



# Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 8D6G  
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Deposited on : 2022-06-06  
Resolution : 2.20 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](mailto: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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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)  
Xtrriage (Phenix) : 1.13  
EDS : 2.35.1  
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) : 2.35.1

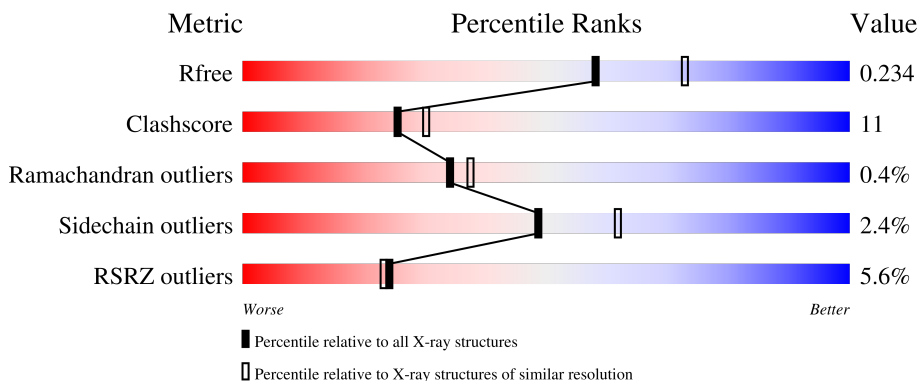
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.20 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	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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 of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	854	

## 2 Entry composition [i](#)

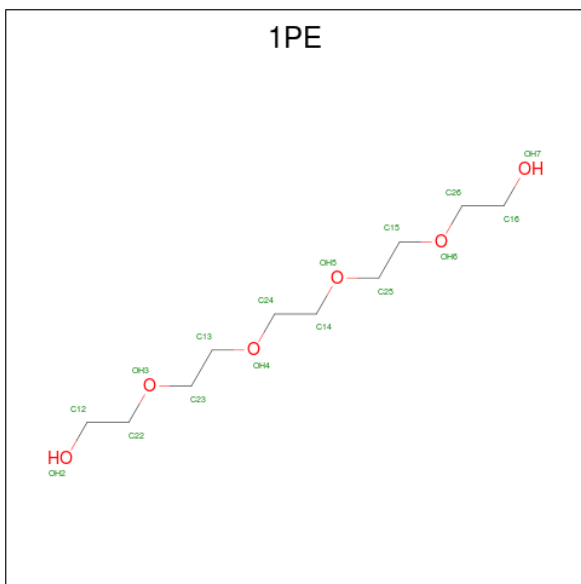
There are 3 unique types of molecules in this entry. The entry contains 6656 atoms, of which 22 are hydrogens and 0 are deuteriums.

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 Saxiphilin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	821	6373	3989	1102	1223	59	0	1	0

- Molecule 2 is PENTAETHYLENE GLYCOL (three-letter code: 1PE) (formula: C<sub>10</sub>H<sub>22</sub>O<sub>6</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	H	O		
2	A	1	38	10	22	6	0	0
2	A	1	16	10	6	6	0	0

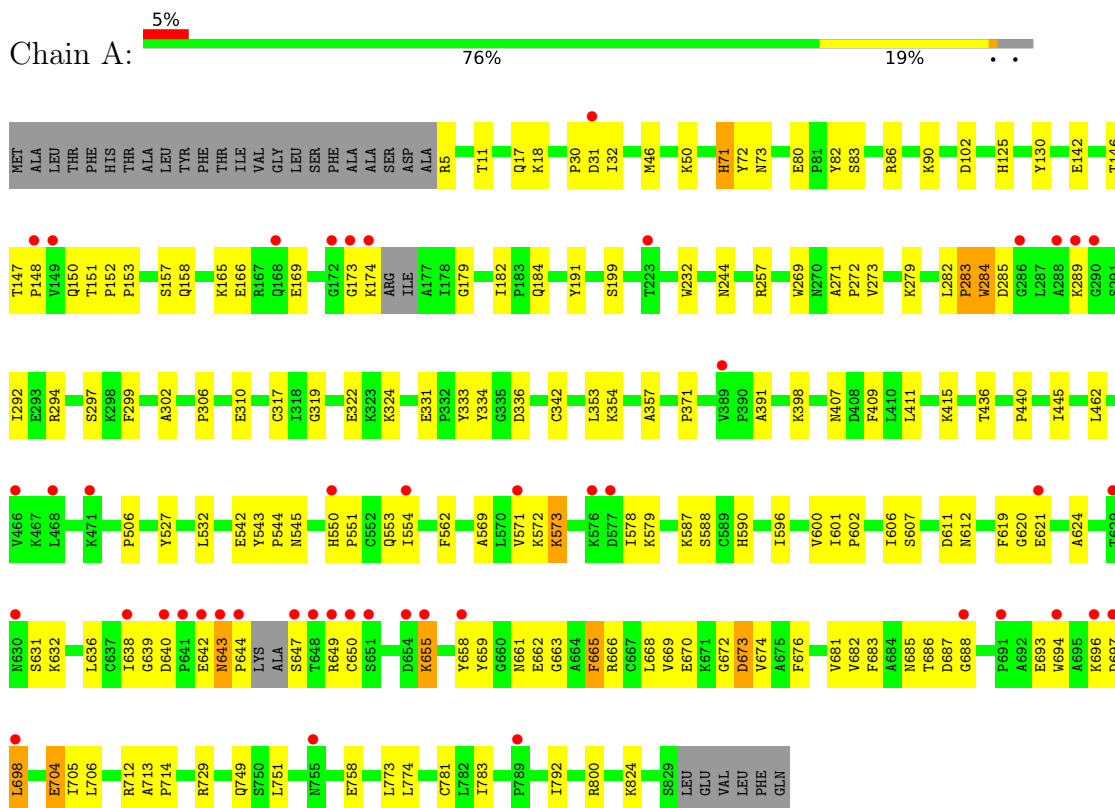
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	229	Total	O	0	0
			229	229		

### 3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: Saxiphilin



## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	229.05Å 229.05Å 67.43Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	43.29 – 2.20 43.29 – 2.20	Depositor EDS
% Data completeness (in resolution range)	98.6 (43.29-2.20) 98.6 (43.29-2.20)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.28 (at 2.20Å)	Xtrriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, $R_{free}$	0.189 , 0.235 0.188 , 0.234	Depositor DCC
$R_{free}$ test set	3252 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	52.7	Xtrriage
Anisotropy	0.270	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 46.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	0.005 for h,-h-k,-l	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	6656	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	66.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.85% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: 1PE

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	0.36	0/6518	0.56	0/8815

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 [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	6373	0	6164	134	0
2	A	32	22	44	1	0
3	A	229	0	0	2	0
All	All	6634	22	6208	134	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 11.

All (134) 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:90:LYS:NZ	3:A:1001:HOH:O	1.89	1.05

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:662:GLU:HG3	1:A:685:ASN:HD21	1.26	0.95
1:A:693:GLU:OE2	1:A:693:GLU:N	2.03	0.90
1:A:571:VAL:HG12	1:A:673:ASP:O	1.74	0.88
1:A:292:ILE:HD12	1:A:462:LEU:HD22	1.55	0.87
1:A:292:ILE:CD1	1:A:462:LEU:HD22	2.06	0.86
1:A:662:GLU:HG3	1:A:685:ASN:ND2	1.95	0.81
1:A:661:ASN:HB3	1:A:681:VAL:HG22	1.65	0.79
1:A:130:TYR:CE1	1:A:147:THR:HG22	2.22	0.74
1:A:693:GLU:HA	1:A:696:LYS:HG3	1.66	0.74
1:A:80:GLU:OE2	1:A:436:THR:HG22	1.86	0.74
1:A:86:ARG:HB2	1:A:232:TRP:CD1	2.24	0.72
1:A:638:ILE:HG13	1:A:658:TYR:CE1	2.25	0.72
1:A:682:VAL:O	1:A:686:THR:HG23	1.90	0.71
1:A:279:LYS:HE3	1:A:371:PRO:O	1.92	0.69
1:A:30:PRO:O	1:A:31:ASP:HB2	1.92	0.69
1:A:602:PRO:O	1:A:606:ILE:HD12	1.93	0.69
1:A:147:THR:HB	1:A:148:PRO:HD2	1.73	0.68
1:A:179:GLY:HA2	1:A:199:SER:OG	1.92	0.68
1:A:693:GLU:CB	1:A:696:LYS:HE3	2.24	0.67
1:A:147:THR:H	1:A:150:GLN:NE2	1.95	0.64
1:A:542:GLU:OE1	1:A:729:ARG:HD3	1.97	0.64
1:A:407:ASN:ND2	1:A:440:PRO:HB3	2.13	0.64
1:A:661:ASN:HB3	1:A:681:VAL:CG2	2.29	0.63
1:A:687:ASP:N	1:A:698:LEU:O	2.33	0.62
1:A:607:SER:HB2	1:A:612:ASN:HA	1.80	0.61
1:A:165:LYS:NZ	1:A:169:GLU:HG2	2.15	0.61
1:A:319:GLY:O	1:A:324:LYS:HB2	2.01	0.61
1:A:151:THR:HB	1:A:152:PRO:HD2	1.83	0.60
1:A:553:GLN:HB3	1:A:554:ILE:HD12	1.82	0.60
1:A:621:GLU:HB3	1:A:636:LEU:HD13	1.86	0.58
1:A:693:GLU:HB2	1:A:696:LYS:HE3	1.85	0.58
1:A:590:HIS:HD2	1:A:624:ALA:HB2	1.66	0.58
1:A:506:PRO:HG2	1:A:527:TYR:CE2	2.39	0.58
1:A:294:ARG:HG3	1:A:310:GLU:OE2	2.04	0.57
1:A:661:ASN:O	1:A:681:VAL:HG11	2.04	0.57
1:A:682:VAL:HG13	1:A:683:PHE:N	2.20	0.57
1:A:571:VAL:HG21	1:A:706:LEU:HD21	1.86	0.57
1:A:640:ASP:OD2	1:A:642:GLU:HG3	2.05	0.57
1:A:317:CYS:HB3	1:A:331:GLU:OE1	2.05	0.56
1:A:284:TRP:HA	1:A:289:LYS:HD2	1.87	0.56
1:A:166:GLU:OE1	1:A:191:TYR:OH	2.14	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:666:ARG:HH21	1:A:694:TRP:HB3	1.71	0.55
1:A:411:LEU:HD23	1:A:411:LEU:O	2.06	0.55
1:A:621:GLU:HB3	1:A:636:LEU:CD1	2.37	0.55
1:A:704:GLU:OE2	1:A:712:ARG:NH2	2.40	0.54
1:A:587:LYS:HA	1:A:621:GLU:O	2.08	0.54
1:A:543:TYR:CZ	1:A:545:ASN:HB3	2.44	0.53
1:A:642:GLU:O	1:A:642:GLU:CD	2.47	0.52
1:A:102:ASP:OD2	1:A:800:ARG:NE	2.37	0.51
1:A:639:GLY:HA3	1:A:649:ARG:HG2	1.91	0.51
1:A:640:ASP:OD2	1:A:655:LYS:NZ	2.43	0.51
1:A:590:HIS:CD2	1:A:624:ALA:HB2	2.45	0.51
1:A:411:LEU:HD23	1:A:411:LEU:C	2.31	0.51
1:A:257:ARG:HA	1:A:302:ALA:O	2.11	0.50
1:A:283:PRO:O	1:A:284:TRP:HB2	2.10	0.50
1:A:50:LYS:HE2	1:A:71:HIS:O	2.12	0.50
1:A:173:GLY:O	1:A:174:LYS:HG3	2.12	0.50
1:A:32:ILE:HG21	1:A:409:PHE:HB2	1.94	0.50
1:A:147:THR:H	1:A:150:GLN:HE21	1.60	0.50
1:A:644:PRO:HG2	1:A:647:SER:HA	1.93	0.50
1:A:130:TYR:CZ	1:A:147:THR:HG22	2.46	0.49
1:A:151:THR:HB	1:A:152:PRO:CD	2.42	0.49
1:A:125:HIS:ND1	1:A:611:ASP:OD1	2.37	0.49
1:A:554:ILE:HD12	1:A:554:ILE:N	2.28	0.49
1:A:146:THR:OG1	1:A:153:PRO:HD3	2.13	0.48
1:A:665:PHE:CE1	1:A:685:ASN:HB3	2.48	0.48
1:A:682:VAL:HG13	1:A:683:PHE:H	1.78	0.48
1:A:543:TYR:CE2	1:A:551:PRO:HG2	2.48	0.48
1:A:688:GLY:CA	1:A:697:ASP:HA	2.44	0.48
1:A:306:PRO:HG2	1:A:334:TYR:HA	1.96	0.47
1:A:620:GLY:O	1:A:632:LYS:HE2	2.14	0.47
1:A:336:ASP:HB3	1:A:357:ALA:HB2	1.96	0.47
1:A:569:ALA:O	1:A:705:ILE:HA	2.14	0.47
1:A:82:TYR:CE2	1:A:391:ALA:HB2	2.50	0.47
1:A:165:LYS:NZ	1:A:169:GLU:CG	2.78	0.46
1:A:668:LEU:HD12	1:A:672:GLY:O	2.15	0.46
1:A:553:GLN:HB3	1:A:554:ILE:CD1	2.45	0.46
1:A:619:PHE:O	1:A:632:LYS:NZ	2.41	0.46
1:A:336:ASP:HB3	1:A:357:ALA:CB	2.46	0.46
1:A:532:LEU:HD22	1:A:792:ILE:HG13	1.97	0.46
1:A:749:GLN:OE1	1:A:781:CYS:HB2	2.15	0.46
1:A:269:TRP:O	1:A:272:PRO:HD2	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:284:TRP:HA	1:A:289:LYS:CD	2.45	0.46
1:A:751:LEU:HD22	1:A:758:GLU:HG3	1.98	0.46
1:A:142:GLU:OE1	1:A:142:GLU:HA	2.14	0.46
1:A:173:GLY:O	1:A:174:LYS:CB	2.63	0.46
1:A:573:LYS:HD2	1:A:704:GLU:CD	2.37	0.45
1:A:294:ARG:O	1:A:297:SER:HB3	2.17	0.45
1:A:590:HIS:CE1	1:A:676:PHE:CD2	3.04	0.45
1:A:544:PRO:HD2	1:A:562:PHE:O	2.17	0.45
1:A:596:ILE:HA	1:A:600:VAL:HB	1.99	0.45
1:A:573:LYS:HD2	1:A:704:GLU:OE2	2.17	0.45
1:A:18:LYS:HE2	1:A:18:LYS:HB2	1.76	0.44
1:A:658:TYR:OH	1:A:670:GLU:OE2	2.23	0.44
1:A:685:ASN:HD22	1:A:694:TRP:HH2	1.64	0.44
1:A:572:LYS:HE3	1:A:669:VAL:HA	1.99	0.44
1:A:11:THR:HG21	1:A:17:GLN:HA	1.99	0.44
1:A:588:SER:HA	1:A:674:VAL:O	2.18	0.44
1:A:578:ILE:O	1:A:578:ILE:HG23	2.18	0.44
1:A:46:MET:HE2	1:A:72:TYR:CD2	2.52	0.43
1:A:73:ASN:OD1	1:A:398:LYS:HE3	2.18	0.43
1:A:284:TRP:CZ2	1:A:292:ILE:HG13	2.53	0.43
1:A:285:ASP:N	1:A:285:ASP:OD1	2.50	0.43
1:A:244:ASN:ND2	3:A:1004:HOH:O	2.31	0.43
1:A:273:VAL:HG11	1:A:292:ILE:HG21	2.00	0.43
1:A:587:LYS:HD3	1:A:621:GLU:HB2	1.99	0.43
1:A:655:LYS:HA	1:A:655:LYS:HD2	1.72	0.43
1:A:282:LEU:HD11	1:A:299:PHE:HB2	2.01	0.43
1:A:310:GLU:HA	1:A:310:GLU:OE1	2.19	0.42
1:A:550:HIS:N	1:A:551:PRO:HD2	2.34	0.42
1:A:80:GLU:OE2	1:A:436:THR:CG2	2.62	0.42
1:A:572:LYS:NZ	1:A:669:VAL:HA	2.34	0.42
1:A:659:TYR:O	1:A:663:GLY:HA3	2.20	0.42
1:A:269:TRP:C	1:A:272:PRO:HD2	2.40	0.42
1:A:693:GLU:CA	1:A:696:LYS:HG3	2.45	0.42
1:A:693:GLU:O	1:A:696:LYS:HG3	2.19	0.42
1:A:182:ILE:O	1:A:184:GLN:NE2	2.49	0.42
1:A:562:PHE:H	2:A:902:1PE:H252	1.84	0.42
1:A:322:GLU:OE2	1:A:322:GLU:N	2.49	0.42
1:A:407:ASN:HD21	1:A:440:PRO:HB3	1.84	0.42
1:A:642:GLU:O	1:A:643:ASN:C	2.59	0.41
1:A:5:ARG:HB2	1:A:31:ASP:O	2.19	0.41
1:A:333:TYR:CG	1:A:342:CYS:HB2	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:601:ILE:HB	1:A:602:PRO:HD3	2.01	0.41
1:A:665:PHE:CE2	1:A:669:VAL:HG13	2.56	0.41
1:A:551:PRO:HB2	1:A:783:ILE:HD11	2.03	0.41
1:A:773:LEU:O	1:A:774:LEU:HB2	2.21	0.41
1:A:165:LYS:HA	1:A:165:LYS:HD2	1.81	0.41
1:A:665:PHE:C	1:A:665:PHE:CD2	2.94	0.41
1:A:713:ALA:HB1	1:A:714:PRO:HD2	2.02	0.41
1:A:271:ALA:HB3	1:A:272:PRO:HD3	2.03	0.41
1:A:554:ILE:O	1:A:554:ILE:HG22	2.21	0.40
1:A:336:ASP:OD2	1:A:354:LYS:HG2	2.21	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	816/854 (96%)	774 (95%)	39 (5%)	3 (0%)	34 37

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	284	TRP
1	A	283	PRO
1	A	643	ASN

### 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	706/731 (97%)	689 (98%)	17 (2%)	49 62

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	71	HIS
1	A	83	SER
1	A	157	SER
1	A	158	GLN
1	A	353	LEU
1	A	415	LYS
1	A	445	ILE
1	A	573	LYS
1	A	579	LYS
1	A	631	SER
1	A	650	CYS
1	A	655	LYS
1	A	665	PHE
1	A	673	ASP
1	A	698	LEU
1	A	704	GLU
1	A	824	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	150	GLN
1	A	407	ASN
1	A	590	HIS
1	A	608	ASN
1	A	685	ASN

### 5.3.3 RNA [i](#)

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](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

2 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	1PE	A	902	-	15,15,15	0.53	0	14,14,14	0.33	0
2	1PE	A	901	-	15,15,15	0.49	0	14,14,14	0.27	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	1PE	A	902	-	-	5/13/13/13	-
2	1PE	A	901	-	-	4/13/13/13	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	901	1PE	OH7-C16-C26-OH6
2	A	901	1PE	OH5-C14-C24-OH4
2	A	902	1PE	OH4-C13-C23-OH3
2	A	902	1PE	OH7-C16-C26-OH6
2	A	901	1PE	OH6-C15-C25-OH5
2	A	901	1PE	C25-C15-OH6-C26

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Mol	Chain	Res	Type	Atoms
2	A	902	1PE	C24-C14-OH5-C25
2	A	902	1PE	OH5-C14-C24-OH4
2	A	902	1PE	C14-C24-OH4-C13

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	902	1PE	1	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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	821/854 (96%)	0.15	46 (5%) <span style="border: 1px solid red; padding: 2px;">24</span> <span style="border: 1px solid red; padding: 2px;">23</span>	39, 59, 113, 141	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	174	LYS	8.1
1	A	691	PRO	7.0
1	A	644	PRO	5.7
1	A	697	ASP	5.6
1	A	638	ILE	5.2
1	A	640	ASP	5.0
1	A	641	PRO	4.7
1	A	466	VAL	4.6
1	A	696	LYS	4.4
1	A	290	GLY	4.3
1	A	648	THR	4.1
1	A	173	GLY	4.0
1	A	698	LEU	3.9
1	A	647	SER	3.8
1	A	288	ALA	3.7
1	A	149	VAL	3.7
1	A	755	ASN	3.7
1	A	649	ARG	3.7
1	A	694	TRP	3.6
1	A	289	LYS	3.5
1	A	688	GLY	3.4
1	A	168	GLN	3.4
1	A	172	GLY	3.2
1	A	789	PRO	3.0
1	A	643	ASN	3.0
1	A	576	LYS	2.9
1	A	630	ASN	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	629	THR	2.9
1	A	577	ASP	2.9
1	A	286	GLY	2.8
1	A	148	PRO	2.8
1	A	31	ASP	2.7
1	A	650	CYS	2.6
1	A	554	ILE	2.6
1	A	471	LYS	2.5
1	A	642	GLU	2.4
1	A	223	THR	2.4
1	A	621	GLU	2.2
1	A	550	HIS	2.2
1	A	389	VAL	2.2
1	A	655	LYS	2.2
1	A	651	SER	2.2
1	A	658	TYR	2.1
1	A	654	ASP	2.1
1	A	571	VAL	2.1
1	A	468	LEU	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 monosaccharides 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, 95<sup>th</sup> 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(Å <sup>2</sup> )	Q<0.9
2	1PE	A	901	16/16	0.81	0.18	65,86,108,109	0
2	1PE	A	902	16/16	0.83	0.17	74,90,99,100	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.