



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:30 PM UTC

PDB ID : 3GUU / pdb_00003guu
Title : X-ray structure of Candida Antarctica lipase A
Authors : Brandt, A.-M.; Li, X.-G.; Nymalm-Rejstrom, Y.; Airene, T.; Kanerva, L.T.;
Salminen, T.A.
Deposited on : 2009-03-30
Resolution : 2.10 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 types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

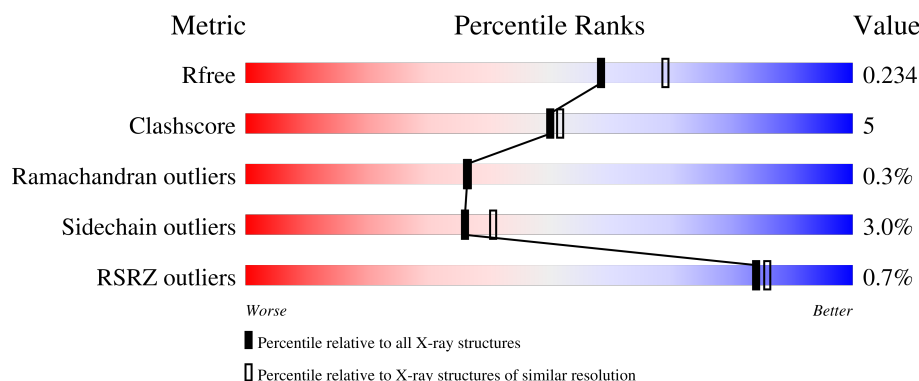
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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}	180053	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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 ≥ 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	462	
1	B	462	

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
3	GOL	A	443	-	-	X	-

2 Entry composition [i](#)

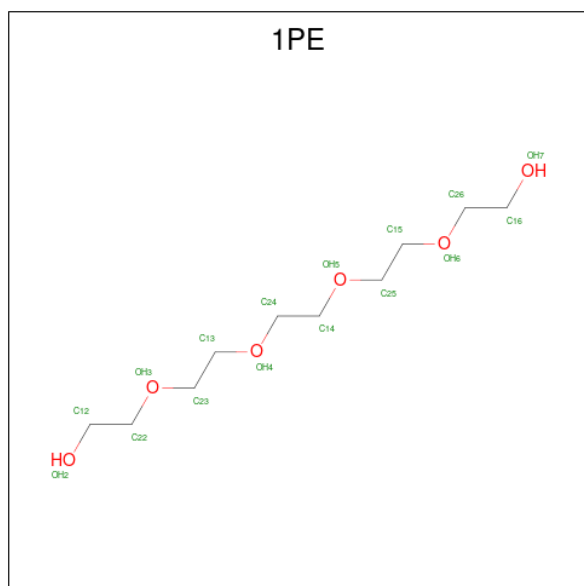
There are 5 unique types of molecules in this entry. The entry contains 7077 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Lipase A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	431	Total	C	N	O	S	0	1	0
			3272	2107	536	623	6			
1	B	431	Total	C	N	O	S	0	0	0
			3263	2102	535	620	6			

- Molecule 2 is PENTAETHYLENE GLYCOL (CCD ID: 1PE) (formula: $C_{10}H_{22}O_6$).



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

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).

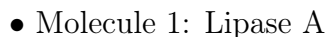


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	249	Total 249	O 249	0	0
5	B	250	Total 250	O 250	0	0

- Molecule 1: Lipase A



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	92.10Å 92.10Å 300.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.92 – 2.10 19.92 – 2.10	Depositor EDS
% Data completeness (in resolution range)	100.0 (19.92-2.10) 98.9 (19.92-2.10)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.66 (at 2.09Å)	Xtriage
Refinement program	REFMAC 5.3.0040	Depositor
R, R_{free}	0.189 , 0.234 0.196 , 0.234	Depositor DCC
R_{free} test set	3784 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	14.7	Xtriage
Anisotropy	0.095	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 49.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	7077	wwPDB-VP
Average B, all atoms (Å ²)	13.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.05% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 1PE, GOL, 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	1.21	14/3363 (0.4%)	1.12	8/4605 (0.2%)
1	B	1.24	11/3354 (0.3%)	1.10	6/4593 (0.1%)
All	All	1.22	25/6717 (0.4%)	1.11	14/9198 (0.2%)

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	320	SER	C-N	-10.09	1.22	1.33
1	A	332	ILE	C-O	-9.21	1.12	1.24
1	A	246	PRO	C-N	-8.56	1.22	1.34
1	A	120	VAL	C-O	-8.22	1.14	1.24
1	B	121	LEU	C-O	-7.88	1.14	1.24
1	B	122	ASP	C-O	-7.23	1.15	1.24
1	B	24	THR	C-O	-7.18	1.14	1.24
1	A	375	VAL	CA-CB	6.56	1.57	1.54
1	B	120	VAL	C-O	-6.46	1.16	1.24
1	B	92	VAL	CA-CB	6.10	1.62	1.54
1	A	23	THR	C-O	-6.00	1.16	1.23
1	B	405	THR	C-O	-5.94	1.16	1.24
1	A	202	PRO	C-O	-5.80	1.16	1.24
1	A	121	LEU	C-O	-5.60	1.17	1.24
1	A	401	ILE	CA-CB	5.59	1.61	1.54
1	B	184	SER	C-O	-5.48	1.19	1.24
1	B	207	VAL	CA-CB	5.39	1.60	1.54
1	B	150	ILE	CA-CB	5.36	1.61	1.54
1	A	247	ASP	C-N	-5.21	1.26	1.33
1	A	381	ILE	CA-CB	5.06	1.60	1.54
1	A	61	THR	C-O	-5.04	1.19	1.24
1	B	74	THR	CA-CB	5.04	1.61	1.53
1	A	336	ILE	N-CA	5.04	1.50	1.46
1	A	191	VAL	CA-CB	5.03	1.60	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	273	CYS	C-O	-5.02	1.16	1.24

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	247	ASP	O-C-N	-8.37	112.42	122.22
1	B	321	VAL	CA-C-N	7.75	127.81	119.90
1	B	321	VAL	C-N-CA	7.75	127.81	119.90
1	A	152	GLY	N-CA-C	7.13	121.29	112.73
1	A	367	LEU	CA-CB-CG	-6.74	92.70	116.30
1	B	367	LEU	CA-CB-CG	-6.62	93.14	116.30
1	A	414	GLY	N-CA-C	-6.46	104.55	112.68
1	A	228	GLY	CA-C-N	6.33	126.24	119.28
1	A	228	GLY	C-N-CA	6.33	126.24	119.28
1	B	336	ILE	N-CA-C	6.07	116.42	111.62
1	B	321	VAL	O-C-N	-5.87	114.41	121.10
1	A	406	THR	CA-C-N	-5.64	113.80	120.11
1	A	406	THR	C-N-CA	-5.64	113.80	120.11
1	B	60	ARG	CD-NE-CZ	5.19	131.66	124.40

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	3272	0	3209	40	0
1	B	3263	0	3204	29	0
2	A	16	0	22	1	0
2	B	16	0	22	3	0
3	A	6	0	8	8	0
4	B	5	0	0	0	0
5	A	249	0	0	4	0
5	B	250	0	0	6	0
All	All	7077	0	6465	69	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 5.

All (69) 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:B:127:ILE:HG13	5:B:631:HOH:O	1.33	1.25
1:A:247:ASP:HB2	3:A:443:GOL:O1	1.67	0.92
1:A:304:ILE:HD11	3:A:443:GOL:H11	1.51	0.91
1:B:238:VAL:HG22	2:B:442:1PE:H222	1.59	0.83
1:A:198:GLU:HG3	5:A:507:HOH:O	1.78	0.83
1:A:245:HIS:HD1	3:A:443:GOL:HO3	1.20	0.83
1:B:127:ILE:CD1	5:B:631:HOH:O	2.24	0.80
1:A:245:HIS:ND1	3:A:443:GOL:O3	2.16	0.71
1:B:330:HIS:HD2	5:B:447:HOH:O	1.73	0.70
1:A:247:ASP:HB2	3:A:443:GOL:C1	2.22	0.69
1:B:127:ILE:HG23	1:B:137:VAL:HG11	1.75	0.69
1:B:367:LEU:HD22	1:B:429:SER:CB	2.23	0.68
1:A:83:SER:HA	1:A:84:PRO:C	2.20	0.67
1:A:365:GLU:OE1	5:A:453:HOH:O	2.13	0.67
1:B:238:VAL:HG22	2:B:442:1PE:C22	2.25	0.66
1:B:33:LYS:HA	1:B:60:ARG:HD2	1.83	0.60
1:B:65:GLN:OE1	5:B:673:HOH:O	2.15	0.59
1:B:127:ILE:CG1	5:B:631:HOH:O	2.05	0.57
1:B:281:TYR:N	1:B:282:PRO:HD3	2.20	0.57
1:A:116:LYS:HE2	5:A:602:HOH:O	2.04	0.56
1:B:259:LYS:HE2	1:B:289:LEU:HD23	1.88	0.56
1:A:155:GLU:OE1	1:A:188:HIS:ND1	2.39	0.55
1:A:210:SER:HB2	1:A:381:ILE:HD11	1.88	0.55
1:A:304:ILE:HD11	3:A:443:GOL:C1	2.32	0.54
1:A:367:LEU:HD13	1:A:433:LYS:HB2	1.89	0.54
1:A:159:ILE:HD11	1:A:189:ALA:HB1	1.88	0.54
1:B:367:LEU:HD13	1:B:433:LYS:HB2	1.89	0.53
1:B:222:PHE:HB2	2:B:442:1PE:H131	1.91	0.53
1:A:222:PHE:HB2	2:A:442:1PE:H231	1.92	0.52
1:A:428:GLN:NE2	1:A:432:GLY:C	2.68	0.51
1:A:242:SER:HB3	1:A:249:GLU:OE1	2.12	0.50
1:B:159:ILE:HD11	1:B:189:ALA:HB1	1.94	0.49
1:B:300:PRO:O	1:B:304:ILE:HG12	2.14	0.48
1:A:149:PHE:O	1:A:150:ILE:HG22	2.14	0.47
1:A:235:LEU:HD21	1:A:263:THR:HG22	1.96	0.47
1:A:245:HIS:HB3	3:A:443:GOL:H2	1.97	0.47
1:B:281:TYR:N	1:B:282:PRO:CD	2.77	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:292:ASP:OD2	1:A:298:GLU:OE1	2.33	0.47
1:A:274:LEU:C	1:A:274:LEU:HD23	2.39	0.46
1:A:181:GLU:HA	1:A:210:SER:O	2.15	0.46
1:A:122:ASP:OD2	1:A:370:GLU:OE2	2.34	0.46
1:A:217:SER:HB3	1:A:220:ASP:HB2	1.97	0.45
1:B:365:GLU:H	1:B:368:THR:HG1	1.62	0.45
1:B:432:GLY:C	5:B:684:HOH:O	2.59	0.45
1:A:115:ASN:HD22	1:A:115:ASN:HA	1.65	0.45
1:A:219:LYS:HB2	1:A:305:LEU:HG	1.98	0.45
1:A:290:VAL:HG21	1:A:295:LEU:HD22	1.99	0.45
1:A:326:ARG:HD3	5:A:445:HOH:O	2.16	0.44
1:B:259:LYS:HE2	1:B:289:LEU:CD2	2.48	0.44
1:A:157:MET:HE2	1:A:200:TYR:CD1	2.53	0.44
1:B:295:LEU:HG	1:B:301:ILE:HD12	1.99	0.44
1:A:428:GLN:HE22	1:A:432:GLY:C	2.26	0.43
1:A:248:MET:H	3:A:443:GOL:H12	1.83	0.43
1:B:212:GLY:HA2	1:B:329:TRP:O	2.18	0.43
1:A:13:LEU:HB3	1:A:14:PRO:HD2	2.01	0.43
1:B:60:ARG:HD3	1:B:61:THR:O	2.19	0.43
1:B:95:ASP:HB2	1:B:274:LEU:HD12	2.00	0.42
1:B:181:GLU:HA	1:B:210:SER:O	2.18	0.42
1:A:54:SER:HB3	1:A:76:TRP:CD1	2.55	0.42
1:A:367:LEU:HD22	1:A:429:SER:CB	2.50	0.42
1:A:428:GLN:NE2	1:A:432:GLY:O	2.51	0.42
1:A:127:ILE:HG23	1:A:137:VAL:HG11	2.03	0.41
1:B:292:ASP:HB3	1:B:295:LEU:HD13	2.02	0.41
1:A:326:ARG:NH2	1:A:345:TYR:OH	2.54	0.41
1:B:54:SER:HB3	1:B:76:TRP:CD1	2.55	0.41
1:B:367:LEU:HD22	1:B:429:SER:OG	2.21	0.41
1:B:167:LYS:HB3	1:B:167:LYS:HE2	1.64	0.41
1:A:376:PRO:HA	1:A:406:THR:HG21	2.04	0.40
1:A:177:LYS:HG2	1:A:207:VAL:HG11	2.03	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	430/462 (93%)	413 (96%)	16 (4%)	1 (0%)	43	44
1	B	429/462 (93%)	416 (97%)	11 (3%)	2 (0%)	24	22
All	All	859/924 (93%)	829 (96%)	27 (3%)	3 (0%)	36	36

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	150	ILE
1	B	150	ILE
1	B	281	TYR

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	347/368 (94%)	337 (97%)	10 (3%)	37	42
1	B	346/368 (94%)	335 (97%)	11 (3%)	34	38
All	All	693/736 (94%)	672 (97%)	21 (3%)	36	41

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	60	ARG
1	A	115	ASN
1	A	150	ILE
1	A	210	SER
1	A	259	LYS
1	A	295	LEU
1	A	296	LEU
1	A	305	LEU
1	A	411	GLN

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Mol	Chain	Res	Type
1	A	415	SER
1	B	38	GLN
1	B	60	ARG
1	B	150	ILE
1	B	253	GLU
1	B	264	LEU
1	B	265	LYS
1	B	306	LYS
1	B	382	LYS
1	B	383	GLN
1	B	398	ILE
1	B	401	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	ASN
1	A	115	ASN
1	A	297	ASN
1	A	330	HIS
1	B	38	GLN
1	B	65	GLN
1	B	132	GLN
1	B	330	HIS
1	B	383	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

4 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	442	-	15,15,15	0.49	0	14,14,14	0.61	0
2	1PE	B	442	-	15,15,15	0.61	0	14,14,14	0.51	0
3	GOL	A	443	-	5,5,5	0.73	0	5,5,5	0.99	0
4	SO4	B	443	-	4,4,4	0.27	0	6,6,6	0.78	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	442	-	-	6/13/13/13	-
2	1PE	B	442	-	-	5/13/13/13	-
3	GOL	A	443	-	-	3/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	442	1PE	OH7-C16-C26-OH6
2	B	442	1PE	OH4-C13-C23-OH3
3	A	443	GOL	O1-C1-C2-C3
3	A	443	GOL	C1-C2-C3-O3
2	A	442	1PE	OH2-C12-C22-OH3
2	B	442	1PE	OH7-C16-C26-OH6
3	A	443	GOL	O2-C2-C3-O3
2	A	442	1PE	C12-C22-OH3-C23

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Mol	Chain	Res	Type	Atoms
2	A	442	1PE	C13-C23-OH3-C22
2	A	442	1PE	OH5-C14-C24-OH4
2	B	442	1PE	C13-C23-OH3-C22
2	B	442	1PE	C23-C13-OH4-C24
2	B	442	1PE	OH5-C14-C24-OH4
2	A	442	1PE	OH4-C13-C23-OH3

There are no ring outliers.

3 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	442	1PE	1	0
2	B	442	1PE	3	0
3	A	443	GOL	8	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	431/462 (93%)	-0.41	3 (0%)	84 86	5, 12, 24, 40	1 (0%)
1	B	431/462 (93%)	-0.46	3 (0%)	84 86	6, 12, 22, 39	0
All	All	862/924 (93%)	-0.44	6 (0%)	84 86	5, 12, 23, 40	1 (0%)

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	198	GLU	3.6
1	A	247	ASP	3.0
1	B	441	PRO	2.7
1	A	432	GLY	2.3
1	B	171	ASN	2.1
1	B	295	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 oligosaccharides 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($\text{\AA}^2$)	Q<0.9
3	GOL	A	443	6/6	0.73	0.18	29,29,33,34	0
2	1PE	A	442	16/16	0.92	0.08	17,23,29,30	0
2	1PE	B	442	16/16	0.93	0.09	14,22,36,38	0
4	SO4	B	443	5/5	0.94	0.13	32,32,36,37	0

6.5 Other polymers [i](#)

There are no such residues in this entry.