



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 21, 2020 – 11:57 AM BST

PDB ID : 4OKH
Title : Crystal structure of calpain-3 penta-EF-hand domain
Authors : Karunan Partha, S.; Ravulapalli, R.; Campbell, R.L.; Allingham, J.S.; Davies, P.L.
Deposited on : 2014-01-22
Resolution : 2.45 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
buster-report : 1.1.7 (2018)
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) : 2.13.1

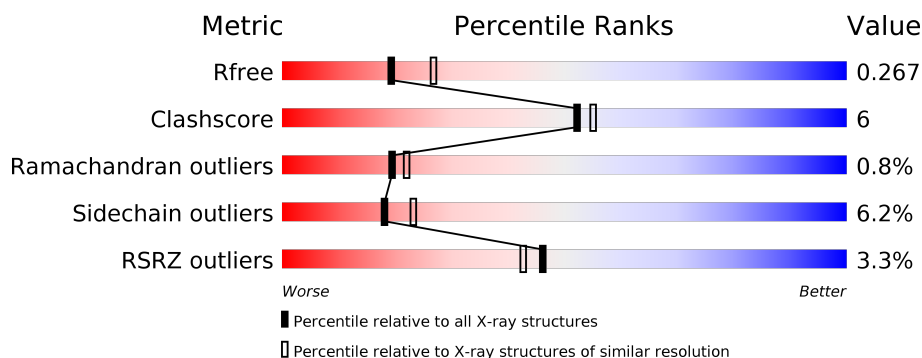
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 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	1544 (2.48-2.44)
Clashscore	141614	1613 (2.48-2.44)
Ramachandran outliers	138981	1598 (2.48-2.44)
Sidechain outliers	138945	1598 (2.48-2.44)
RSRZ outliers	127900	1523 (2.48-2.44)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	189	<div> <div>2%</div> <div> <div></div> <div>79%</div> <div>10%</div> <div>••</div> <div>10%</div> </div> </div>
1	B	189	<div> <div>2%</div> <div> <div></div> <div>80%</div> <div>9%</div> <div>•</div> <div>10%</div> </div> </div>
1	C	189	<div> <div>4%</div> <div> <div></div> <div>59%</div> <div>14%</div> <div>5%</div> <div>22%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4097 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 Calpain-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	171	Total	C	N	O	S	0	0	0
			1409	895	245	257	12			
1	B	170	Total	C	N	O	S	0	0	0
			1394	884	237	261	12			
1	C	147	Total	C	N	O	S	0	0	0
			1202	765	204	221	12			

There are 27 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	641	MET	-	INITIATING METHIONINE	UNP P20807
A	822	LEU	-	EXPRESSION TAG	UNP P20807
A	823	GLU	-	EXPRESSION TAG	UNP P20807
A	824	HIS	-	EXPRESSION TAG	UNP P20807
A	825	HIS	-	EXPRESSION TAG	UNP P20807
A	826	HIS	-	EXPRESSION TAG	UNP P20807
A	827	HIS	-	EXPRESSION TAG	UNP P20807
A	828	HIS	-	EXPRESSION TAG	UNP P20807
A	829	HIS	-	EXPRESSION TAG	UNP P20807
B	641	MET	-	INITIATING METHIONINE	UNP P20807
B	822	LEU	-	EXPRESSION TAG	UNP P20807
B	823	GLU	-	EXPRESSION TAG	UNP P20807
B	824	HIS	-	EXPRESSION TAG	UNP P20807
B	825	HIS	-	EXPRESSION TAG	UNP P20807
B	826	HIS	-	EXPRESSION TAG	UNP P20807
B	827	HIS	-	EXPRESSION TAG	UNP P20807
B	828	HIS	-	EXPRESSION TAG	UNP P20807
B	829	HIS	-	EXPRESSION TAG	UNP P20807
C	641	MET	-	INITIATING METHIONINE	UNP P20807
C	822	LEU	-	EXPRESSION TAG	UNP P20807
C	823	GLU	-	EXPRESSION TAG	UNP P20807
C	824	HIS	-	EXPRESSION TAG	UNP P20807
C	825	HIS	-	EXPRESSION TAG	UNP P20807

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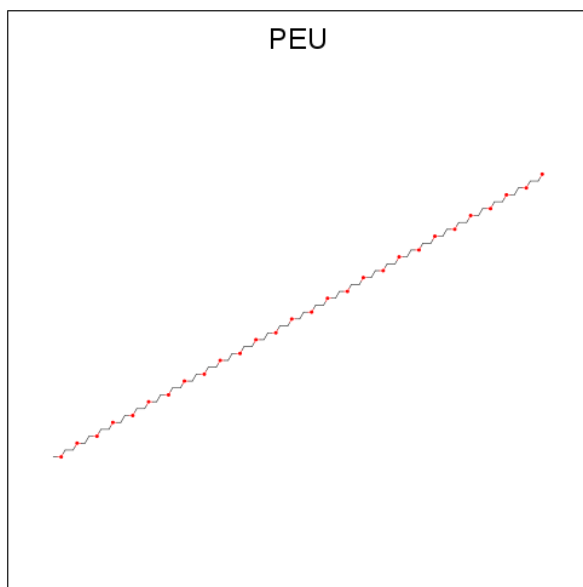
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Chain	Residue	Modelled	Actual	Comment	Reference
C	826	HIS	-	EXPRESSION TAG	UNP P20807
C	827	HIS	-	EXPRESSION TAG	UNP P20807
C	828	HIS	-	EXPRESSION TAG	UNP P20807
C	829	HIS	-	EXPRESSION TAG	UNP P20807

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	4	Total Ca 4 4	0	0
2	A	4	Total Ca 4 4	0	0
2	C	3	Total Ca 3 3	0	0

- Molecule 3 is 2,5,8,11,14,17,20,23,26,29,32,35,38,41,44,47,50,53,56,59,62,65,68,71,74,77,80-HEPTACOSAOXADOCTACONTAN-82-OL (three-letter code: PEU) (formula: C₅₅H₁₁₂O₂₈).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 22 14 8	0	0
3	B	1	Total C O 13 8 5	0	0

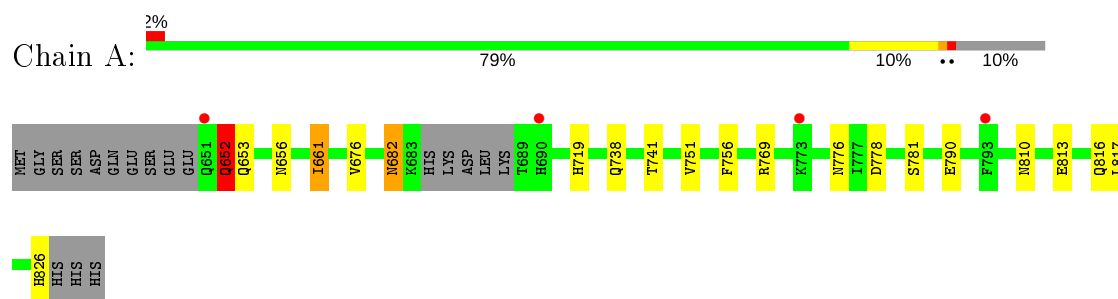
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	17	Total 17	O 17	0	0
4	B	17	Total 17	O 17	0	0
4	C	12	Total 12	O 12	0	0

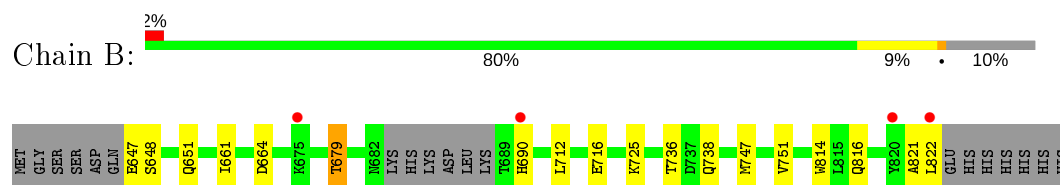
3 Residue-property plots [i](#)

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.

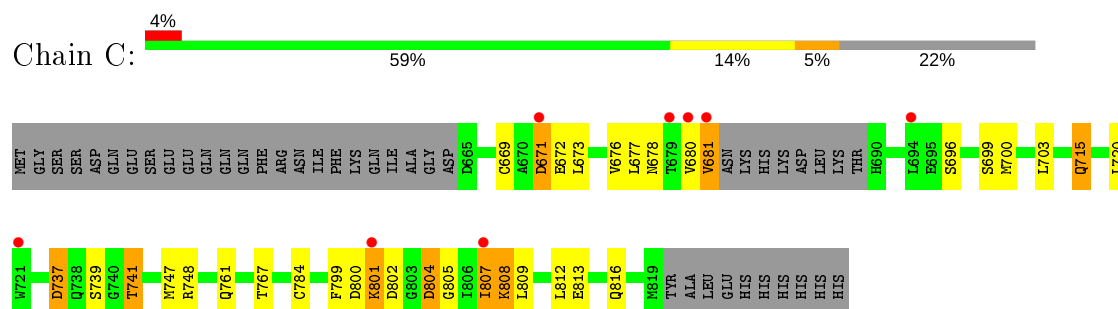
• Molecule 1: Calpain-3



• Molecule 1: Calpain-3



• Molecule 1: Calpain-3



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	107.00 Å 107.00 Å 96.71 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.85 – 2.45 47.85 – 2.45	Depositor EDS
% Data completeness (in resolution range)	100.0 (47.85-2.45) 100.0 (47.85-2.45)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.39 (at 2.45 Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.198 , 0.269 0.201 , 0.267	Depositor DCC
R_{free} test set	1098 reflections (5.16%)	wwPDB-VP
Wilson B-factor (Å ²)	46.9	Xtriage
Anisotropy	0.127	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 51.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4097	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.28% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PEU, CA

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	0.65	0/1440	0.76	0/1935
1	B	0.61	0/1422	0.76	0/1911
1	C	0.57	0/1227	0.75	0/1649
All	All	0.61	0/4089	0.76	0/5495

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	B	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	679	THR	Peptide

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	1409	0	1352	12	0
1	B	1394	0	1336	10	0
1	C	1202	0	1159	23	0
2	A	4	0	0	0	0
2	B	4	0	0	0	0
2	C	3	0	0	0	0
3	A	22	0	29	3	0
3	B	13	0	17	0	0
4	A	17	0	0	1	0
4	B	17	0	0	0	0
4	C	12	0	0	2	0
All	All	4097	0	3893	44	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 6.

All (44) 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:C:739:SER:OG	1:C:741:THR:HG22	1.75	0.85
1:A:653:GLN:HA	1:A:656:ASN:HD22	1.59	0.67
1:C:748:ARG:HD2	4:C:1003:HOH:O	1.93	0.67
1:B:661:ILE:C	1:B:661:ILE:HD12	2.17	0.65
1:A:719:HIS:HB2	4:A:1006:HOH:O	1.97	0.64
1:C:802:ASP:OD2	1:C:804:ASP:HB3	2.00	0.62
1:B:821:ALA:C	1:B:822:LEU:HD12	2.22	0.59
1:A:816:GLN:HB3	3:A:905:PEU:HCW2	1.84	0.58
1:B:647:GLU:OE1	1:B:647:GLU:N	2.38	0.57
1:B:647:GLU:HG2	1:B:648:SER:H	1.70	0.57
1:A:682:ASN:OD1	1:A:682:ASN:C	2.43	0.56
1:C:673:LEU:O	1:C:677:LEU:HB2	2.09	0.52
1:C:801:LYS:HA	1:C:801:LYS:HE2	1.92	0.51
1:C:812:LEU:O	1:C:816:GLN:HG3	2.10	0.51
1:C:799:PHE:HB3	1:C:807:ILE:HD11	1.93	0.50
1:C:680:VAL:O	1:C:681:VAL:HG13	2.12	0.50
1:B:725:LYS:HA	1:B:725:LYS:HE2	1.94	0.49
1:C:739:SER:HG	1:C:741:THR:HG22	1.77	0.49
1:A:741:THR:HB	1:A:776:ASN:HB3	1.95	0.48
1:C:677:LEU:HD21	1:C:700:MET:HE1	1.96	0.47
1:A:813:GLU:HG3	3:A:905:PEU:CCK	2.45	0.47
1:C:807:ILE:HA	1:C:808:LYS:HD3	1.96	0.47
1:C:715:GLN:HB3	1:C:715:GLN:HE21	1.60	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:747:MET:CG	1:C:767:THR:HG22	2.46	0.46
1:C:747:MET:HG2	1:C:767:THR:HG22	1.99	0.45
1:C:800:ASP:OD2	1:C:805:GLY:N	2.46	0.45
1:C:813:GLU:HA	1:C:816:GLN:HG3	1.98	0.44
1:C:802:ASP:OD2	1:C:804:ASP:CB	2.66	0.43
1:A:817:LEU:HG	1:B:814:TRP:CH2	2.54	0.43
1:B:747:MET:O	1:B:751:VAL:HG23	2.17	0.43
1:A:661:ILE:HG12	1:A:676:VAL:HB	2.01	0.42
1:C:671:ASP:OD1	1:C:671:ASP:N	2.52	0.42
1:A:810:ASN:HD21	3:A:905:PEU:CCK	2.33	0.42
1:B:712:LEU:HA	1:B:716:GLU:OE2	2.20	0.42
1:C:696:SER:HB3	1:C:784:CYS:SG	2.59	0.42
1:C:741:THR:HG23	4:C:1004:HOH:O	2.19	0.42
1:A:751:VAL:HG13	1:A:756:PHE:HB2	2.01	0.42
1:A:652:GLN:N	1:A:652:GLN:OE1	2.52	0.41
1:B:821:ALA:O	1:B:822:LEU:HD12	2.20	0.41
1:B:679:THR:O	1:B:679:THR:HG22	2.20	0.41
1:C:672:GLU:O	1:C:676:VAL:HG12	2.20	0.41
1:C:804:ASP:O	1:C:804:ASP:OD1	2.39	0.41
1:A:778:ASP:OD1	1:A:781:SER:HB2	2.20	0.41
1:C:673:LEU:HD21	1:C:700:MET:HE2	2.04	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	167/189 (88%)	163 (98%)	2 (1%)	2 (1%)	13	12
1	B	166/189 (88%)	159 (96%)	6 (4%)	1 (1%)	25	29
1	C	143/189 (76%)	140 (98%)	2 (1%)	1 (1%)	22	25
All	All	476/567 (84%)	462 (97%)	10 (2%)	4 (1%)	19	22

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	737	ASP
1	A	738	GLN
1	A	652	GLN
1	B	690	HIS

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	152/169 (90%)	146 (96%)	6 (4%)	32	42
1	B	151/169 (89%)	146 (97%)	5 (3%)	38	49
1	C	131/169 (78%)	115 (88%)	16 (12%)	5	4
All	All	434/507 (86%)	407 (94%)	27 (6%)	18	23

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	652	GLN
1	A	661	ILE
1	A	682	ASN
1	A	769	ARG
1	A	790	GLU
1	A	826	HIS
1	B	651	GLN
1	B	664	ASP
1	B	736	THR
1	B	738	GLN
1	B	816	GLN
1	C	669	CYS
1	C	671	ASP
1	C	678	ASN
1	C	681	VAL
1	C	699	SER
1	C	703	LEU

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Mol	Chain	Res	Type
1	C	715	GLN
1	C	720	LEU
1	C	737	ASP
1	C	741	THR
1	C	761	GLN
1	C	801	LYS
1	C	804	ASP
1	C	807	ILE
1	C	808	LYS
1	C	809	LEU

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

Mol	Chain	Res	Type
1	A	651	GLN
1	A	656	ASN
1	A	719	HIS
1	B	690	HIS
1	C	810	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](#)

Of 13 ligands modelled in this entry, 11 are monoatomic - leaving 2 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	PEU	A	905	-	21,21,82	0.86	0	20,20,81	0.97	1 (5%)
3	PEU	B	905	-	12,12,82	0.59	0	11,11,81	0.35	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
3	PEU	A	905	-	-	10/19/19/80	-
3	PEU	B	905	-	-	5/10/10/80	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	905	PEU	OCP-CCQ-CCR	2.69	122.52	110.39

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	905	PEU	OCM-CCN-CCO-OCP
3	A	905	PEU	OCY-CCZ-CDA-ODB
3	B	905	PEU	OCY-CCZ-CDA-ODB
3	B	905	PEU	OCS-CCT-CCU-OCV
3	A	905	PEU	ODB-CDC-CDD-ODE
3	B	905	PEU	ODB-CDC-CDD-ODE
3	A	905	PEU	OCS-CCT-CCU-OCV
3	B	905	PEU	OCV-CCW-CCX-OCY
3	A	905	PEU	CCW-CCX-OCY-CCZ
3	B	905	PEU	CCX-CCW-OCV-CCU
3	A	905	PEU	CCK-CCL-OCM-CCN
3	A	905	PEU	CDD-CDC-ODB-CDA
3	A	905	PEU	OCJ-CCK-CCL-OCM
3	A	905	PEU	CCR-CCQ-OCP-CCO

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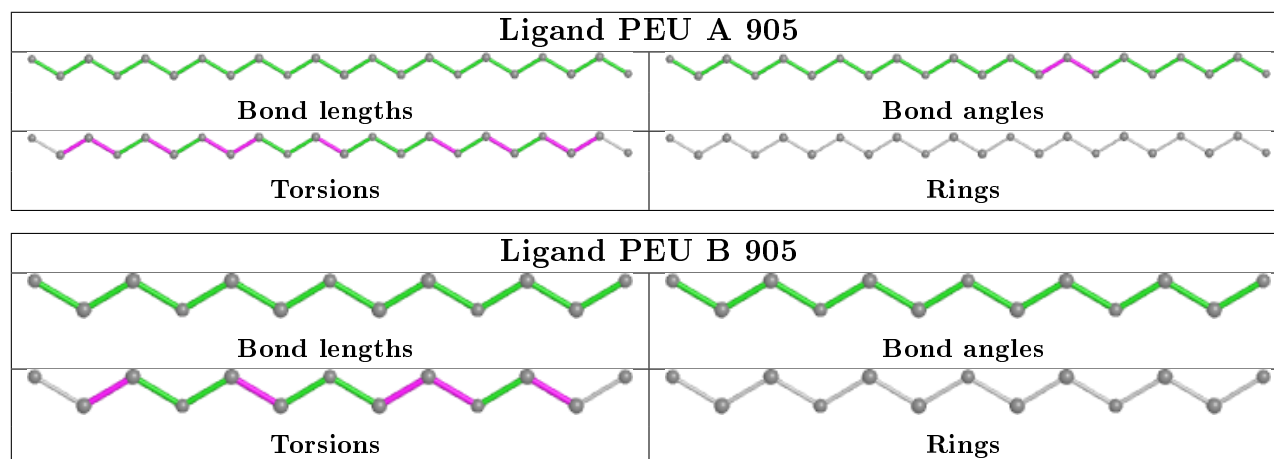
Mol	Chain	Res	Type	Atoms
3	A	905	PEU	OCV-CCW-CCX-OCY

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	905	PEU	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	171/189 (90%)	-0.11	4 (2%) 60 56	32, 49, 88, 110	0
1	B	170/189 (89%)	-0.10	4 (2%) 59 54	34, 53, 84, 115	0
1	C	147/189 (77%)	0.27	8 (5%) 25 23	36, 70, 113, 137	0
All	All	488/567 (86%)	0.01	16 (3%) 46 43	32, 55, 101, 137	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	820	TYR	4.7
1	C	680	VAL	3.9
1	C	679	THR	3.6
1	C	681	VAL	3.3
1	B	822	LEU	3.2
1	C	801	LYS	3.0
1	A	651	GLN	2.7
1	A	690	HIS	2.7
1	C	694	LEU	2.7
1	C	807	ILE	2.3
1	A	793	PHE	2.2
1	B	690	HIS	2.2
1	C	721	TRP	2.1
1	B	675	LYS	2.1
1	A	773	LYS	2.1
1	C	671	ASP	2.0

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 [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

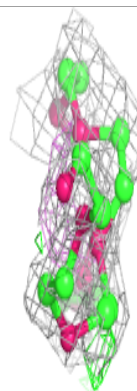
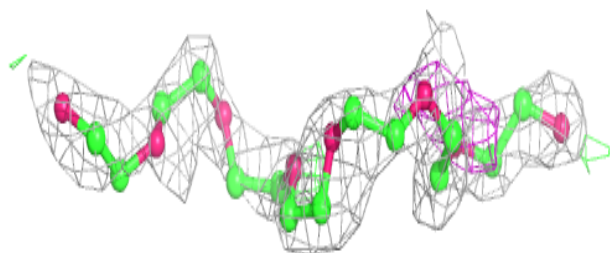
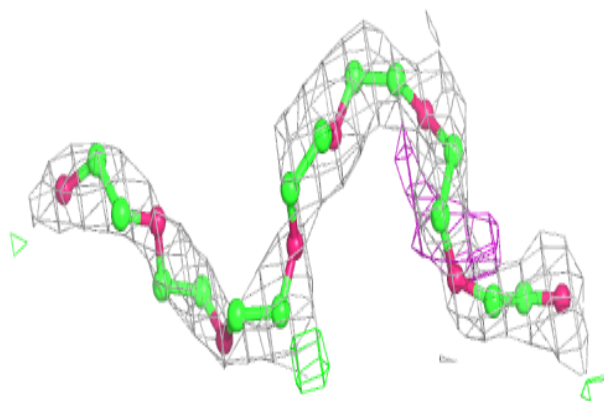
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	PEU	A	905	22/83	0.78	0.31	45,71,77,79	0
3	PEU	B	905	13/83	0.82	0.24	57,93,107,109	0
2	CA	C	903	1/1	0.87	0.07	97,97,97,97	0
2	CA	B	903	1/1	0.90	0.07	86,86,86,86	0
2	CA	C	901	1/1	0.92	0.05	80,80,80,80	0
2	CA	A	903	1/1	0.93	0.09	83,83,83,83	0
2	CA	A	901	1/1	0.97	0.11	53,53,53,53	0
2	CA	C	902	1/1	0.97	0.08	76,76,76,76	0
2	CA	B	902	1/1	0.97	0.14	48,48,48,48	0
2	CA	B	901	1/1	0.97	0.10	58,58,58,58	0
2	CA	A	904	1/1	0.98	0.09	38,38,38,38	0
2	CA	B	904	1/1	0.98	0.14	51,51,51,51	0
2	CA	A	902	1/1	0.99	0.13	49,49,49,49	0

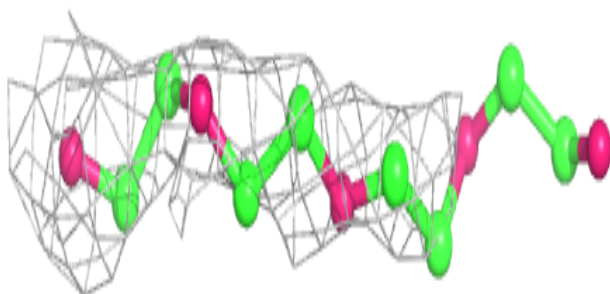
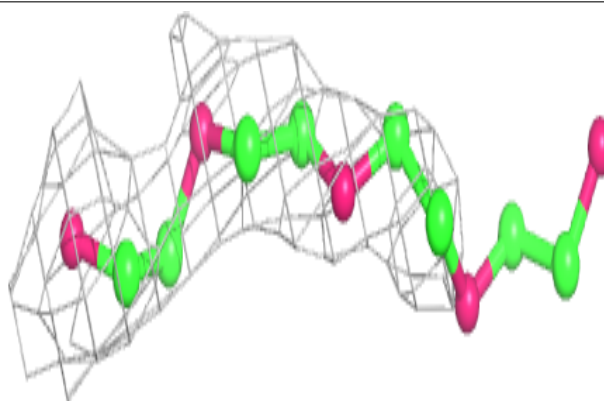
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around PEU A 905:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around PEU B 905:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.