



Full wwPDB X-ray Structure Validation Report ⓘ

May 5, 2022 – 12:03 PM EDT

PDB ID : 7URY
Title : Tetradecameric hub domain of CaMKII beta
Authors : Ozden, C.; Saha, S.; Samkutty, A.; Stratton, M.M.; Garman, S.C.; Perry, S.L.
Deposited on : 2022-04-22
Resolution : 2.64 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 : 2.28.1
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) : 2.28.1

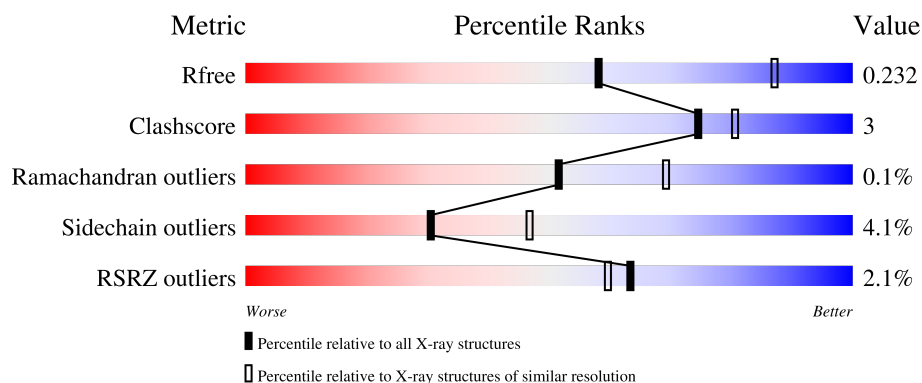
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.64 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	1426 (2.66-2.62)
Clashscore	141614	1472 (2.66-2.62)
Ramachandran outliers	138981	1446 (2.66-2.62)
Sidechain outliers	138945	1446 (2.66-2.62)
RSRZ outliers	127900	1408 (2.66-2.62)

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	137	<div> <div>81%</div> <div>13%</div> <div>..</div> </div>
1	B	137	<div> <div>82%</div> <div>9%</div> <div>9%</div> </div>
1	C	137	<div> <div>3%</div> <div>90%</div> <div>5%</div> <div>..</div> </div>
1	D	137	<div> <div>3%</div> <div>84%</div> <div>9%</div> <div>6%</div> </div>
1	E	137	<div> <div>85%</div> <div>12%</div> <div>..</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	137	<div> <div>4%</div> <div>87%</div> <div>6%</div> <div>6%</div> </div>
1	G	137	<div> <div>%</div> <div>83%</div> <div>7%</div> <div>7%</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7072 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 Calcium/calmodulin-dependent protein kinase type II subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	G	127	Total	C	N	O	S	0	0	0
			979	621	169	184	5			
1	A	131	Total	C	N	O	S	0	0	0
			1021	647	178	191	5			
1	B	125	Total	C	N	O	S	0	0	0
			973	614	170	184	5			
1	C	131	Total	C	N	O	S	0	0	0
			997	635	170	188	4			
1	D	129	Total	C	N	O	S	0	0	0
			1012	640	178	189	5			
1	E	134	Total	C	N	O	S	0	0	0
			1049	665	181	198	5			
1	F	129	Total	C	N	O	S	0	0	0
			1005	637	174	189	5			

There are 28 discrepancies between the modelled and reference sequences:

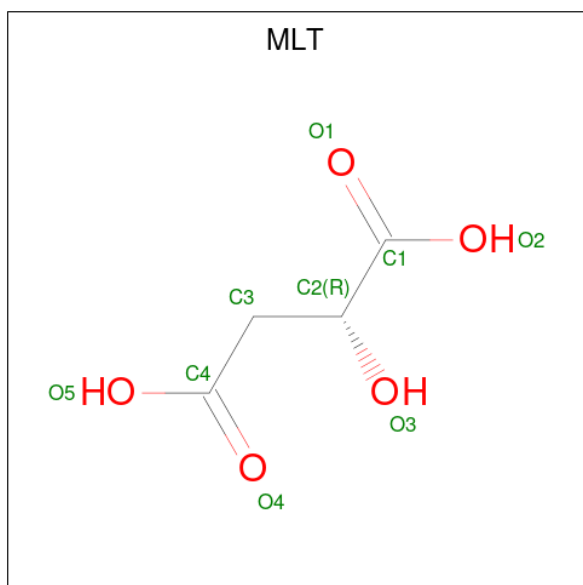
Chain	Residue	Modelled	Actual	Comment	Reference
G	530	GLY	-	expression tag	UNP Q13554
G	531	PRO	-	expression tag	UNP Q13554
G	532	HIS	-	expression tag	UNP Q13554
G	533	MET	-	expression tag	UNP Q13554
A	530	GLY	-	expression tag	UNP Q13554
A	531	PRO	-	expression tag	UNP Q13554
A	532	HIS	-	expression tag	UNP Q13554
A	533	MET	-	expression tag	UNP Q13554
B	530	GLY	-	expression tag	UNP Q13554
B	531	PRO	-	expression tag	UNP Q13554
B	532	HIS	-	expression tag	UNP Q13554
B	533	MET	-	expression tag	UNP Q13554
C	530	GLY	-	expression tag	UNP Q13554
C	531	PRO	-	expression tag	UNP Q13554

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Chain	Residue	Modelled	Actual	Comment	Reference
C	532	HIS	-	expression tag	UNP Q13554
C	533	MET	-	expression tag	UNP Q13554
D	530	GLY	-	expression tag	UNP Q13554
D	531	PRO	-	expression tag	UNP Q13554
D	532	HIS	-	expression tag	UNP Q13554
D	533	MET	-	expression tag	UNP Q13554
E	530	GLY	-	expression tag	UNP Q13554
E	531	PRO	-	expression tag	UNP Q13554
E	532	HIS	-	expression tag	UNP Q13554
E	533	MET	-	expression tag	UNP Q13554
F	530	GLY	-	expression tag	UNP Q13554
F	531	PRO	-	expression tag	UNP Q13554
F	532	HIS	-	expression tag	UNP Q13554
F	533	MET	-	expression tag	UNP Q13554

- Molecule 2 is D-MALATE (three-letter code: MLT) (formula: $C_4H_6O_5$).

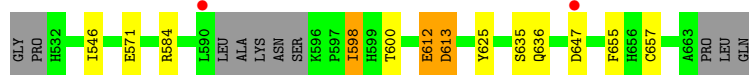
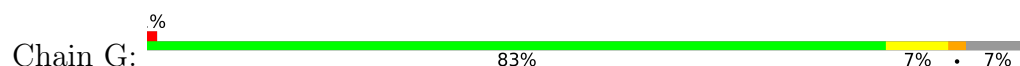


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			9	4	5		
2	D	1	Total	C	O	0	0
			9	4	5		
2	E	1	Total	C	O	0	0
			9	4	5		
2	F	1	Total	C	O	0	0
			9	4	5		

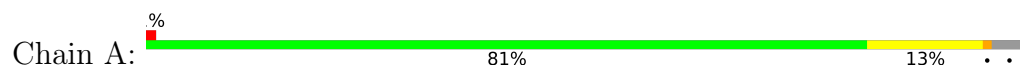
3 Residue-property plots [i](#)

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.

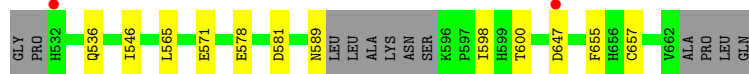
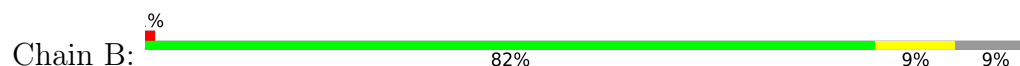
- Molecule 1: Calcium/calmodulin-dependent protein kinase type II subunit beta



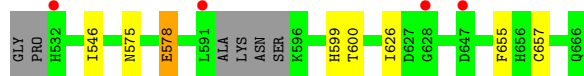
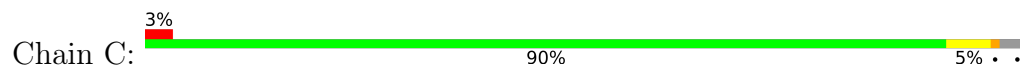
- Molecule 1: Calcium/calmodulin-dependent protein kinase type II subunit beta



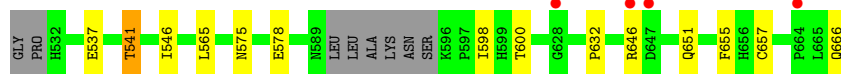
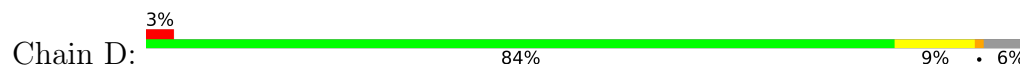
- Molecule 1: Calcium/calmodulin-dependent protein kinase type II subunit beta



- Molecule 1: Calcium/calmodulin-dependent protein kinase type II subunit beta



- Molecule 1: Calcium/calmodulin-dependent protein kinase type II subunit beta



- Molecule 1: Calcium/calmodulin-dependent protein kinase type II subunit beta

Chain E:

85%

12%

..



● Molecule 1: Calcium/calmodulin-dependent protein kinase type II subunit beta

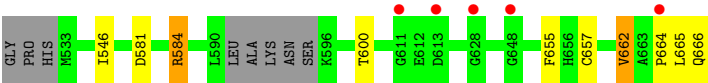
Chain F:

4%

87%

6%

6%



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	105.71Å 182.79Å 110.52Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.23 – 2.64 38.20 – 2.64	Depositor EDS
% Data completeness (in resolution range)	88.7 (38.23-2.64) 88.8 (38.20-2.64)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.92 (at 2.65Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.204 , 0.235 0.205 , 0.232	Depositor DCC
R_{free} test set	1452 reflections (5.14%)	wwPDB-VP
Wilson B-factor (Å ²)	54.7	Xtriage
Anisotropy	0.074	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 42.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.013 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.021 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7072	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.90% 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: MLT

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	0.64	0/1048	0.77	0/1427
1	B	0.65	0/999	0.73	0/1361
1	C	0.65	0/1024	0.77	0/1400
1	D	0.65	0/1038	0.76	0/1411
1	E	0.64	0/1076	0.76	1/1464 (0.1%)
1	F	0.64	0/1031	0.74	0/1404
1	G	0.65	0/1005	0.73	0/1370
All	All	0.64	0/7221	0.75	1/9837 (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	C	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	666	GLN	CA-C-O	5.59	131.83	120.10

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	578	GLU	Peptide

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	1021	0	937	12	0
1	B	973	0	878	5	0
1	C	997	0	896	5	0
1	D	1012	0	939	7	0
1	E	1049	0	979	7	0
1	F	1005	0	926	7	0
1	G	979	0	885	5	0
2	A	9	0	4	0	0
2	D	9	0	4	0	0
2	E	9	0	4	0	0
2	F	9	0	4	0	0
All	All	7072	0	6456	43	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 3.

All (43) 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:D:537:GLU:O	1:D:541:THR:HG22	1.69	0.93
1:E:591:LEU:H	1:E:594:ASN:HD22	1.44	0.65
1:A:597:PRO:HD2	1:A:626:ILE:HG12	1.80	0.62
1:A:623:THR:OG1	1:A:662:VAL:HG21	1.99	0.62
1:B:565:LEU:O	1:B:578:GLU:O	2.18	0.61
1:E:565:LEU:O	1:E:578:GLU:O	2.18	0.61
1:D:565:LEU:O	1:D:578:GLU:O	2.18	0.61
1:A:546:ILE:HG21	1:A:600:THR:HG21	1.83	0.60
1:E:546:ILE:HG21	1:E:600:THR:HG21	1.84	0.60
1:F:546:ILE:HG21	1:F:600:THR:HG21	1.83	0.59
1:B:546:ILE:HG21	1:B:600:THR:HG21	1.85	0.59
1:G:546:ILE:HG21	1:G:600:THR:HG21	1.87	0.56
1:C:546:ILE:HG21	1:C:600:THR:HG21	1.87	0.55
1:F:581:ASP:OD1	1:F:584:ARG:NH1	2.39	0.55
1:G:598:ILE:HG23	1:G:625:TYR:HB3	1.89	0.55
1:D:546:ILE:HG21	1:D:600:THR:HG21	1.87	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:554:PHE:CZ	1:A:584:ARG:HG2	2.43	0.54
1:A:623:THR:OG1	1:A:662:VAL:CG2	2.56	0.53
1:A:601:THR:HG22	1:A:603:LEU:HD13	1.92	0.51
1:A:636:GLN:NE2	1:C:575:ASN:HD21	2.09	0.50
1:G:636:GLN:NE2	1:D:575:ASN:HD21	2.09	0.50
1:G:612:GLU:O	1:G:613:ASP:CB	2.60	0.50
1:D:632:PRO:HG3	1:F:665:LEU:O	2.12	0.49
1:A:597:PRO:HD2	1:A:626:ILE:CG1	2.42	0.48
1:D:646:ARG:HH22	1:D:651:GLN:HE21	1.63	0.47
1:E:571:GLU:OE1	1:E:661:PRO:HB3	2.15	0.47
1:A:655:PHE:CZ	1:A:657:CYS:HB2	2.50	0.46
1:A:597:PRO:HD2	1:A:626:ILE:CD1	2.46	0.46
1:B:655:PHE:CZ	1:B:657:CYS:HB2	2.51	0.46
1:G:655:PHE:CZ	1:G:657:CYS:HB2	2.52	0.45
1:D:655:PHE:CZ	1:D:657:CYS:HB2	2.52	0.45
1:E:655:PHE:CZ	1:E:657:CYS:HB2	2.51	0.45
1:F:655:PHE:CZ	1:F:657:CYS:HB2	2.52	0.45
1:A:662:VAL:HG22	1:A:662:VAL:O	2.16	0.44
1:B:589:ASN:HB3	1:C:626:ILE:HD11	1.98	0.43
1:C:655:PHE:CZ	1:C:657:CYS:HB2	2.53	0.43
1:F:664:PRO:O	1:F:665:LEU:HB3	2.18	0.43
1:E:613:ASP:N	1:E:613:ASP:OD1	2.52	0.42
1:F:662:VAL:O	1:F:662:VAL:CG2	2.68	0.42
1:E:554:PHE:CZ	1:E:584:ARG:HG2	2.55	0.42
1:A:613:ASP:N	1:A:613:ASP:OD1	2.53	0.41
1:B:581:ASP:HB3	1:C:599:HIS:CE1	2.55	0.41
1:F:664:PRO:O	1:F:665:LEU:CB	2.68	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	127/137 (93%)	126 (99%)	1 (1%)	0	100	100
1	B	121/137 (88%)	120 (99%)	1 (1%)	0	100	100
1	C	127/137 (93%)	125 (98%)	2 (2%)	0	100	100
1	D	125/137 (91%)	124 (99%)	1 (1%)	0	100	100
1	E	132/137 (96%)	130 (98%)	2 (2%)	0	100	100
1	F	125/137 (91%)	124 (99%)	1 (1%)	0	100	100
1	G	123/137 (90%)	121 (98%)	1 (1%)	1 (1%)	19	28
All	All	880/959 (92%)	870 (99%)	9 (1%)	1 (0%)	51	69

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	613	ASP

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	103/117 (88%)	95 (92%)	8 (8%)	12	19
1	B	98/117 (84%)	94 (96%)	4 (4%)	30	47
1	C	98/117 (84%)	97 (99%)	1 (1%)	76	86
1	D	103/117 (88%)	100 (97%)	3 (3%)	42	60
1	E	107/117 (92%)	103 (96%)	4 (4%)	34	51
1	F	102/117 (87%)	99 (97%)	3 (3%)	42	60
1	G	97/117 (83%)	91 (94%)	6 (6%)	18	28
All	All	708/819 (86%)	679 (96%)	29 (4%)	30	47

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	G	571	GLU

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Mol	Chain	Res	Type
1	G	584	ARG
1	G	598	ILE
1	G	612	GLU
1	G	635	SER
1	G	647	ASP
1	A	533	MET
1	A	534	ARG
1	A	547	GLU
1	A	571	GLU
1	A	598	ILE
1	A	603	LEU
1	A	613	ASP
1	A	644	HIS
1	B	536	GLN
1	B	571	GLU
1	B	598	ILE
1	B	647	ASP
1	C	578	GLU
1	D	541	THR
1	D	598	ILE
1	D	666	GLN
1	E	596	LYS
1	E	613	ASP
1	E	645	ARG
1	E	647	ASP
1	F	584	ARG
1	F	662	VAL
1	F	666	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	G	583	HIS
1	G	599	HIS
1	G	624	GLN
1	G	636	GLN
1	G	652	ASN
1	A	544	GLN
1	A	636	GLN
1	C	652	ASN
1	D	544	GLN
1	D	651	GLN

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Mol	Chain	Res	Type
1	E	594	ASN
1	E	651	GLN
1	F	651	GLN
1	F	666	GLN

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](#)

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	MLT	D	701	-	2,8,8	0.55	0	3,10,10	0.60	0
2	MLT	F	701	-	2,8,8	0.51	0	3,10,10	0.82	0
2	MLT	A	701	-	2,8,8	0.55	0	3,10,10	0.72	0
2	MLT	E	701	-	2,8,8	1.04	0	3,10,10	1.77	1 (33%)

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	MLT	D	701	-	-	0/2/8/8	-
2	MLT	F	701	-	-	0/2/8/8	-
2	MLT	A	701	-	-	2/2/8/8	-
2	MLT	E	701	-	-	2/2/8/8	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	701	MLT	C3-C2-C1	2.68	114.50	111.10

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	701	MLT	C1-C2-C3-C4
2	A	701	MLT	O3-C2-C3-C4
2	E	701	MLT	C1-C2-C3-C4
2	E	701	MLT	O3-C2-C3-C4

There are no ring outliers.

No monomer is involved in short contacts.

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	131/137 (95%)	-0.15	2 (1%) 73 71	31, 50, 90, 115	0
1	B	125/137 (91%)	-0.25	2 (1%) 72 69	31, 53, 80, 135	0
1	C	131/137 (95%)	-0.05	4 (3%) 49 45	36, 63, 102, 137	0
1	D	129/137 (94%)	-0.12	4 (3%) 49 45	37, 55, 87, 117	0
1	E	134/137 (97%)	-0.44	0 100 100	31, 50, 84, 123	0
1	F	129/137 (94%)	-0.07	5 (3%) 39 36	36, 55, 92, 117	0
1	G	127/137 (92%)	-0.14	2 (1%) 72 69	38, 59, 95, 117	0
All	All	906/959 (94%)	-0.17	19 (2%) 63 60	31, 54, 94, 137	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	664	PRO	4.4
1	B	532	HIS	4.1
1	C	532	HIS	4.0
1	C	591	LEU	3.8
1	A	664	PRO	3.4
1	D	646	ARG	3.4
1	C	647	ASP	3.4
1	D	664	PRO	3.3
1	F	648	GLY	3.0
1	C	628	GLY	2.9
1	G	647	ASP	2.8
1	A	665	LEU	2.4
1	F	613	ASP	2.4
1	B	647	ASP	2.3
1	F	628	GLY	2.2
1	D	628	GLY	2.2
1	F	611	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	647	ASP	2.0
1	G	590	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, 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
2	MLT	E	701	9/9	0.80	0.27	78,89,92,99	0
2	MLT	F	701	9/9	0.85	0.33	69,83,86,89	0
2	MLT	A	701	9/9	0.88	0.27	73,85,89,90	0
2	MLT	D	701	9/9	0.89	0.25	68,87,93,93	0

6.5 Other polymers [i](#)

There are no such residues in this entry.