420	SGA	O2S-S	3.55	1.58	1.45
2	A	420	SGA	O1S-S	3.59	1.58	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	410	MAN	O2-C2	3.60	1.51	1.43
2	A	400	GLC	C2-C3	4.59	1.58	1.52
2	A	400	GLC	O5-C5	4.73	1.53	1.43
2	A	400	GLC	C1-C2	4.74	1.63	1.52
2	A	410	MAN	C2-C3	4.88	1.59	1.52
2	A	400	GLC	O5-C1	5.35	1.52	1.43

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	410	MAN	O5-C1-C2	-5.09	102.60	110.86
2	A	420	SGA	O2S-S-O1S	-4.73	91.84	112.46
2	A	420	SGA	C1-C2-C3	-4.70	103.47	109.24
2	A	420	SGA	O3S-S-O1S	-4.53	91.54	108.56
2	A	410	MAN	C2-C3-C4	-3.80	104.59	111.04
2	A	410	MAN	C6-C5-C4	-3.23	105.04	113.02
2	A	420	SGA	O3S-S-O2S	-2.21	100.26	108.56
2	A	400	GLC	O5-C1-C2	-2.15	107.37	110.86
2	A	400	GLC	C2-C3-C4	-2.08	107.51	111.04
2	A	420	SGA	C6-C5-C4	-2.06	107.94	113.02
2	A	410	MAN	O2-C2-C3	2.12	114.38	110.12
2	A	420	SGA	O5-C5-C6	2.34	112.42	107.35
2	A	410	MAN	C1-C2-C3	2.36	112.33	109.54
2	A	420	SGA	O3-S-O1S	5.41	124.41	106.86
2	A	410	MAN	C1-O5-C5	6.11	120.00	112.25
2	A	420	SGA	O3-S-O2S	6.40	127.63	106.86
2	A	410	MAN	O6-C6-C5	10.09	144.67	111.33
2	A	420	SGA	C3-O3-S	11.07	139.86	118.77

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	420	SGA	S-O3-C3-C2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	410	MAN	1	0

5.6 Ligand geometry (i)

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	$\#\|Z\| > 2$
3	RET	A	250	1	19,20,21	2.06	4 (21%)	27,27,28	1.73	11 (40%)
4	L3P	A	260	-	58,58,58	1.18	8 (13%)	64,73,73	1.23	6 (9%)
5	L2P	A	270	2	45,45,45	1.69	11 (24%)	50,53,53	1.19	5 (10%)
6	L1P	A	280	-	39,39,49	1.95	10 (25%)	43,48,60	1.02	2 (4%)
6	L1P	A	290	-	49,49,49	1.73	14 (28%)	55,60,60	1.44	7 (12%)
6	L1P	A	300	-	49,49,49	1.62	11 (22%)	55,60,60	1.18	6 (10%)

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	RET	A	250	1	-	0/13/30/31	0/1/1/1
4	L3P	A	260	-	-	0/67/67/67	0/0/0/0
5	L2P	A	270	2	-	0/51/51/51	0/0/0/0
6	L1P	A	280	-	-	0/43/43/55	0/0/0/0
6	L1P	A	290	-	-	0/55/55/55	0/0/0/0
6	L1P	A	300	-	-	0/55/55/55	0/0/0/0

All (58) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	250	RET	C2-C3	-2.44	1.45	1.52
4	A	260	L3P	P1-O2P	-2.20	1.45	1.54
4	A	260	L3P	P1-O4	2.01	1.68	1.59
6	A	290	L1P	C54-C53	2.02	1.59	1.52
4	A	260	L3P	C6-C5	2.14	1.59	1.51
6	A	300	L1P	C14-C13	2.14	1.60	1.52
6	A	300	L1P	C19-C18	2.15	1.60	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	290	L1P	C25-C23	2.16	1.63	1.52
6	A	290	L1P	C56-C55	2.18	1.62	1.52
6	A	280	L1P	C42-C41	2.19	1.58	1.50
5	A	270	L2P	C1-C2	2.21	1.57	1.50
6	A	280	L1P	C52-C53	2.24	1.64	1.52
6	A	300	L1P	C26-C27	2.25	1.62	1.52
6	A	290	L1P	C26-C27	2.27	1.62	1.52
6	A	290	L1P	C19-C18	2.29	1.60	1.52
5	A	270	L2P	C22-C23	2.30	1.64	1.52
6	A	290	L1P	C14-C13	2.31	1.60	1.52
5	A	270	L2P	C25-C23	2.31	1.64	1.52
6	A	280	L1P	P-O3	2.34	1.68	1.60
6	A	300	L1P	O1-C11	2.36	1.52	1.42
5	A	270	L2P	C50-C48	2.36	1.65	1.52
6	A	300	L1P	C55-C53	2.36	1.65	1.52
6	A	300	L1P	C17-C18	2.36	1.65	1.52
6	A	280	L1P	O1-C1	2.38	1.49	1.42
6	A	290	L1P	C17-C18	2.38	1.65	1.52
4	A	260	L3P	C42-C41	2.39	1.59	1.50
6	A	280	L1P	C56-C55	2.41	1.63	1.52
5	A	270	L2P	O2-C2	2.45	1.50	1.43
6	A	290	L1P	C55-C53	2.49	1.65	1.52
4	A	260	L3P	P2-O6	2.51	1.67	1.60
3	A	250	RET	C7-C6	2.53	1.55	1.45
6	A	300	L1P	C15-C13	2.57	1.66	1.52
6	A	290	L1P	O1-C11	2.60	1.53	1.42
5	A	270	L2P	C24-C23	2.62	1.61	1.52
4	A	260	L3P	P1-O3	2.67	1.71	1.59
6	A	290	L1P	C15-C13	2.71	1.66	1.52
4	A	260	L3P	O2-C2	2.78	1.51	1.43
6	A	280	L1P	O2-C41	2.81	1.50	1.42
6	A	300	L1P	C12-C11	2.88	1.61	1.50
6	A	290	L1P	C12-C11	2.93	1.61	1.50
5	A	270	L2P	C42-C41	3.03	1.61	1.50
5	A	270	L2P	O2-C41	3.08	1.51	1.42
6	A	300	L1P	P-O3P	3.10	1.65	1.54
4	A	260	L3P	O2-C41	3.13	1.51	1.42
5	A	270	L2P	C47-C48	3.13	1.69	1.52
6	A	290	L1P	P-O3P	3.17	1.66	1.54
6	A	300	L1P	C3-C2	3.17	1.59	1.50
6	A	290	L1P	C3-C2	3.43	1.60	1.50
5	A	270	L2P	C3-C2	3.46	1.59	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	270	L2P	C26-C25	3.70	1.69	1.52
6	A	280	L1P	C1-C2	3.76	1.61	1.50
6	A	300	L1P	P-O3	4.11	1.74	1.60
6	A	280	L1P	C3-C2	4.11	1.62	1.50
6	A	290	L1P	P-O3	4.23	1.74	1.60
6	A	280	L1P	P-O3P	4.40	1.70	1.54
3	A	250	RET	C5-C6	4.98	1.42	1.34
6	A	280	L1P	O2-C2	5.24	1.57	1.43
3	A	250	RET	C1-C6	5.97	1.62	1.53

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	A	270	L2P	C24-C23-C22	-3.65	97.05	111.08
5	A	270	L2P	C22-C21-C20	-3.30	96.62	112.99
4	A	260	L3P	C46-C45-C43	-2.80	106.22	115.49
3	A	250	RET	C8-C9-C10	-2.63	114.74	118.98
5	A	270	L2P	C26-C25-C23	-2.58	106.92	115.49
4	A	260	L3P	C26-C25-C23	-2.53	107.09	115.49
3	A	250	RET	C1-C6-C5	-2.47	119.03	122.66
6	A	300	L1P	C47-C46-C45	-2.42	100.97	112.99
6	A	290	L1P	C47-C46-C45	-2.39	101.11	112.99
4	A	260	L3P	C51-C50-C48	-2.32	107.81	115.49
3	A	250	RET	C7-C6-C5	-2.24	116.24	121.37
3	A	250	RET	C16-C1-C2	-2.08	101.32	108.79
3	A	250	RET	C3-C4-C5	2.03	117.08	113.87
6	A	290	L1P	O1-C11-C12	2.04	115.45	109.52
3	A	250	RET	C17-C1-C6	2.06	113.53	110.30
3	A	250	RET	C19-C9-C8	2.10	121.59	118.10
4	A	260	L3P	O6P-P2-O6	2.12	110.53	105.35
6	A	290	L1P	C16-C17-C18	2.16	122.64	115.49
6	A	290	L1P	C21-C20-C18	2.24	122.92	115.49
6	A	300	L1P	O3-P-O1P	2.26	112.89	107.14
3	A	250	RET	C16-C1-C6	2.26	113.85	110.30
6	A	300	L1P	C21-C20-C18	2.31	123.15	115.49
6	A	300	L1P	C16-C17-C18	2.33	123.21	115.49
6	A	290	L1P	O3-P-O1P	2.34	113.11	107.14
3	A	250	RET	C2-C3-C4	2.36	117.47	111.53
6	A	280	L1P	O1-C1-C2	2.46	112.86	108.71
5	A	270	L2P	C21-C20-C18	2.48	123.70	115.49
3	A	250	RET	C2-C1-C6	2.60	114.48	110.36
6	A	300	L1P	O3P-P-O3	2.66	114.23	106.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	A	290	L1P	O3P-P-O3	2.73	114.42	106.56
5	A	270	L2P	C21-C22-C23	3.08	125.69	115.49
3	A	250	RET	C1-C6-C7	3.18	124.72	115.82
4	A	260	L3P	C11-O1-C1	3.37	124.42	113.44
6	A	280	L1P	C51-C50-C48	3.65	127.59	115.49
6	A	300	L1P	O1-C1-C2	4.26	115.89	108.71
4	A	260	L3P	C41-O2-C2	5.03	126.96	115.50
6	A	290	L1P	O1-C1-C2	7.64	121.60	108.71

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	250	RET	4	0
4	A	260	L3P	5	0
5	A	270	L2P	7	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, 95^{th} 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	227/262 (86%)	0.42	24 (10%) 8 8	33, 44, 66, 88	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	5	THR	12.3
1	A	72	GLY	5.9
1	A	74	GLU	4.5
1	A	196	ALA	4.2
1	A	6	GLY	3.7
1	A	73	GLY	3.7
1	A	203	ILE	3.3
1	A	198	ILE	2.7
1	A	71	PHE	2.7
1	A	12	TRP	2.7
1	A	227	ARG	2.6
1	A	192	GLY	2.4
1	A	158	SER	2.4
1	A	7	ARG	2.4
1	A	231	GLY	2.3
1	A	84	ALA	2.3
1	A	33	GLY	2.3
1	A	191	ILE	2.2
1	A	76	ASN	2.2
1	A	199	VAL	2.2
1	A	225	ARG	2.1
1	A	230	PHE	2.1
1	A	133	TYR	2.0
1	A	75	GLN	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)

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, 95^{th} 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
2	GLC	A	400	11/12	0.81	0.29	2.01	89,90,91,92	0
2	MAN	A	410	11/12	0.80	0.26	1.30	95,97,101,104	0
2	SGA	A	420	15/16	0.09	0.79	-	106,107,113,113	0

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, 95^{th} 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
5	L2P	A	270	46/46	0.69	0.65	11.55	67,86,114,114	0
4	L3P	A	260	59/59	0.59	0.49	7.31	67,92,104,105	13
6	L1P	A	300	50/50	0.43	0.47	5.56	112,122,126,126	0
6	L1P	A	280	40/50	0.37	0.49	5.02	108,124,128,129	0
6	L1P	A	290	50/50	0.27	0.53	3.73	120,130,140,141	0
3	RET	A	250	20/21	0.92	0.19	1.51	34,40,43,44	0
7	XE	A	638	1/1	0.97	0.09	-2.80	44,44,44,44	1

6.5 Other polymers (i)

There are no such residues in this entry.