



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:15 AM GMT

PDB ID : 2CFH
Title : STRUCTURE OF THE BET3-TPC6B CORE OF TRAPP
Authors : Kummel, D.; Muller, J.J.; Roske, Y.; Henke, N.; Heinemann, U.
Deposited on : 2006-02-21
Resolution : 2.30 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
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

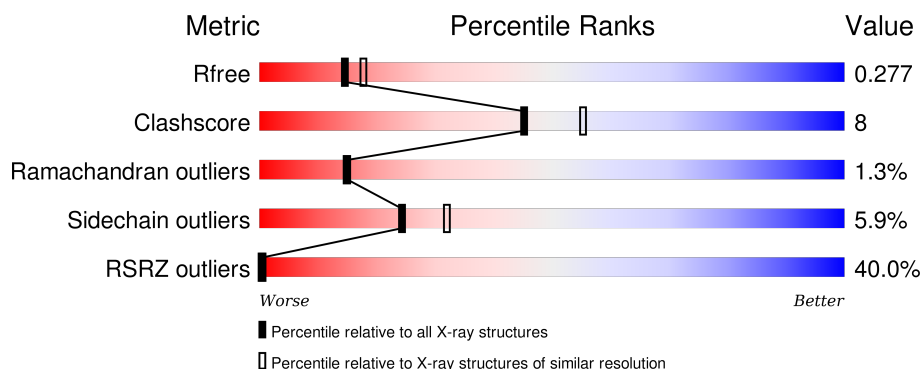
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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. 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	194	<div> <div>32%</div> <div>70% 11% • 16%</div> </div>
1	B	194	<div> <div>36%</div> <div>71% 11% • 16%</div> </div>
2	C	158	<div> <div>37%</div> <div>70% 20% 6% •</div> </div>
2	D	158	<div> <div>37%</div> <div>70% 22% • •</div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5049 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.

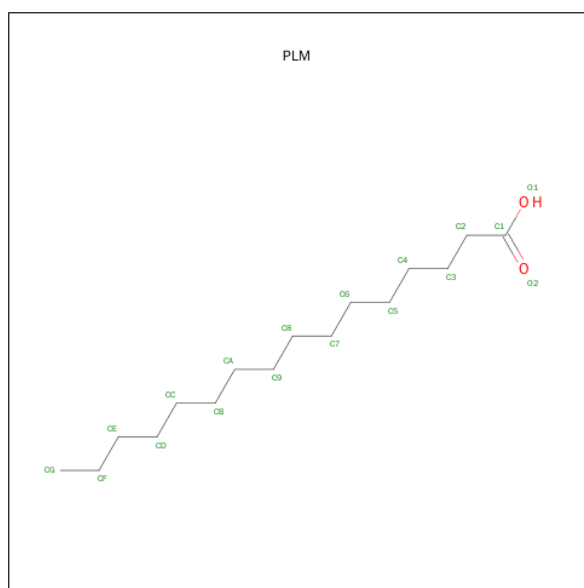
- Molecule 1 is a protein called TRAFFICKING PROTEIN PARTICLE COMPLEX SUB-UNIT 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	162	Total	C	N	O	S	0	0	1
			1276	807	215	246	8			
1	B	162	Total	C	N	O	S	0	0	1
			1276	807	215	246	8			

- Molecule 2 is a protein called TRAFFICKING PROTEIN PARTICLE COMPLEX SUB-UNIT 6B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	151	Total	C	N	O	S	0	0	1
			1199	764	206	220	9			
2	D	152	Total	C	N	O	S	0	0	1
			1207	769	207	221	10			

- Molecule 3 is PALMITIC ACID (three-letter code: PLM) (formula: C₁₆H₃₂O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			17	16	1		
3	B	1	Total	C	O	0	0
			17	16	1		

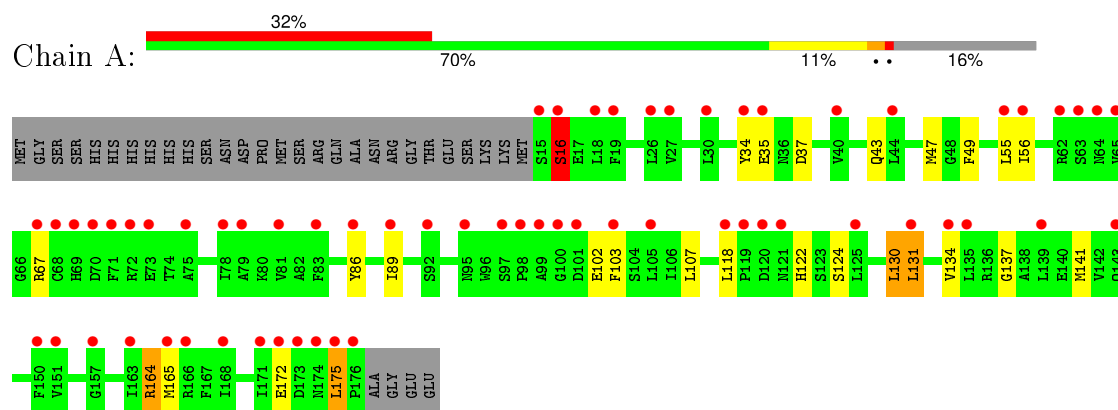
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	18	Total	O	0	0
			18	18		
4	B	12	Total	O	0	0
			12	12		
4	C	10	Total	O	0	0
			10	10		
4	D	17	Total	O	0	0
			17	17		

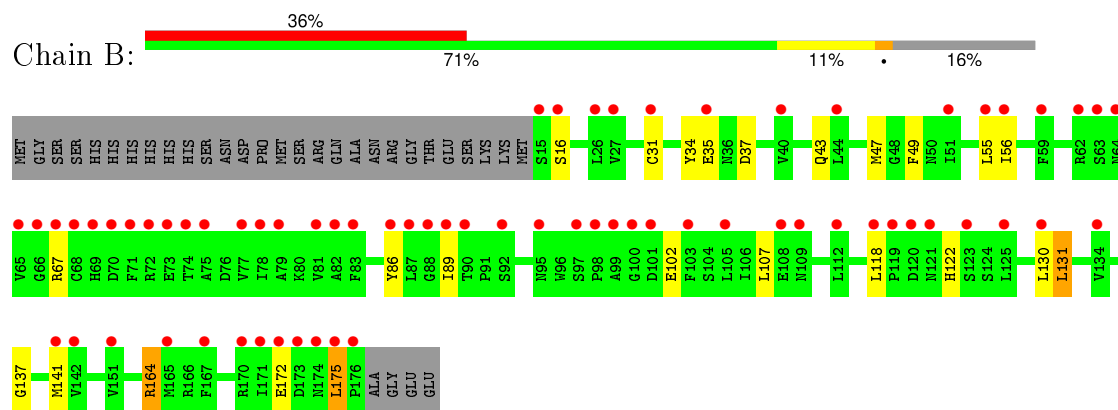
3 Residue-property plots

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.

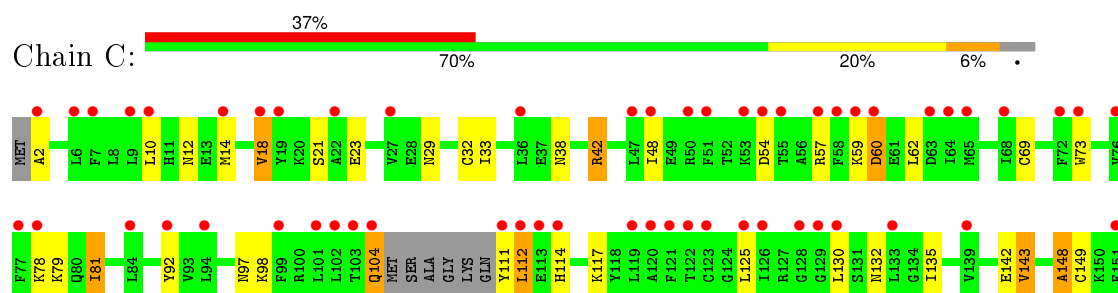
• Molecule 1: TRAFFICKING PROTEIN PARTICLE COMPLEX SUBUNIT 3



• Molecule 1: TRAFFICKING PROTEIN PARTICLE COMPLEX SUBUNIT 3

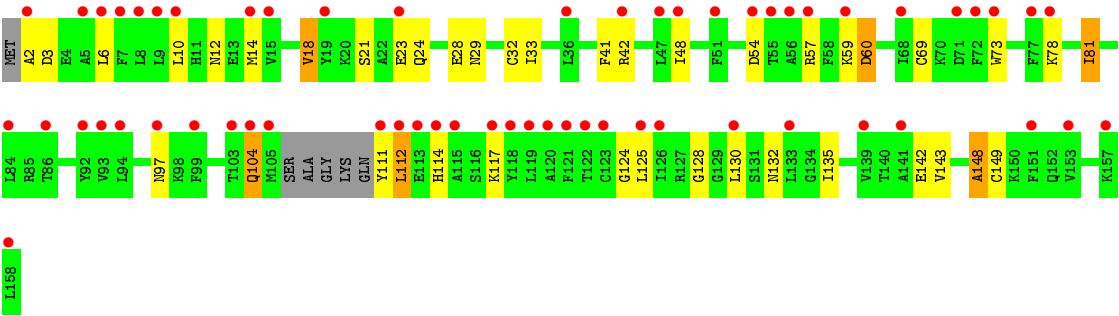


• Molecule 2: TRAFFICKING PROTEIN PARTICLE COMPLEX SUBUNIT 6B





● Molecule 2: TRAFFICKING PROTEIN PARTICLE COMPLEX SUBUNIT 6B



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	69.45Å 69.59Å 144.01Å 90.00° 91.55° 90.00°	Depositor
Resolution (Å)	20.00 – 2.30 19.93 – 2.30	Depositor EDS
% Data completeness (in resolution range)	100.0 (20.00-2.30) 98.7 (19.93-2.30)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.85 (at 2.30Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.221 , 0.269 0.230 , 0.277	Depositor DCC
R_{free} test set	1513 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	46.6	Xtriage
Anisotropy	0.121	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 61.5	EDS
Estimated twinning fraction	0.000 for k,h,-l 0.000 for -k,-h,-l 0.035 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 30262 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5049	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 34.02 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 7.2959e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: CSO, PLM

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	0.70	1/1296 (0.1%)	0.72	1/1752 (0.1%)
1	B	0.70	2/1296 (0.2%)	0.72	1/1752 (0.1%)
2	C	0.98	3/1209 (0.2%)	0.87	2/1619 (0.1%)
2	D	0.79	0/1217	0.80	0/1629
All	All	0.80	6/5018 (0.1%)	0.78	4/6752 (0.1%)

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
2	C	0	1
2	D	0	1
All	All	0	4

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	157	LYS	CE-NZ	20.81	2.01	1.49
2	C	157	LYS	CD-CE	8.59	1.72	1.51
2	C	157	LYS	CG-CD	6.87	1.75	1.52
1	B	175	LEU	C-N	-5.48	1.23	1.34
1	A	175	LEU	C-N	-5.40	1.24	1.34
1	B	31	CYS	CB-SG	-5.07	1.73	1.81

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	157	LYS	CD-CE-NZ	-15.16	76.83	111.70
1	B	67	ARG	C-N-CA	7.19	139.68	121.70
1	A	67	ARG	C-N-CA	7.09	139.43	121.70
2	C	157	LYS	CG-CD-CE	-5.18	96.37	111.90

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	16	SER	Peptide
1	B	16	SER	Peptide
2	C	148	ALA	Peptide
2	D	148	ALA	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	1276	0	1263	18	0
1	B	1276	0	1263	13	0
2	C	1199	0	1208	30	0
2	D	1207	0	1217	23	0
3	A	17	0	31	1	0
3	B	17	0	31	2	0
4	A	18	0	0	0	0
4	B	12	0	0	0	0
4	C	10	0	0	0	0
4	D	17	0	0	0	0
All	All	5049	0	5013	78	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 8.

All (78) 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)
2:C:157:LYS:CD	2:C:157:LYS:CG	1.75	1.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:32:CSO:SG	2:C:32:CSO:OD	1.92	1.28
2:D:32:CSO:OD	2:D:32:CSO:SG	1.93	1.25
2:C:157:LYS:NZ	2:C:157:LYS:CE	2.01	1.24
2:C:157:LYS:CD	2:C:157:LYS:NZ	2.33	0.92
2:C:69:CYS:HA	2:C:81:ILE:HD11	1.52	0.91
2:D:69:CYS:HA	2:D:81:ILE:HD11	1.51	0.91
2:C:157:LYS:CG	2:C:157:LYS:CE	2.59	0.80
2:C:157:LYS:CD	2:C:157:LYS:CB	2.60	0.79
2:D:21:SER:OG	2:D:29:ASN:ND2	2.27	0.67
2:D:69:CYS:HA	2:D:81:ILE:CD1	2.25	0.65
2:C:21:SER:OG	2:C:29:ASN:ND2	2.31	0.64
2:C:69:CYS:HA	2:C:81:ILE:CD1	2.28	0.62
1:A:16:SER:HB3	2:C:2:ALA:HB3	1.82	0.61
1:B:34:TYR:O	1:B:35:GLU:HG2	2.01	0.61
1:B:55:LEU:HD11	1:B:86:TYR:CE1	2.39	0.57
1:A:89:ILE:HD12	1:A:107:LEU:HD22	1.88	0.56
1:B:55:LEU:HD11	1:B:86:TYR:CZ	2.42	0.55
1:A:56:ILE:HG21	1:A:141:MET:HB2	1.87	0.54
1:A:55:LEU:HD11	1:A:86:TYR:CE1	2.41	0.54
1:B:56:ILE:HG21	1:B:141:MET:HB2	1.89	0.54
2:C:143:VAL:HG23	2:C:148:ALA:O	2.08	0.54
2:D:97:ASN:HA	2:D:148:ALA:HB2	1.90	0.53
1:B:86:TYR:HA	2:D:3:ASP:HB2	1.91	0.51
1:A:16:SER:HB2	2:C:2:ALA:O	2.09	0.51
2:D:69:CYS:SG	2:D:81:ILE:HG13	2.50	0.51
2:C:97:ASN:HA	2:C:148:ALA:HB2	1.92	0.51
1:B:43:GLN:HE21	1:B:47:MET:HG3	1.77	0.50
1:B:89:ILE:HD12	1:B:107:LEU:HD22	1.93	0.50
1:A:34:TYR:O	1:A:35:GLU:HG2	2.12	0.50
1:A:102:GLU:OE1	1:A:164:ARG:NH1	2.45	0.50
2:C:130:LEU:HB3	2:C:135:ILE:HB	1.93	0.50
1:B:37:ASP:OD2	1:B:122:HIS:HA	2.12	0.49
1:A:55:LEU:HD11	1:A:86:TYR:CZ	2.49	0.47
1:B:49:PHE:CE1	1:B:137:GLY:HA2	2.48	0.47
1:A:49:PHE:CE1	1:A:137:GLY:HA2	2.50	0.47
2:C:73:TRP:CG	2:C:81:ILE:HD13	2.49	0.47
1:B:102:GLU:OE1	1:B:164:ARG:NH1	2.48	0.46
2:C:62:LEU:HD22	2:C:92:TYR:OH	2.15	0.46
2:D:143:VAL:HG23	2:D:148:ALA:O	2.15	0.46
2:C:69:CYS:SG	2:C:81:ILE:HG13	2.56	0.46
2:C:79:LYS:NZ	2:C:98:LYS:O	2.43	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:14:MET:HA	2:D:18:VAL:HG23	1.98	0.46
2:D:111:TYR:CD2	2:D:112:LEU:N	2.84	0.46
2:D:130:LEU:HB3	2:D:135:ILE:HB	1.97	0.46
2:D:73:TRP:CG	2:D:81:ILE:HD13	2.52	0.45
1:A:37:ASP:OD2	1:A:122:HIS:HA	2.16	0.45
2:C:111:TYR:CD2	2:C:112:LEU:N	2.85	0.45
2:D:48:ILE:HG21	2:D:132:ASN:HB3	1.99	0.45
2:C:48:ILE:HG21	2:C:132:ASN:HB3	1.98	0.44
2:D:24:GLN:HB2	2:D:28:GLU:OE1	2.18	0.44
1:A:47:MET:CE	2:C:18:VAL:CG2	2.95	0.44
2:C:114:HIS:HB3	2:C:117:LYS:HD2	1.99	0.44
1:A:16:SER:CB	2:C:2:ALA:HB3	2.48	0.44
2:C:38:ASN:O	2:C:42:ARG:NE	2.51	0.43
1:A:131:LEU:HD21	3:A:1068:PLM:HC2	1.98	0.43
2:D:12:ASN:HB2	2:D:104:GLN:HG3	2.00	0.43
2:C:60:ASP:OD1	2:C:60:ASP:C	2.56	0.43
2:D:33:ILE:HD13	2:D:33:ILE:HA	1.85	0.43
1:A:103:PHE:CZ	1:A:165:MET:HG3	2.53	0.43
2:C:59:LYS:O	2:C:60:ASP:HB3	2.19	0.43
2:C:33:ILE:HA	2:C:33:ILE:HD13	1.86	0.43
2:C:14:MET:HA	2:C:18:VAL:HG23	2.01	0.42
2:C:12:ASN:HB2	2:C:104:GLN:HG3	2.00	0.42
2:D:41:PHE:CE1	2:D:128:GLY:HA2	2.55	0.42
2:C:23:GLU:HA	2:C:29:ASN:HD21	1.85	0.42
2:D:23:GLU:HA	2:D:29:ASN:HD21	1.84	0.41
1:B:86:TYR:O	2:D:2:ALA:HA	2.21	0.41
1:A:130:LEU:O	1:A:134:VAL:HG23	2.20	0.41
2:D:41:PHE:HA	2:D:124:GLY:O	2.20	0.41
1:B:131:LEU:HD21	3:B:1068:PLM:HC2	2.02	0.41
3:B:1068:PLM:HF1	2:D:6:LEU:HD11	2.02	0.41
2:D:59:LYS:O	2:D:60:ASP:HB3	2.21	0.41
1:A:37:ASP:HB2	1:A:124:SER:HB2	2.02	0.41
1:A:43:GLN:NE2	1:A:43:GLN:HA	2.35	0.40
1:A:164:ARG:HA	1:A:164:ARG:HD3	1.84	0.40
2:D:114:HIS:HB3	2:D:117:LYS:HD2	2.03	0.40
1:B:164:ARG:HA	1:B:164:ARG:HD3	1.85	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	160/194 (82%)	154 (96%)	5 (3%)	1 (1%)	30	36
1	B	160/194 (82%)	154 (96%)	5 (3%)	1 (1%)	30	36
2	C	146/158 (92%)	132 (90%)	11 (8%)	3 (2%)	9	7
2	D	147/158 (93%)	131 (89%)	13 (9%)	3 (2%)	9	7
All	All	613/704 (87%)	571 (93%)	34 (6%)	8 (1%)	15	15

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	C	60	ASP
2	C	149	CYS
2	D	60	ASP
2	D	149	CYS
2	C	112	LEU
2	D	112	LEU
1	A	175	LEU
1	B	175	LEU

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	142/170 (84%)	136 (96%)	6 (4%)	36	49
1	B	142/170 (84%)	137 (96%)	5 (4%)	43	58

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	C	129/136 (95%)	118 (92%)	11 (8%)	13	16
2	D	130/136 (96%)	120 (92%)	10 (8%)	16	20
All	All	543/612 (89%)	511 (94%)	32 (6%)	24	32

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	SER
1	A	118	LEU
1	A	130	LEU
1	A	131	LEU
1	A	164	ARG
1	A	172	GLU
1	B	118	LEU
1	B	130	LEU
1	B	131	LEU
1	B	164	ARG
1	B	172	GLU
2	C	10	LEU
2	C	18	VAL
2	C	42	ARG
2	C	54	ASP
2	C	57	ARG
2	C	78	LYS
2	C	81	ILE
2	C	104	GLN
2	C	125	LEU
2	C	142	GLU
2	C	143	VAL
2	D	10	LEU
2	D	18	VAL
2	D	42	ARG
2	D	54	ASP
2	D	57	ARG
2	D	78	LYS
2	D	81	ILE
2	D	104	GLN
2	D	125	LEU
2	D	142	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	GLN
1	B	43	GLN
2	C	29	ASN
2	C	38	ASN
2	C	83	ASN
2	C	95	GLN
2	D	29	ASN
2	D	38	ASN
2	D	83	ASN
2	D	95	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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	CSO	C	32	2	3,6,7	0.50	0	1,6,8	1.13	0
2	CSO	D	32	2	3,6,7	0.37	0	1,6,8	1.25	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	CSO	C	32	2	-	0/1/5/7	0/0/0/0
2	CSO	D	32	2	-	0/1/5/7	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	32	CSO	1	0
2	D	32	CSO	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates 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	$\# Z > 2$
3	PLM	A	1068	1	16,16,17	0.56	0	14,15,17	0.60	0
3	PLM	B	1068	1	16,16,17	0.60	0	14,15,17	0.63	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	PLM	A	1068	1	-	0/13/14/15	0/0/0/0
3	PLM	B	1068	1	-	0/13/14/15	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1068	PLM	1	0
3	B	1068	PLM	2	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, 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	162/194 (83%)	1.98	63 (38%) 0 0	57, 59, 60, 61	0
1	B	162/194 (83%)	2.16	69 (42%) 0 0	57, 59, 61, 62	0
2	C	150/158 (94%)	1.88	59 (39%) 0 0	54, 59, 61, 64	0
2	D	151/158 (95%)	2.03	59 (39%) 0 0	54, 59, 61, 65	0
All	All	625/704 (88%)	2.02	250 (40%) 0 0	54, 59, 61, 65	0

All (250) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	175	LEU	12.6
2	D	105	MET	11.2
1	B	176	PRO	10.2
1	B	99	ALA	10.1
2	C	111	TYR	9.8
2	D	56	ALA	8.4
2	D	158	LEU	8.1
1	B	78	ILE	7.8
1	A	98	PRO	7.6
1	A	176	PRO	7.5
1	A	99	ALA	7.3
2	D	57	ARG	6.9
1	B	65	VAL	6.9
2	C	54	ASP	6.8
1	B	67	ARG	6.8
1	B	66	GLY	6.7
2	D	111	TYR	6.7
1	B	68	CYS	6.6
2	D	54	ASP	6.6
1	A	16	SER	6.5
2	C	112	LEU	6.5

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Mol	Chain	Res	Type	RSRZ
1	B	100	GLY	6.4
2	C	103	THR	6.0
2	D	72	PHE	6.0
1	B	64	ASN	6.0
1	A	92	SER	5.9
2	C	55	THR	5.9
1	A	100	GLY	5.8
2	C	57	ARG	5.8
1	B	173	ASP	5.6
2	D	125	LEU	5.6
1	A	71	PHE	5.5
1	A	68	CYS	5.5
1	A	67	ARG	5.5
1	A	69	HIS	5.4
2	C	125	LEU	5.3
2	D	48	ILE	5.2
1	B	71	PHE	5.1
2	C	10	LEU	5.1
1	B	73	GLU	5.0
1	B	90	THR	4.8
1	B	88	GLY	4.8
1	B	97	SER	4.8
1	B	69	HIS	4.8
1	B	75	ALA	4.7
2	D	120	ALA	4.7
2	C	94	LEU	4.6
2	D	126	ILE	4.6
1	B	175	LEU	4.5
1	A	70	ASP	4.5
1	B	105	LEU	4.4
2	C	114	HIS	4.4
2	C	48	ILE	4.4
2	D	7	PHE	4.3
2	C	122	THR	4.3
1	A	174	ASN	4.3
1	B	82	ALA	4.3
1	A	143	GLN	4.2
1	B	83	PHE	4.2
2	D	84	LEU	4.2
1	A	27	VAL	4.2
2	D	10	LEU	4.2
2	D	114	HIS	4.1

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Mol	Chain	Res	Type	RSRZ
1	B	119	PRO	4.1
1	A	30	LEU	4.0
2	C	68	ILE	4.0
1	A	118	LEU	4.0
2	C	72	PHE	4.0
1	B	174	ASN	4.0
2	D	104	GLN	4.0
2	D	103	THR	3.9
2	C	104	GLN	3.9
1	B	125	LEU	3.9
1	B	98	PRO	3.9
1	A	78	ILE	3.9
2	D	122	THR	3.9
2	C	63	ASP	3.8
1	A	131	LEU	3.8
2	D	68	ILE	3.8
2	C	126	ILE	3.8
1	A	89	ILE	3.7
2	D	8	LEU	3.7
1	B	70	ASP	3.7
1	A	101	ASP	3.7
2	C	60	ASP	3.6
2	D	6	LEU	3.6
2	C	121	PHE	3.6
1	A	105	LEU	3.6
1	A	172	GLU	3.6
1	B	74	THR	3.6
1	A	139	LEU	3.6
2	D	113	GLU	3.6
1	B	16	SER	3.5
2	D	9	LEU	3.5
1	B	95	ASN	3.5
2	C	102	LEU	3.5
1	A	103	PHE	3.5
1	B	165	MET	3.4
1	B	170	ARG	3.4
2	D	119	LEU	3.4
1	B	59	PHE	3.3
2	D	99	PHE	3.3
1	A	151	VAL	3.3
2	C	53	LYS	3.3
1	B	167	PHE	3.3

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Mol	Chain	Res	Type	RSRZ
2	C	157	LYS	3.3
2	C	6	LEU	3.3
2	D	153	VAL	3.3
1	A	168	ILE	3.2
2	D	2	ALA	3.2
2	D	93	VAL	3.2
1	B	171	ILE	3.2
2	C	158	LEU	3.2
1	A	15	SER	3.2
1	B	103	PHE	3.2
1	B	123	SER	3.2
1	A	63	SER	3.2
2	D	59	LYS	3.2
2	C	51	PHE	3.1
1	B	87	LEU	3.1
2	D	121	PHE	3.1
1	B	56	ILE	3.1
2	D	47	LEU	3.1
2	C	59	LYS	3.1
2	C	64	ILE	3.1
1	B	26	LEU	3.0
1	B	86	TYR	3.0
2	C	123	CYS	3.0
1	B	81	VAL	3.0
1	B	172	GLU	3.0
2	C	113	GLU	3.0
1	B	79	ALA	3.0
2	D	130	LEU	3.0
1	A	119	PRO	3.0
2	C	2	ALA	3.0
2	D	36	LEU	3.0
1	B	89	ILE	2.9
1	A	173	ASP	2.9
2	C	120	ALA	2.9
2	D	157	LYS	2.9
2	C	151	PHE	2.9
2	D	133	LEU	2.9
2	D	5	ALA	2.9
2	C	58	PHE	2.9
2	D	123	CYS	2.9
1	B	77	VAL	2.9
2	C	36	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
2	D	15	VAL	2.8
1	A	95	ASN	2.8
2	D	112	LEU	2.8
2	D	51	PHE	2.8
2	C	101	LEU	2.8
1	B	142	VAL	2.8
1	A	26	LEU	2.8
2	C	73	TRP	2.8
2	D	151	PHE	2.8
1	A	35	GLU	2.8
2	D	139	VAL	2.8
1	A	83	PHE	2.8
1	A	135	LEU	2.8
1	B	151	VAL	2.7
1	A	171	ILE	2.7
2	C	9	LEU	2.7
1	B	44	LEU	2.7
1	B	27	VAL	2.7
1	B	51	ILE	2.7
1	A	55	LEU	2.7
2	D	92	TYR	2.7
1	A	166	ARG	2.7
2	C	19	TYR	2.7
1	B	118	LEU	2.6
2	C	84	LEU	2.6
1	A	73	GLU	2.6
2	D	55	THR	2.6
2	D	14	MET	2.6
1	A	134	VAL	2.6
1	B	108	GLU	2.6
2	D	86	THR	2.6
1	B	121	ASN	2.6
2	C	50	ARG	2.6
1	A	44	LEU	2.6
2	D	78	LYS	2.5
2	D	118	TYR	2.5
1	A	65	VAL	2.5
2	D	77	PHE	2.5
1	B	55	LEU	2.5
1	A	64	ASN	2.5
2	D	94	LEU	2.5
1	A	19	PHE	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	165	MET	2.5
2	D	23	GLU	2.5
1	B	72	ARG	2.4
1	A	121	ASN	2.4
2	C	130	LEU	2.4
1	B	92	SER	2.4
1	A	79	ALA	2.4
2	D	141	ALA	2.4
1	A	62	ARG	2.4
1	B	62	ARG	2.4
2	C	78	LYS	2.4
2	D	42	ARG	2.4
1	A	157	GLY	2.4
1	B	63	SER	2.4
1	B	112	LEU	2.4
2	C	47	LEU	2.4
2	C	139	VAL	2.3
2	D	73	TRP	2.3
1	B	15	SER	2.3
2	D	19	TYR	2.3
1	B	134	VAL	2.3
2	C	22	ALA	2.3
2	C	99	PHE	2.3
1	B	101	ASP	2.3
1	B	120	ASP	2.3
2	D	71	ASP	2.3
2	C	92	TYR	2.2
2	C	129	GLY	2.2
1	A	72	ARG	2.2
1	A	97	SER	2.2
1	A	56	ILE	2.2
1	A	34	TYR	2.2
2	C	14	MET	2.2
1	A	163	ILE	2.2
1	A	40	VAL	2.2
1	A	86	TYR	2.2
1	A	81	VAL	2.1
1	B	40	VAL	2.1
2	C	27	VAL	2.1
2	C	77	PHE	2.1
1	B	130	LEU	2.1
1	B	31	CYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	120	ASP	2.1
1	A	75	ALA	2.1
1	A	150	PHE	2.1
1	A	18	LEU	2.1
2	D	115	ALA	2.1
2	C	18	VAL	2.1
1	A	125	LEU	2.1
2	C	128	GLY	2.1
2	D	117	LYS	2.1
2	C	76	VAL	2.1
1	B	109	ASN	2.0
2	C	119	LEU	2.0
2	C	133	LEU	2.0
2	C	153	VAL	2.0
2	C	7	PHE	2.0
1	B	35	GLU	2.0
2	D	97	ASN	2.0
1	B	141	MET	2.0
2	C	65	MET	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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. 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	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	CSO	D	32	7/8	0.94	0.19	-	48,57,57,57	0
2	CSO	C	32	7/8	0.81	0.21	-	48,57,57,58	0

6.3 Carbohydrates [i](#)

There are no carbohydrates 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. 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	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	PLM	B	1068	17/18	0.59	0.37	0.44	58,59,61,61	0
3	PLM	A	1068	17/18	0.67	0.27	0.11	58,59,61,61	0

6.5 Other polymers

There are no such residues in this entry.