



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 09:45 AM GMT

PDB ID : 3DDL
Title : Crystallographic Structure of Xanthorhodopsin, a Light-Driven Ion Pump with Dual Chromophore
Authors : Stagno, J.; Luecke, H.; Schobert, B.; Lanyi, J.K.; Imasheva, E.S.; Wang, J.M.; Balashov, S.P.
Deposited on : 2008-06-05
Resolution : 1.90 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>

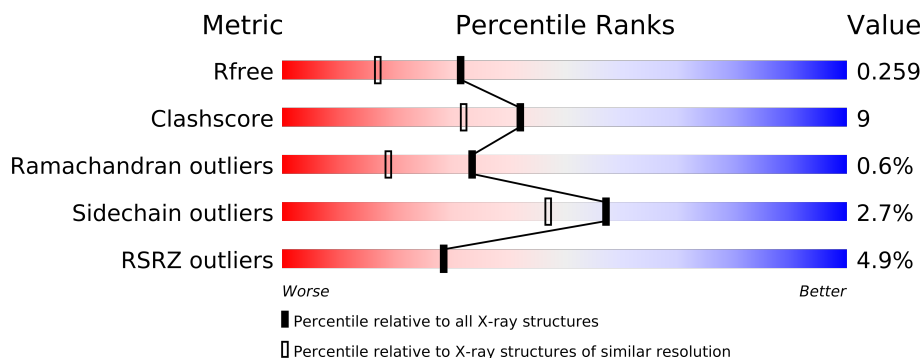
The following versions of software and data (see [references](#)) were used in the production of this report:

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 1.90 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	3684 (1.90-1.90)
Clashscore	79885	4465 (1.90-1.90)
Ramachandran outliers	78287	4413 (1.90-1.90)
Sidechain outliers	78261	4414 (1.90-1.90)
RSRZ outliers	66119	3686 (1.90-1.90)

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	273	
1	B	273	

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

Mol	Type	Chain	Res	Geometry	Electron density
2	SXN	A	1401	-	X
2	SXN	B	1401	-	X
3	UNL	A	1402	-	X
3	UNL	A	1403	-	X
3	UNL	A	1404	-	X
3	UNL	A	1405	-	X

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Mol	Type	Chain	Res	Geometry	Electron density
3	UNL	A	1406	-	X
3	UNL	A	1407	-	X
3	UNL	A	1408	-	X
3	UNL	A	1409	-	X
3	UNL	B	1402	-	X
3	UNL	B	1403	-	X
3	UNL	B	1404	-	X
3	UNL	B	1405	-	X
3	UNL	B	1406	-	X
3	UNL	B	1409	-	X
3	UNL	B	1411	-	X
3	UNL	B	1412	-	X
3	UNL	B	1413	-	X
3	UNL	B	1414	-	X
6	PCW	B	1416	-	X

2 Entry composition i

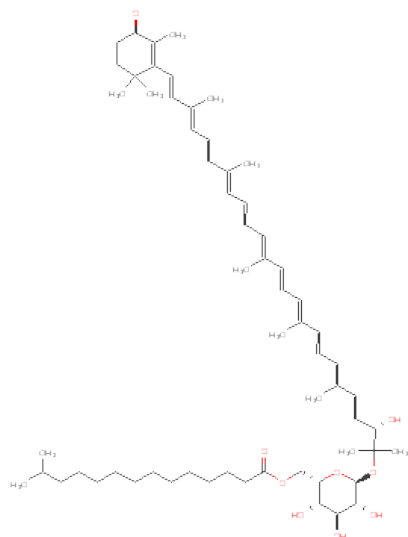
There are 7 unique types of molecules in this entry. The entry contains 4417 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Xanthorhodopsin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	252	Total	C	N	O	S	0	0	0
			1971	1307	310	345	9			
1	B	250	Total	C	N	O	S	0	0	0
			1954	1297	308	340	9			

- Molecule 2 is SALINIXANTHIN (three-letter code: SXN) (formula: $C_{61}H_{92}O_9$).

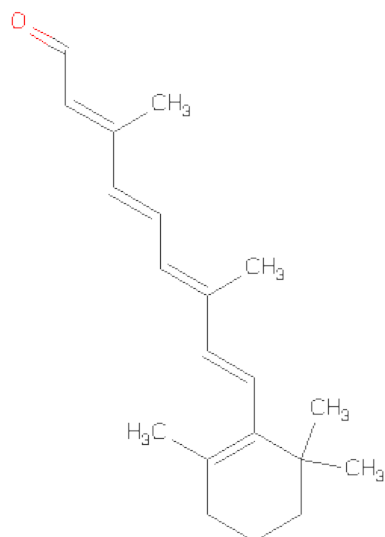


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			70	61	9		
2	B	1	Total	C	O	0	0
			70	61	9		

- Molecule 3 is UNKNOWN LIGAND (three-letter code: UNL) (formula:).

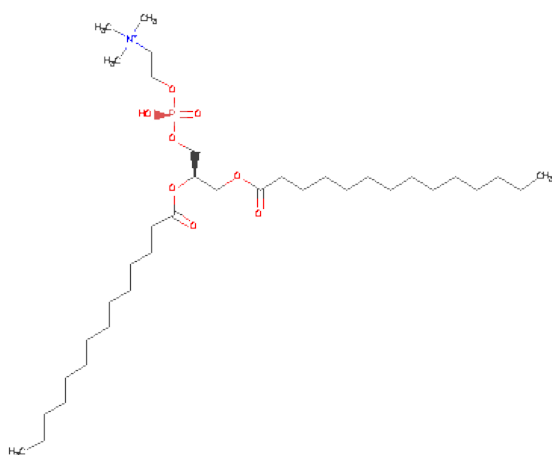
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	13	Total	C	0	0
			114	114		
3	A	8	Total	C	0	0
			62	62		

- Molecule 4 is RETINAL (three-letter code: RET) (formula: $C_{20}H_{28}O$).



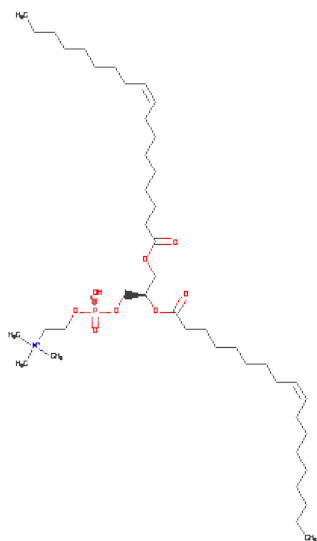
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	C	0	0
			20	20		
4	B	1	Total	C	0	0
			20	20		

- Molecule 5 is 1,2-DIMYRISTOYL-SN-GLYCERO-3-PHOSPHOCHOLINE (three-letter code: PX4) (formula: $C_{36}H_{73}NO_8P$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	B	1	Total	C	O	P	0	0
			34	25	8	1		

- Molecule 6 is 1,2-DIOLEOYL-SN-GLYCERO-3-PHOSPHOCHOLINE (three-letter code: PCW) (formula: $C_{44}H_{85}NO_8P$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	B	1	Total	C	O	P	0	0
			40	31	8	1		

- Molecule 7 is water.

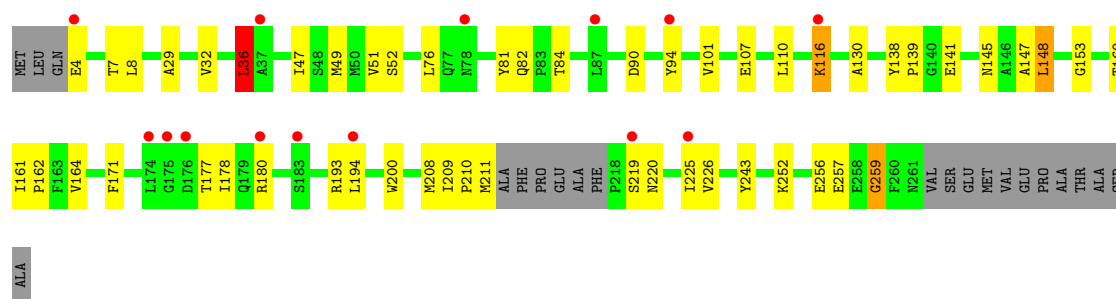
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	34	Total 34	O 34	0	0
7	B	28	Total 28	O 28	0	0

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.

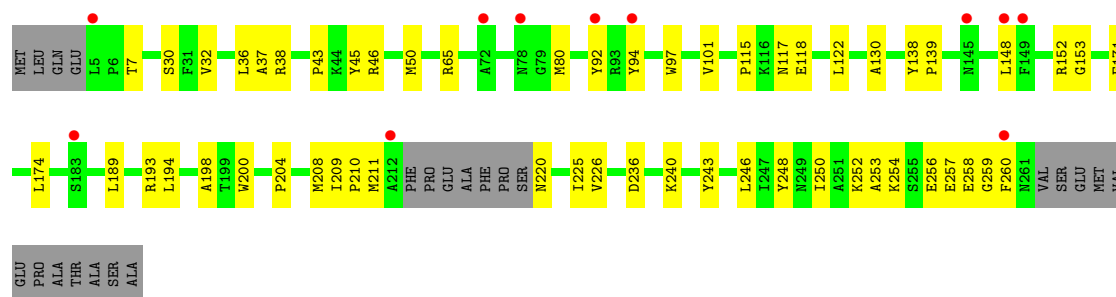
• Molecule 1: Xanthorhodopsin

Chain A: 



• Molecule 1: Xanthorhodopsin

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	52.74Å 59.49Å 59.72Å 76.35° 74.93° 64.08°	Depositor
Resolution (Å)	45.10 – 1.90 45.10 – 1.90	Depositor EDS
% Data completeness (in resolution range)	94.1 (45.10-1.90) 94.1 (45.10-1.90)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.61 (at 1.89Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.247 , 0.265 0.238 , 0.259	Depositor DCC
R_{free} test set	2333 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	35.1	Xtriage
Anisotropy	0.478	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 68.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 46284 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4417	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.24% 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: SXN, UNL, PCW, RET, PX4

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	0.54	0/2024	0.64	1/2761 (0.0%)
1	B	0.55	0/2006	0.62	0/2737
All	All	0.54	0/4030	0.63	1/5498 (0.0%)

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
1	B	0	1
All	All	0	2

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	36	LEU	CA-CB-CG	5.89	128.85	115.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	243	TYR	Sidechain
1	B	243	TYR	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	1971	0	1996	40	0
1	B	1954	0	1982	34	0
2	A	70	0	76	3	0
2	B	70	0	76	4	0
3	A	62	0	0	0	0
3	B	114	0	0	0	0
4	A	20	0	27	1	0
4	B	20	0	27	1	0
5	B	34	0	44	1	0
6	B	40	0	53	2	0
7	A	34	0	0	1	0
7	B	28	0	0	0	0
All	All	4417	0	4281	78	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 (78) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:101:VAL:HG12	1:B:130:ALA:HB1	1.41	0.98
1:A:208:MET:HA	1:A:211:MET:HE3	1.50	0.93
1:A:101:VAL:HG12	1:A:130:ALA:HB1	1.49	0.91
1:A:8:LEU:HD21	1:A:225:ILE:HD11	1.68	0.75
1:A:194:LEU:HD21	2:A:1401:SVN:H8'	1.70	0.74
1:B:252:LYS:O	1:B:256:GLU:HG3	1.88	0.72
1:A:194:LEU:HD21	2:A:1401:SVN:H10'	1.74	0.70
1:A:101:VAL:HG13	1:A:200:TRP:CH2	2.30	0.67
1:A:138:TYR:HB3	1:A:139:PRO:HD3	1.76	0.66
1:B:32:VAL:O	1:B:36:LEU:HD13	1.94	0.66
1:B:198:ALA:HB1	2:B:1401:SVN:H31A	1.78	0.66
1:B:115:PRO:HD2	1:B:118:GLU:OE1	1.97	0.65
1:A:116:LYS:NZ	1:A:116:LYS:HB2	2.13	0.64
1:A:94:TYR:OH	1:A:138:TYR:HB2	1.98	0.64
1:B:101:VAL:HG13	1:B:200:TRP:CH2	2.33	0.64

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:94:TYR:OH	1:B:138:TYR:HB2	2.00	0.62
1:B:38:ARG:O	1:B:46:ARG:HG2	2.00	0.61
1:B:101:VAL:CG1	1:B:130:ALA:HB1	2.24	0.61
1:A:76:LEU:HD13	1:A:81:TYR:CE1	2.35	0.61
1:B:253:ALA:O	1:B:257:GLU:HG3	2.03	0.59
1:A:194:LEU:HD23	1:A:194:LEU:O	2.03	0.58
1:B:236:ASP:O	1:B:240:LYS:HD2	2.02	0.58
1:B:138:TYR:HB3	1:B:139:PRO:HD3	1.86	0.57
1:B:46:ARG:O	1:B:50:MET:HG3	2.07	0.55
1:A:29:ALA:O	1:A:32:VAL:HG22	2.07	0.55
1:B:117:ASN:OD1	1:B:118:GLU:HG3	2.06	0.55
1:B:148:LEU:O	1:B:153:GLY:HA3	2.06	0.55
1:A:101:VAL:CG1	1:A:130:ALA:HB1	2.30	0.55
1:B:148:LEU:HD13	1:B:148:LEU:C	2.28	0.54
1:B:152:ARG:HG2	1:B:211:MET:CE	2.38	0.53
1:A:209:ILE:HB	1:A:210:PRO:HD3	1.89	0.53
1:A:145:ASN:HD21	1:A:147:ALA:HB3	1.75	0.52
1:A:171:PHE:CZ	1:A:193:ARG:HG2	2.46	0.51
1:A:180:ARG:O	1:A:180:ARG:HD2	2.11	0.51
1:A:84:THR:O	1:A:84:THR:HG22	2.10	0.51
1:A:178:ILE:HD12	1:A:178:ILE:H	1.76	0.51
1:B:122:LEU:HD21	1:B:174:LEU:HD11	1.92	0.50
1:B:258:GLU:OE2	1:B:260:PHE:HE2	1.94	0.50
1:A:52:SER:HA	1:A:107:GLU:OE2	2.12	0.50
5:B:1415:PX4:H47	5:B:1415:PX4:H20	1.94	0.50
1:A:145:ASN:HD21	1:A:147:ALA:CB	2.25	0.49
1:A:225:ILE:HG13	1:A:226:VAL:N	2.27	0.49
1:B:208:MET:O	1:B:211:MET:HG2	2.13	0.49
1:B:65:ARG:NH2	1:B:92:TYR:OH	2.47	0.48
1:A:178:ILE:HD12	1:A:178:ILE:N	2.28	0.48
1:B:171:PHE:CZ	1:B:193:ARG:HG2	2.49	0.47
1:B:209:ILE:N	1:B:210:PRO:HD2	2.29	0.47
1:A:252:LYS:O	1:A:256:GLU:HG3	2.15	0.47
1:A:47:ILE:HG23	1:A:110:LEU:HD13	1.96	0.47
1:A:116:LYS:HZ3	1:A:116:LYS:HB2	1.79	0.46
1:B:97:TRP:CD1	4:B:1301:RET:H14	2.49	0.46
1:B:37:ALA:HB1	1:B:248:TYR:CE1	2.51	0.46
1:A:180:ARG:HH21	1:A:180:ARG:HG3	1.82	0.45
2:B:1401:SXN:O4*	6:B:1416:PCW:H283	2.17	0.45
1:B:152:ARG:HG2	1:B:211:MET:HE1	1.98	0.45
1:A:49:MET:CE	1:A:49:MET:HA	2.46	0.45
1:B:45:TYR:OH	1:B:254:LYS:HE3	2.17	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:189:LEU:HD23	1:B:250:ILE:HD13	1.99	0.44
1:A:32:VAL:O	1:A:36:LEU:HB2	2.17	0.44
1:A:148:LEU:O	1:A:153:GLY:HA3	2.18	0.44
1:A:194:LEU:CD2	2:A:1401:SNX:H8'	2.43	0.43
1:A:47:ILE:O	1:A:51:VAL:HG23	2.18	0.43
2:B:1401:SNX:H19	2:B:1401:SNX:H7	1.79	0.43
1:A:7:THR:HB	1:A:82:GLN:OE1	2.18	0.43
1:B:200:TRP:O	1:B:204:PRO:HD2	2.18	0.43
1:B:258:GLU:HB2	1:B:260:PHE:CD2	2.53	0.43
1:B:7:THR:HG22	1:B:80:MET:HE3	1.99	0.43
1:A:145:ASN:ND2	1:A:147:ALA:HB3	2.34	0.42
1:B:246:LEU:CD2	6:B:1416:PCW:H261	2.49	0.42
1:A:160:THR:O	1:A:164:VAL:HG13	2.20	0.42
1:A:180:ARG:NH2	1:A:180:ARG:HG3	2.35	0.41
1:A:194:LEU:C	1:A:194:LEU:HD23	2.41	0.41
4:A:1301:RET:H7	4:A:1301:RET:H181	1.78	0.41
1:A:257:GLU:C	1:A:259:GLY:H	2.23	0.41
1:A:7:THR:HG22	7:A:430:HOH:O	2.21	0.41
1:A:161:ILE:HB	1:A:162:PRO:CD	2.51	0.41
1:B:194:LEU:CD2	2:B:1401:SNX:H6'	2.52	0.40
1:B:225:ILE:HG23	1:B:226:VAL:N	2.35	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	248/273 (91%)	237 (96%)	9 (4%)	2 (1%)	27	12
1	B	246/273 (90%)	241 (98%)	4 (2%)	1 (0%)	43	29
All	All	494/546 (90%)	478 (97%)	13 (3%)	3 (1%)	33	19

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	219	SER
1	A	259	GLY
1	B	259	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	207/223 (93%)	199 (96%)	8 (4%)	43	30
1	B	204/223 (92%)	201 (98%)	3 (2%)	76	73
All	All	411/446 (92%)	400 (97%)	11 (3%)	57	47

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	GLU
1	A	36	LEU
1	A	90	ASP
1	A	116	LYS
1	A	141	GLU
1	A	148	LEU
1	A	177	THR
1	A	220	ASN
1	B	30	SER
1	B	43	PRO
1	B	220	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	ASN
1	A	77	GLN
1	A	145	ASN
1	A	220	ASN
1	B	39	ASN
1	B	77	GLN

5.3.3 RNA ⓘ

There are no RNA chains 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 27 ligands modelled in this entry, 21 are unknown - 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$
4	RET	A	1301	1	19,20,21	1.72	6 (31%)	26,27,28	2.26	10 (38%)
2	SXN	A	1401	-	56,57,71	3.35	18 (32%)	66,70,94	1.97	13 (19%)
4	RET	B	1301	1	19,20,21	1.70	4 (21%)	26,27,28	2.40	12 (46%)
2	SXN	B	1401	-	56,57,71	3.42	17 (30%)	66,70,94	1.97	12 (18%)
5	PX4	B	1415	-	33,33,45	2.04	4 (12%)	38,38,53	1.22	3 (7%)
6	PCW	B	1416	-	38,39,53	1.30	3 (7%)	43,44,61	1.27	5 (11%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	RET	A	1301	1	-	0/13/30/31	0/1/1/1
2	SXN	A	1401	-	-	1/51/72/105	0/1/1/2
4	RET	B	1301	1	-	0/13/30/31	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SXN	B	1401	-	-	1/51/72/105	0/1/1/2
5	PX4	B	1415	-	1/1/4/5	0/35/35/49	0/0/0/0
6	PCW	B	1416	-	1/1/4/9	0/41/41/57	0/0/0/0

All (52) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1401	SXN	O37-C22	17.67	1.23	1.11
2	A	1401	SXN	O37-C22	17.06	1.23	1.11
5	B	1415	PX4	C16-C15	-8.80	1.51	1.55
2	A	1401	SXN	C10'-C9'	6.80	1.44	1.35
2	B	1401	SXN	C14'-C13'	6.61	1.44	1.35
2	A	1401	SXN	C6'-C5'	6.51	1.44	1.35
2	B	1401	SXN	C10'-C9'	6.37	1.44	1.35
2	B	1401	SXN	C6'-C5'	6.23	1.44	1.35
2	A	1401	SXN	C14'-C13'	6.13	1.43	1.35
2	B	1401	SXN	C14-C13	5.57	1.43	1.35
2	A	1401	SXN	C14-C13	5.53	1.43	1.35
2	B	1401	SXN	C10-C9	4.99	1.42	1.35
2	A	1401	SXN	C10-C9	4.97	1.42	1.35
2	B	1401	SXN	C6-C5	4.82	1.42	1.35
2	A	1401	SXN	C6-C5	4.81	1.42	1.35
5	B	1415	PX4	O5-C9	4.64	1.47	1.33
2	B	1401	SXN	C1-C6	4.58	1.60	1.53
6	B	1416	PCW	O2-C31	4.53	1.48	1.34
6	B	1416	PCW	O3-C11	4.44	1.47	1.33
5	B	1415	PX4	O7-C23	4.37	1.47	1.34
2	B	1401	SXN	C4-C5	-4.24	1.39	1.48
4	B	1301	RET	C1-C6	4.15	1.59	1.53
2	A	1401	SXN	C4-C5	-4.07	1.39	1.48
6	B	1416	PCW	C20-C19	3.98	1.53	1.31
2	B	1401	SXN	C16'-C1'	3.79	1.55	1.46
2	A	1401	SXN	C16'-C1'	3.69	1.55	1.46
5	B	1415	PX4	P1-O2	3.68	1.62	1.50
2	A	1401	SXN	C4'-C3'	3.49	1.41	1.32
4	A	1301	RET	C1-C6	3.46	1.58	1.53
2	A	1401	SXN	C1-C6	3.35	1.58	1.53
2	B	1401	SXN	C17'-C1'	3.17	1.54	1.46
2	A	1401	SXN	C17'-C1'	3.16	1.54	1.46
2	B	1401	SXN	C4'-C3'	3.13	1.40	1.32
4	A	1301	RET	C17-C1	3.07	1.60	1.53
4	B	1301	RET	C17-C1	3.01	1.60	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1401	SXN	C7'-C8'	2.99	1.42	1.34
2	A	1401	SXN	C11'-C12'	2.96	1.42	1.34
2	B	1401	SXN	C11'-C12'	2.90	1.42	1.34
2	B	1401	SXN	C8-C7	2.85	1.41	1.32
4	A	1301	RET	C5-C6	2.85	1.38	1.34
2	A	1401	SXN	C8-C7	2.85	1.41	1.32
2	B	1401	SXN	C7'-C8'	2.71	1.41	1.34
4	A	1301	RET	C10-C9	2.43	1.38	1.35
4	B	1301	RET	C5-C6	2.42	1.38	1.34
2	A	1401	SXN	C11-C12	2.34	1.40	1.34
2	B	1401	SXN	C11-C12	2.29	1.40	1.34
2	A	1401	SXN	C23-C22	2.27	1.54	1.47
4	A	1301	RET	C16-C1	2.27	1.58	1.53
4	A	1301	RET	C11-C12	2.22	1.40	1.34
4	B	1301	RET	C16-C1	2.19	1.58	1.53
2	A	1401	SXN	C2'-C3'	2.17	1.59	1.49
2	B	1401	SXN	C2'-C3'	2.09	1.58	1.49

All (55) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1401	SXN	C17'-C1'-C2'	-7.83	110.11	121.14
2	A	1401	SXN	C17'-C1'-C2'	-7.72	110.27	121.14
2	B	1401	SXN	C16'-C1'-C2'	-6.51	111.97	121.14
2	A	1401	SXN	C16'-C1'-C2'	-6.42	112.09	121.14
4	B	1301	RET	C11-C10-C9	-5.54	119.31	127.29
4	B	1301	RET	C18-C5-C6	5.18	130.38	124.51
4	A	1301	RET	C11-C10-C9	-4.72	120.49	127.29
4	A	1301	RET	C18-C5-C6	4.67	129.81	124.51
2	A	1401	SXN	C11-C10-C9	-4.13	121.34	127.29
4	A	1301	RET	C8-C9-C10	-3.92	112.95	118.97
4	B	1301	RET	C8-C9-C10	-3.77	113.18	118.97
2	A	1401	SXN	C7-C8-C9	-3.69	120.69	126.22
2	B	1401	SXN	C11-C10-C9	-3.68	122.00	127.29
5	B	1415	PX4	O7-C23-C24	3.63	119.51	111.56
4	A	1301	RET	C7-C6-C5	-3.61	112.94	121.57
6	B	1416	PCW	O2-C31-C32	3.56	119.36	111.56
2	B	1401	SXN	C7-C8-C9	-3.55	120.90	126.22
4	B	1301	RET	C7-C6-C5	-3.41	113.42	121.57
2	A	1401	SXN	C16'-C1'-C17'	-3.31	108.29	119.11
4	A	1301	RET	C1-C6-C7	3.23	124.65	115.69
2	B	1401	SXN	C16'-C1'-C17'	-3.22	108.58	119.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1401	SXN	C15-C14-C13	-3.21	122.67	127.29
4	B	1301	RET	C1-C6-C7	3.11	124.30	115.69
2	A	1401	SXN	C18-C5-C6	-3.08	120.42	124.22
2	B	1401	SXN	C18-C5-C6	-3.05	120.46	124.22
2	B	1401	SXN	C3-C4-C5	3.03	121.74	118.76
2	B	1401	SXN	C15-C14-C13	-2.96	123.03	127.29
2	A	1401	SXN	C15'-C14'-C13'	-2.92	123.08	127.29
2	B	1401	SXN	C15'-C14'-C13'	-2.86	123.17	127.29
6	B	1416	PCW	C18-C19-C20	-2.83	109.39	125.43
4	B	1301	RET	C20-C13-C12	2.79	122.61	118.09
2	A	1401	SXN	C3-C4-C5	2.73	121.44	118.76
4	A	1301	RET	C20-C13-C12	2.64	122.36	118.09
6	B	1416	PCW	C21-C20-C19	-2.64	110.49	125.43
5	B	1415	PX4	O5-C9-C10	2.57	120.02	111.94
6	B	1416	PCW	O3-C11-C12	2.53	119.90	111.94
4	A	1301	RET	C10-C11-C12	-2.44	115.00	123.24
2	B	1401	SXN	C2'-C3'-C4'	-2.41	119.24	124.01
4	B	1301	RET	C19-C9-C8	2.36	121.90	118.09
4	B	1301	RET	C10-C11-C12	-2.35	115.32	123.24
4	B	1301	RET	C17-C1-C6	2.30	114.14	110.33
2	A	1401	SXN	C2'-C3'-C4'	-2.26	119.53	124.01
4	B	1301	RET	C14-C13-C12	-2.26	111.77	119.62
2	B	1401	SXN	C18'-C5'-C4'	2.26	121.74	118.09
2	B	1401	SXN	C19'-C9'-C8'	2.25	121.74	118.09
4	B	1301	RET	C18-C5-C4	-2.24	109.23	113.34
2	A	1401	SXN	C19'-C9'-C8'	2.21	121.67	118.09
4	A	1301	RET	C19-C9-C8	2.18	121.62	118.09
5	B	1415	PX4	C16-C15-C14	-2.15	109.62	114.46
4	A	1301	RET	C14-C13-C12	-2.15	112.15	119.62
2	A	1401	SXN	C18'-C5'-C4'	2.09	121.48	118.09
2	A	1401	SXN	C18-C5-C4	2.05	118.46	115.38
4	B	1301	RET	C11-C12-C13	-2.04	120.55	126.38
6	B	1416	PCW	C2-O2-C31	2.04	122.94	117.92
4	A	1301	RET	C17-C1-C6	2.01	113.67	110.33

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
5	B	1415	PX4	C7
6	B	1416	PCW	C2

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1401	SXN	C3'-C2'-C1'-C16'
2	A	1401	SXN	C3'-C2'-C1'-C16'

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	252/273 (92%)	0.53	14 (5%) 24 24	28, 41, 68, 76	0
1	B	250/273 (91%)	0.46	11 (4%) 33 33	29, 40, 63, 76	0
All	All	502/546 (91%)	0.49	25 (4%) 28 28	28, 40, 65, 76	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	212	ALA	5.1
1	A	176	ASP	4.0
1	A	78	ASN	3.8
1	B	72	ALA	3.7
1	B	94	TYR	3.6
1	A	94	TYR	3.6
1	A	4	GLU	3.5
1	A	37	ALA	3.4
1	B	5	LEU	3.3
1	B	260	PHE	3.1
1	A	225	ILE	3.1
1	B	145	ASN	2.9
1	A	116	LYS	2.8
1	A	219	SER	2.8
1	A	180	ARG	2.8
1	A	194	LEU	2.7
1	B	78	ASN	2.6
1	A	175	GLY	2.5
1	A	183	SER	2.4
1	A	87	LEU	2.4
1	B	149	PHE	2.4
1	B	183	SER	2.4
1	B	92	TYR	2.3
1	B	148	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	174	LEU	2.1

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

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

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	PCW	B	1416	40/54	0.31	12.76	56,77,91,92	0
3	UNL	B	1414	8/-	0.32	9.71	63,65,67,69	0
3	UNL	A	1405	7/-	0.27	8.28	67,68,68,69	0
3	UNL	A	1404	9/-	0.35	8.15	74,74,75,76	0
3	UNL	B	1409	7/-	0.28	8.04	68,68,69,70	0
3	UNL	A	1408	7/-	0.24	7.98	62,62,64,65	0
3	UNL	B	1403	16/-	0.22	7.09	47,50,52,53	0
3	UNL	B	1413	7/-	0.33	6.57	71,73,74,76	0
3	UNL	B	1404	7/-	0.27	6.11	66,69,74,74	0
3	UNL	B	1402	10/-	0.20	5.45	51,54,57,58	0
2	SXN	A	1401	70/70	0.34	4.87	53,79,92,94	0
3	UNL	A	1407	7/-	0.22	3.81	65,66,66,66	0
3	UNL	A	1406	12/-	0.20	3.48	65,67,68,68	0
3	UNL	A	1403	7/-	0.30	3.41	54,58,60,60	0
3	UNL	B	1411	6/-	0.20	3.34	54,56,60,60	0
3	UNL	B	1412	11/-	0.21	3.01	63,65,67,68	0
3	UNL	B	1405	7/-	0.30	2.81	66,67,69,69	0
3	UNL	A	1409	6/-	0.27	2.36	57,57,58,59	0
3	UNL	A	1402	7/-	0.16	2.34	53,53,55,57	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	SXN	B	1401	70/70	0.24	2.32	43,76,90,91	0
3	UNL	B	1406	12/-	0.15	2.02	70,72,74,75	0
5	PX4	B	1415	34/46	0.23	1.85	57,70,81,83	0
3	UNL	B	1408	7/-	0.20	1.53	73,73,74,75	0
3	UNL	B	1410	8/-	0.23	1.22	61,62,63,63	0
4	RET	B	1301	20/21	0.15	0.94	30,32,35,35	0
4	RET	A	1301	20/21	0.15	0.51	29,33,36,37	0
3	UNL	B	1407	8/-	0.45	-	71,72,73,73	0

6.5 Other polymers ⓘ

There are no such residues in this entry.