



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:02 PM GMT

PDB ID : 4K8N  
Title : Crystal structure of human ceramide-1-phosphate transfer protein (CPTP) in complex with 18:1 Ceramide-1-Phosphate (18:1-C1P)  
Authors : Simanshu, D.K.; Brown, R.E.; Patel, D.J.  
Deposited on : 2013-04-18  
Resolution : 3.10 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](mailto: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.

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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 : 1.7 (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

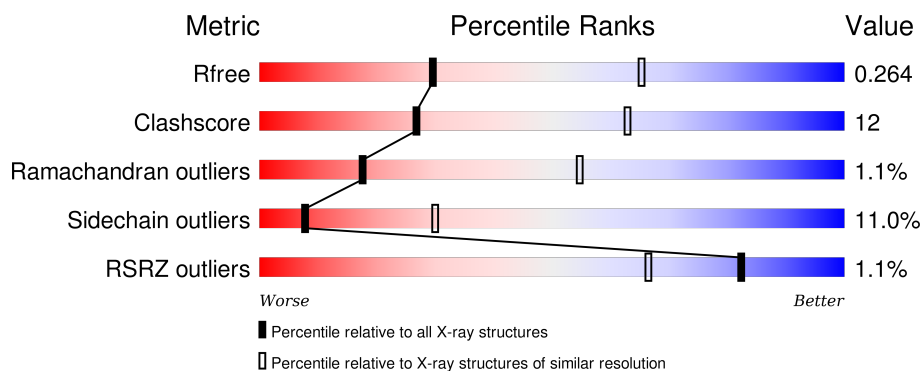
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.10 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))
$R_{free}$	91344	1114 (3.14-3.06)
Clashscore	102246	1222 (3.14-3.06)
Ramachandran outliers	100387	1174 (3.14-3.06)
Sidechain outliers	100360	1174 (3.14-3.06)
RSRZ outliers	91569	1119 (3.14-3.06)

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  $\geq 3$ , 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	215	
1	B	215	
1	C	215	
1	D	215	
1	E	215	

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Mol	Chain	Length	Quality of chain
1	F	215	<div><div></div><div>6%</div><div>53%</div><div>34%</div><div>8%</div><div>5%</div></div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10099 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.

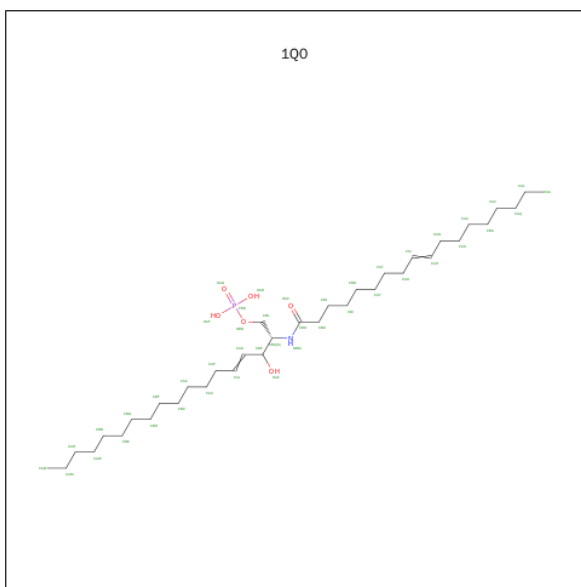
- Molecule 1 is a protein called Glycolipid transfer protein domain-containing protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	207	Total	C	N	O	S	0	0	0
			1659	1060	295	296	8			
1	B	206	Total	C	N	O	S	0	0	0
			1644	1048	293	295	8			
1	C	206	Total	C	N	O	S	0	0	0
			1637	1045	293	291	8			
1	D	206	Total	C	N	O	S	0	0	0
			1651	1051	297	295	8			
1	E	207	Total	C	N	O	S	0	0	0
			1653	1055	296	294	8			
1	F	205	Total	C	N	O	S	0	0	0
			1621	1037	288	288	8			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	EXPRESSION TAG	UNP Q5TA50
B	0	SER	-	EXPRESSION TAG	UNP Q5TA50
C	0	SER	-	EXPRESSION TAG	UNP Q5TA50
D	0	SER	-	EXPRESSION TAG	UNP Q5TA50
E	0	SER	-	EXPRESSION TAG	UNP Q5TA50
F	0	SER	-	EXPRESSION TAG	UNP Q5TA50

- Molecule 2 is (2S,3R,4Z)-3-HYDROXY-2-[(9E)-OCTADEC-9-ENOYLAMINO]OCTADEC-4-EN-1-YL DIHYDROGEN PHOSPHATE (three-letter code: 1Q0) (formula: C<sub>36</sub>H<sub>70</sub>NO<sub>6</sub>P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			44	36	1	6	1		
2	B	1	Total	C	N	O	P	0	0
			27	19	1	6	1		
2	C	1	Total	C	N	O	P	0	0
			44	36	1	6	1		
2	D	1	Total	C	N	O	P	0	0
			44	36	1	6	1		
2	E	1	Total	C	N	O	P	0	0
			44	36	1	6	1		
2	F	1	Total	C	N	O	P	0	0
			16	8	1	6	1		

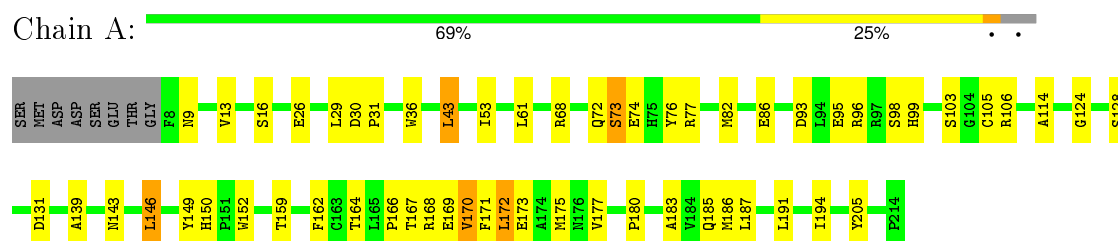
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	4	Total	O	0	0
			4	4		
3	B	3	Total	O	0	0
			3	3		
3	C	3	Total	O	0	0
			3	3		
3	D	2	Total	O	0	0
			2	2		
3	E	3	Total	O	0	0
			3	3		

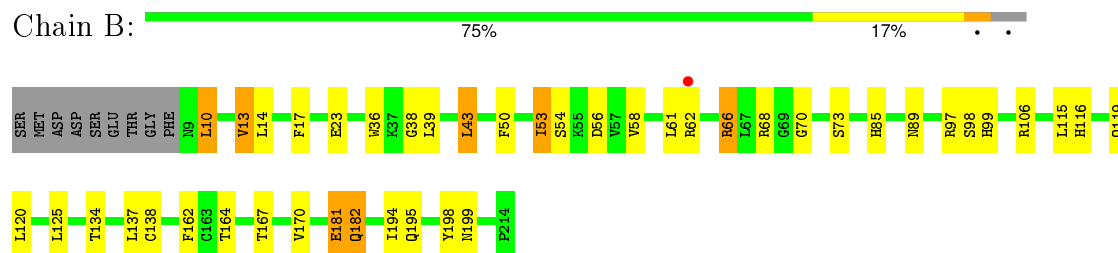
### 3 Residue-property plots

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.

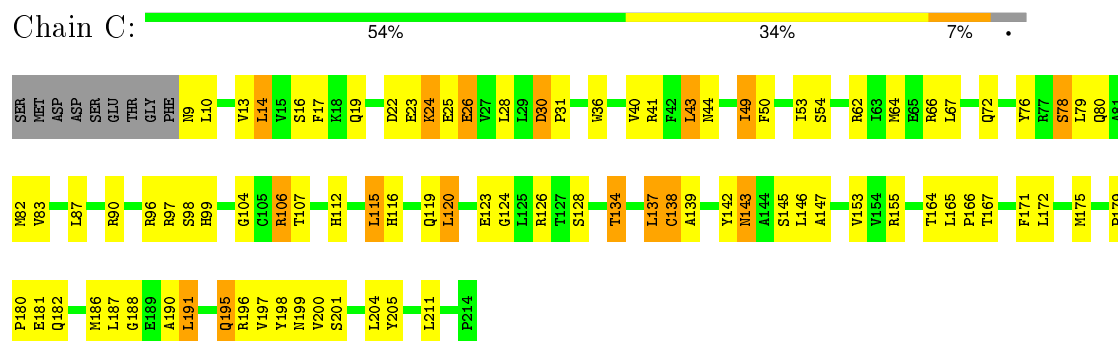
- Molecule 1: Glycolipid transfer protein domain-containing protein 1



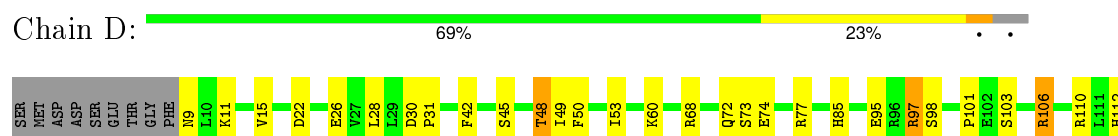
- Molecule 1: Glycolipid transfer protein domain-containing protein 1



- Molecule 1: Glycolipid transfer protein domain-containing protein 1

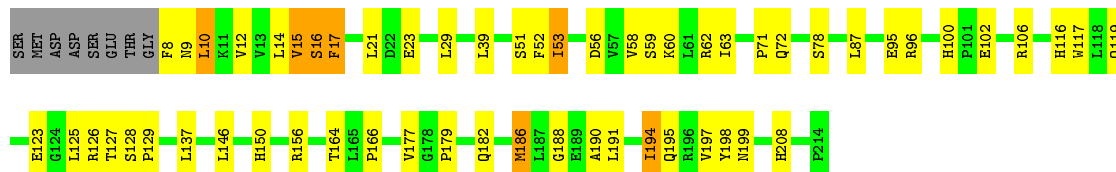


- Molecule 1: Glycolipid transfer protein domain-containing protein 1

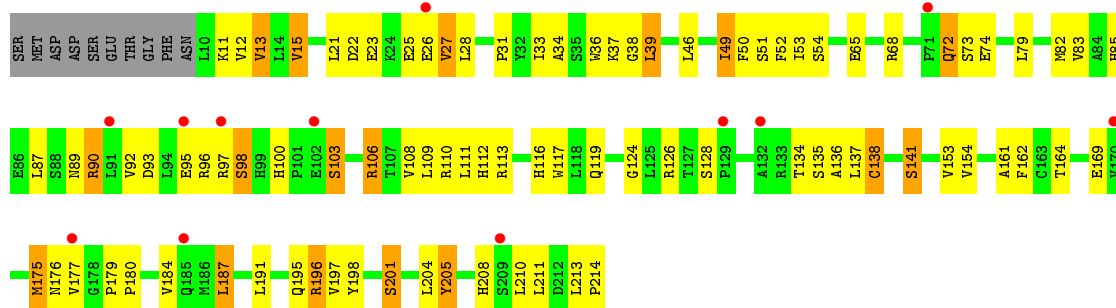




- Molecule 1: Glycolipid transfer protein domain-containing protein 1



- Molecule 1: Glycolipid transfer protein domain-containing protein 1



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	236.98Å 128.76Å 66.66Å 90.00° 93.53° 90.00°	Depositor
Resolution (Å)	30.05 – 3.10 30.05 – 3.10	Depositor EDS
% Data completeness (in resolution range)	98.5 (30.05-3.10) 94.4 (30.05-3.10)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.46 (at 3.11Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, $R_{free}$	0.225 , 0.263 0.226 , 0.264	Depositor DCC
$R_{free}$ test set	1786 reflections (5.49%)	DCC
Wilson B-factor (Å <sup>2</sup> )	66.6	Xtriage
Anisotropy	0.449	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.27 , 38.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	0 of 35756 reflections	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	10099	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	95.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.79% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: 1Q0

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	0.30	0/1697	0.46	0/2301
1	B	0.30	0/1681	0.54	2/2281 (0.1%)
1	C	0.29	0/1674	0.51	1/2272 (0.0%)
1	D	0.28	0/1688	0.47	1/2288 (0.0%)
1	E	0.27	0/1691	0.44	0/2295
1	F	0.27	0/1658	0.51	0/2252
All	All	0.29	0/10089	0.49	4/13689 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	66	ARG	NE-CZ-NH2	9.70	125.15	120.30
1	B	66	ARG	NE-CZ-NH1	-8.51	116.05	120.30
1	C	165	LEU	CB-CA-C	6.71	122.96	110.20
1	D	131	ASP	CB-CA-C	-6.16	98.07	110.40

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	1659	0	1670	31	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1644	0	1650	25	0
1	C	1637	0	1644	52	0
1	D	1651	0	1663	39	0
1	E	1653	0	1655	35	0
1	F	1621	0	1622	54	0
2	A	44	0	68	9	0
2	B	27	0	28	5	0
2	C	44	0	68	1	0
2	D	44	0	68	9	0
2	E	44	0	68	11	0
2	F	16	0	8	3	0
3	A	4	0	0	0	0
3	B	3	0	0	0	0
3	C	3	0	0	0	0
3	D	2	0	0	0	0
3	E	3	0	0	1	0
All	All	10099	0	10212	245	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 12.

All (245) 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:E:301:1Q0:CAO	2:E:301:1Q0:H68	2.09	0.83
1:D:53:ILE:HD12	2:D:301:1Q0:H38	1.65	0.77
1:A:53:ILE:HD12	2:A:301:1Q0:H40	1.66	0.77
1:E:117:TRP:CH2	2:E:301:1Q0:H8	2.23	0.73
1:F:191:LEU:HG	1:F:195:GLN:HE22	1.53	0.73
1:C:25:GLU:OE1	1:C:196:ARG:NH2	2.23	0.70
1:B:58:VAL:HA	1:B:61:LEU:HD12	1.72	0.70
1:D:74:GLU:O	1:D:77:ARG:NH1	2.26	0.69
2:E:301:1Q0:H44	2:E:301:1Q0:H68	1.74	0.69
1:B:195:GLN:O	1:B:199:ASN:ND2	2.25	0.69
1:F:113:ARG:NH2	1:F:211:LEU:O	2.28	0.67
1:B:119:GLN:HG3	1:B:195:GLN:HG2	1.76	0.66
1:A:114:ALA:HB1	2:A:301:1Q0:H42	1.77	0.66
1:D:128:SER:HB3	1:D:168:ARG:HE	1.60	0.65
2:E:301:1Q0:H45	2:E:301:1Q0:H68	1.79	0.65
1:C:143:ASN:HA	1:C:147:ALA:HB3	1.78	0.65
2:A:301:1Q0:H55	2:A:301:1Q0:CAI	2.28	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:79:LEU:HA	1:F:82:MET:HB3	1.80	0.64
1:C:195:GLN:O	1:C:199:ASN:ND2	2.31	0.63
1:C:98:SER:OG	1:C:99:HIS:N	2.30	0.63
1:C:30:ASP:N	1:C:30:ASP:OD1	2.30	0.63
1:B:13:VAL:HG21	1:B:38:GLY:HA3	1.79	0.63
1:F:135:SER:HB2	1:F:162:PHE:HD1	1.63	0.62
1:C:126:ARG:NH1	1:C:187:LEU:O	2.33	0.62
1:F:49:ILE:HG13	1:F:50:PHE:N	2.14	0.62
1:F:37:LYS:NZ	1:F:65:GLU:OE2	2.29	0.62
1:D:48:THR:OG1	1:D:49:ILE:N	2.30	0.62
1:C:139:ALA:O	1:C:143:ASN:ND2	2.32	0.61
1:D:122:LEU:HD21	2:D:301:1Q0:H63	1.83	0.61
1:B:85:HIS:O	1:B:89:ASN:ND2	2.33	0.61
1:C:9:ASN:OD1	1:C:41:ARG:NH2	2.32	0.60
1:A:26:GLU:OE1	1:A:77:ARG:NH2	2.34	0.60
1:A:149:TYR:OH	1:B:182:GLN:NE2	2.25	0.60
2:F:301:1Q0:OAC	2:F:301:1Q0:H29	2.02	0.59
1:C:22:ASP:OD1	1:C:26:GLU:N	2.35	0.58
1:B:98:SER:OG	1:B:99:HIS:N	2.37	0.58
1:C:147:ALA:O	1:C:155:ARG:NE	2.21	0.58
1:B:116:HIS:HB2	1:B:198:TYR:CE1	2.39	0.57
1:E:150:HIS:HE1	2:E:301:1Q0:H16	1.69	0.57
1:B:70:GLY:O	1:B:73:SER:OG	2.23	0.57
1:F:33:ILE:HD13	1:F:65:GLU:HG3	1.87	0.56
1:F:93:ASP:HB3	1:F:96:ARG:O	2.04	0.56
1:F:51:SER:O	1:F:54:SER:OG	2.23	0.56
1:E:116:HIS:HB2	1:E:198:TYR:CE1	2.40	0.56
1:C:40:VAL:HG13	1:C:54:SER:HB3	1.88	0.56
1:B:39:LEU:HD23	2:B:301:1Q0:H50	1.88	0.55
1:B:36:TRP:HB3	1:B:61:LEU:HD21	1.88	0.55
1:B:125:LEU:HA	1:B:137:LEU:HD13	1.88	0.55
2:B:301:1Q0:OBN	2:B:301:1Q0:OAE	2.22	0.55
1:C:13:VAL:O	1:C:16:SER:OG	2.25	0.55
1:D:30:ASP:OD1	1:D:68:ARG:NE	2.40	0.55
1:A:169:GLU:O	1:A:172:LEU:N	2.38	0.55
1:D:146:LEU:HG	1:D:150:HIS:HE1	1.71	0.55
1:D:146:LEU:O	1:D:150:HIS:ND1	2.37	0.55
1:F:72:GLN:HE21	1:F:85:HIS:CE1	2.25	0.54
2:A:301:1Q0:H55	2:A:301:1Q0:H46	1.89	0.54
1:D:97:ARG:HH21	1:D:106:ARG:HH12	1.53	0.54
2:A:301:1Q0:H46	2:A:301:1Q0:CAX	2.38	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:116:HIS:HB2	1:B:198:TYR:CZ	2.43	0.54
1:F:21:LEU:HB2	1:F:197:VAL:HG12	1.90	0.54
1:D:106:ARG:NH2	2:D:301:1Q0:OAF	2.40	0.54
1:A:114:ALA:HB2	2:A:301:1Q0:H70	1.90	0.54
1:C:126:ARG:HG3	1:C:187:LEU:HB2	1.90	0.53
1:A:146:LEU:HB3	1:A:150:HIS:CE1	2.44	0.53
1:E:14:LEU:HG	1:E:190:ALA:HB2	1.89	0.53
1:E:126:ARG:HG3	1:E:127:THR:H	1.73	0.53
1:A:68:ARG:NH1	1:A:73:SER:O	2.42	0.53
1:A:187:LEU:O	1:A:191:LEU:HB2	2.09	0.53
1:F:25:GLU:OE2	1:F:196:ARG:NH2	2.42	0.53
2:A:301:1Q0:CAI	2:A:301:1Q0:CAX	2.87	0.52
1:F:68:ARG:NH1	1:F:73:SER:O	2.42	0.52
1:B:56:ASP:OD2	2:B:301:1Q0:H24	2.09	0.52
1:C:9:ASN:N	1:C:175:MET:O	2.42	0.52
1:E:194:ILE:O	1:E:197:VAL:N	2.42	0.52
1:E:150:HIS:NE2	2:E:301:1Q0:OAC	2.24	0.51
1:C:14:LEU:HB3	1:C:190:ALA:HB1	1.92	0.51
1:E:195:GLN:O	1:E:199:ASN:ND2	2.43	0.51
1:B:56:ASP:OD2	2:B:301:1Q0:NBM	2.44	0.51
1:C:19:GLN:N	1:C:19:GLN:OE1	2.43	0.51
1:C:64:MET:CE	1:C:104:GLY:HA2	2.40	0.51
1:E:129:PRO:HB3	3:E:402:HOH:O	2.10	0.51
1:B:181:GLU:OE2	1:E:96:ARG:NH1	2.43	0.51
1:A:128:SER:OG	1:A:168:ARG:NE	2.42	0.51
1:A:159:THR:HA	1:A:162:PHE:CD2	2.45	0.51
1:D:9:ASN:OD1	1:D:11:LYS:N	2.45	0.50
1:E:39:LEU:HB3	2:E:301:1Q0:H53	1.93	0.50
1:C:10:LEU:HD23	1:C:175:MET:SD	2.52	0.50
1:C:28:LEU:HB2	1:C:31:PRO:HG2	1.94	0.50
1:C:172:LEU:HA	1:C:175:MET:HE3	1.94	0.50
1:E:8:PHE:O	1:E:10:LEU:N	2.45	0.50
1:F:13:VAL:HG21	1:F:38:GLY:HA3	1.94	0.49
1:E:126:ARG:HH21	1:E:188:GLY:HA3	1.77	0.49
1:E:117:TRP:HH2	2:E:301:1Q0:H4	1.78	0.49
1:E:150:HIS:CE1	2:E:301:1Q0:H16	2.47	0.49
1:A:139:ALA:O	1:A:143:ASN:ND2	2.42	0.49
1:C:205:TYR:HB3	1:C:211:LEU:HD23	1.95	0.49
1:F:210:LEU:HD23	1:F:213:LEU:HD11	1.94	0.49
1:D:150:HIS:HB3	1:D:154:VAL:CG2	2.42	0.49
1:B:13:VAL:CG2	1:B:38:GLY:HA3	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:30:ASP:HB2	1:D:31:PRO:HD3	1.95	0.49
1:A:53:ILE:CD1	2:A:301:1Q0:H40	2.40	0.49
1:D:131:ASP:N	1:D:131:ASP:OD1	2.46	0.49
1:F:126:ARG:HD2	1:F:187:LEU:HD23	1.94	0.48
1:F:112:HIS:CE1	1:F:201:SER:HB2	2.48	0.48
1:B:43:LEU:HD22	1:B:50:PHE:HB3	1.95	0.48
1:D:101:PRO:O	1:D:106:ARG:NH1	2.46	0.48
1:D:117:TRP:CH2	2:D:301:1Q0:H66	2.49	0.47
1:D:72:GLN:HG2	1:D:85:HIS:CE1	2.49	0.47
1:F:175:MET:H	1:F:175:MET:HG2	1.54	0.47
1:C:50:PHE:O	1:C:53:ILE:HG12	2.14	0.47
1:F:177:VAL:O	1:F:179:PRO:HD2	2.13	0.47
1:F:116:HIS:O	1:F:119:GLN:HG2	2.14	0.47
1:A:36:TRP:HB3	1:A:61:LEU:HD11	1.96	0.47
1:D:129:PRO:HG2	1:D:132:ALA:HB2	1.97	0.47
1:D:50:PHE:O	1:D:53:ILE:HG12	2.14	0.47
1:F:135:SER:HB2	1:F:162:PHE:CD1	2.48	0.47
1:A:172:LEU:HD23	1:A:180:PRO:HB3	1.95	0.47
1:A:191:LEU:O	1:A:194:ILE:HG22	2.14	0.47
1:A:131:ASP:N	1:A:131:ASP:OD1	2.47	0.47
1:C:64:MET:HE1	1:C:67:LEU:HD12	1.97	0.47
1:E:12:VAL:O	1:E:16:SER:OG	2.32	0.47
1:A:166:PRO:HD2	1:A:171:PHE:CZ	2.50	0.47
2:D:301:1Q0:H44	2:D:301:1Q0:H51	1.71	0.47
1:C:41:ARG:HA	1:C:44:ASN:HB2	1.97	0.47
1:F:28:LEU:HB3	1:F:31:PRO:HD2	1.96	0.47
1:C:134:THR:HA	1:C:137:LEU:HD23	1.96	0.47
1:B:167:THR:OG1	1:B:170:VAL:HG23	2.15	0.46
1:E:126:ARG:NH2	1:E:188:GLY:HA3	2.30	0.46
1:F:134:THR:O	1:F:138:CYS:HB2	2.15	0.46
1:D:153:VAL:HG21	2:F:301:1Q0:H26	1.97	0.46
1:C:62:ARG:O	1:C:66:ARG:HG2	2.15	0.46
1:F:87:LEU:HD12	1:F:92:VAL:HG21	1.98	0.46
1:F:205:TYR:CD1	1:F:210:LEU:HD22	2.51	0.46
1:A:74:GLU:HG3	1:D:74:GLU:HB2	1.97	0.46
1:C:138:CYS:O	1:C:142:TYR:N	2.46	0.46
1:E:125:LEU:HA	1:E:137:LEU:HD13	1.97	0.46
1:D:60:LYS:CE	2:D:301:1Q0:H30	2.46	0.46
1:A:43:LEU:HG	2:A:301:1Q0:H54	1.98	0.46
1:E:123:GLU:O	1:E:126:ARG:HG3	2.16	0.46
1:C:116:HIS:HB2	1:C:198:TYR:CE1	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:42:PHE:HZ	1:D:175:MET:HA	1.81	0.46
1:D:116:HIS:HB2	1:D:198:TYR:CE1	2.51	0.46
1:C:179:PRO:C	1:C:181:GLU:H	2.20	0.45
1:D:112:HIS:CE1	1:D:201:SER:HB2	2.52	0.45
1:F:187:LEU:HG	1:F:191:LEU:HB2	1.97	0.45
1:F:137:LEU:O	1:F:141:SER:OG	2.33	0.45
1:F:103:SER:O	1:F:106:ARG:HG3	2.15	0.45
1:F:50:PHE:CE1	1:F:161:ALA:HB1	2.51	0.45
1:C:115:LEU:HA	1:C:115:LEU:HD12	1.83	0.45
1:B:43:LEU:HA	1:B:43:LEU:HD23	1.85	0.45
1:F:72:GLN:HE21	1:F:85:HIS:HE1	1.65	0.45
1:D:159:THR:O	1:D:162:PHE:HB2	2.17	0.45
2:F:301:1Q0:NBM	2:F:301:1Q0:CAJ	2.78	0.44
1:B:53:ILE:HB	2:B:301:1Q0:CAU	2.46	0.44
1:C:64:MET:HE3	1:C:104:GLY:HA2	1.99	0.44
1:F:22:ASP:HB3	1:F:26:GLU:HG3	1.99	0.44
1:E:63:ILE:HD13	1:E:102:GLU:HB3	1.98	0.44
1:C:78:SER:OG	1:C:79:LEU:N	2.49	0.44
1:F:36:TRP:HA	1:F:39:LEU:HD22	1.99	0.44
2:D:301:1Q0:H36	2:D:301:1Q0:H20	2.00	0.44
1:E:12:VAL:HA	1:E:15:VAL:HG13	1.98	0.44
1:F:106:ARG:NH2	1:F:214:PRO:OXT	2.36	0.44
1:E:52:PHE:CE1	1:E:53:ILE:HG22	2.52	0.44
1:A:124:GLY:O	1:A:128:SER:HB3	2.18	0.44
1:A:9:ASN:HB2	1:A:175:MET:O	2.18	0.44
1:C:24:LYS:HE2	1:C:24:LYS:HB3	1.63	0.44
1:F:93:ASP:CG	1:F:100:HIS:H	2.21	0.44
1:D:146:LEU:HD12	1:D:146:LEU:HA	1.83	0.44
1:E:146:LEU:HG	1:E:150:HIS:CE1	2.52	0.44
1:D:106:ARG:HB2	1:D:106:ARG:HE	1.60	0.44
1:F:49:ILE:O	1:F:52:PHE:N	2.50	0.44
1:F:169:GLU:N	1:F:169:GLU:OE1	2.48	0.44
1:A:76:TYR:OH	1:A:86:GLU:OE2	2.28	0.43
1:E:179:PRO:HB2	1:E:182:GLN:OE1	2.18	0.43
1:C:112:HIS:CE1	1:C:201:SER:HB2	2.54	0.43
1:D:28:LEU:O	1:D:31:PRO:HD2	2.18	0.43
1:B:10:LEU:O	1:B:13:VAL:N	2.48	0.43
1:E:60:LYS:NZ	2:E:301:1Q0:OAF	2.52	0.43
1:A:105:CYS:O	1:A:205:TYR:OH	2.27	0.43
1:F:27:VAL:HG13	1:F:79:LEU:HG	2.00	0.43
1:C:80:GLN:HG3	1:C:204:LEU:HD22	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:114:ALA:HB1	2:D:301:1Q0:H40	2.00	0.43
1:F:82:MET:O	1:F:85:HIS:N	2.51	0.43
1:C:78:SER:OG	1:C:80:GLN:OE1	2.30	0.43
1:E:56:ASP:O	1:E:60:LYS:HG2	2.18	0.43
1:D:117:TRP:CH2	2:D:301:1Q0:CAW	3.02	0.43
1:F:72:GLN:HG2	1:F:72:GLN:H	1.68	0.43
1:C:96:ARG:O	1:C:98:SER:N	2.52	0.43
1:C:43:LEU:HG	2:C:301:1Q0:H54	2.01	0.43
1:C:191:LEU:O	1:C:195:GLN:HB2	2.19	0.42
1:A:93:ASP:HB3	1:A:96:ARG:O	2.19	0.42
1:C:83:VAL:O	1:C:87:LEU:HB2	2.19	0.42
1:A:167:THR:OG1	1:A:170:VAL:HG23	2.19	0.42
1:D:60:LYS:HE3	1:D:110:ARG:HG2	2.02	0.42
1:C:24:LYS:H	1:C:24:LYS:HG2	1.48	0.42
1:C:124:GLY:O	1:C:128:SER:HB3	2.20	0.42
1:C:62:ARG:HG3	1:C:62:ARG:HH11	1.85	0.42
1:F:195:GLN:HA	1:F:198:TYR:HB3	2.02	0.42
1:F:109:LEU:HD22	1:F:213:LEU:HD13	2.02	0.42
1:F:179:PRO:HA	1:F:180:PRO:HD2	1.87	0.42
1:E:8:PHE:O	1:E:10:LEU:HD22	2.19	0.42
1:C:36:TRP:CZ2	1:C:115:LEU:HD13	2.55	0.42
1:F:124:GLY:O	1:F:128:SER:HB3	2.19	0.42
1:D:166:PRO:HD2	1:D:171:PHE:HE1	1.85	0.42
1:D:143:ASN:HA	1:D:147:ALA:HB3	2.02	0.42
1:B:17:PHE:CD2	1:B:194:ILE:HG12	2.54	0.42
1:F:90:ARG:HA	1:F:90:ARG:HD3	1.72	0.42
1:C:119:GLN:O	1:C:123:GLU:HB3	2.20	0.42
2:E:301:1Q0:H42	2:E:301:1Q0:H50	2.02	0.41
1:C:116:HIS:O	1:C:120:LEU:HB2	2.19	0.41
1:C:90:ARG:HD3	1:E:208:HIS:O	2.20	0.41
1:C:64:MET:HE1	1:C:104:GLY:HA2	2.02	0.41
1:E:17:PHE:HA	1:E:17:PHE:HD1	1.73	0.41
1:D:182:GLN:HA	1:D:185:GLN:HB2	2.01	0.41
1:A:98:SER:OG	1:A:99:HIS:N	2.53	0.41
1:A:13:VAL:O	1:A:16:SER:OG	2.25	0.41
1:F:187:LEU:HD21	1:F:191:LEU:HD22	2.01	0.41
1:D:128:SER:O	1:D:168:ARG:NH2	2.53	0.41
1:F:79:LEU:O	1:F:83:VAL:HG23	2.21	0.41
1:D:191:LEU:N	1:D:192:PRO:HD2	2.35	0.41
1:E:59:SER:O	1:E:62:ARG:HB2	2.21	0.41
1:F:135:SER:OG	1:F:136:ALA:N	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:190:ALA:O	1:E:194:ILE:HG13	2.20	0.41
1:A:76:TYR:CE1	1:A:82:MET:HG3	2.54	0.41
1:C:106:ARG:HG3	1:C:107:THR:N	2.34	0.41
1:F:31:PRO:HA	1:F:34:ALA:HB3	2.02	0.41
1:C:142:TYR:O	1:C:145:SER:N	2.39	0.41
1:F:117:TRP:NE1	1:F:141:SER:HB2	2.35	0.41
1:E:58:VAL:O	1:E:62:ARG:HG2	2.21	0.41
1:E:177:VAL:HG11	1:E:186:MET:SD	2.60	0.41
1:A:30:ASP:HB2	1:A:31:PRO:HD3	2.01	0.41
1:E:71:PRO:HB2	1:E:72:GLN:OE1	2.20	0.41
1:F:108:VAL:HA	1:F:111:LEU:HD12	2.02	0.41
1:B:53:ILE:HG13	1:B:54:SER:N	2.35	0.41
1:D:22:ASP:OD1	1:D:26:GLU:N	2.45	0.41
1:E:87:LEU:HD23	1:E:87:LEU:HA	1.82	0.41
1:F:204:LEU:O	1:F:208:HIS:ND1	2.53	0.41
1:F:175:MET:O	1:F:176:ASN:HB2	2.19	0.40
1:F:164:THR:OG1	1:F:164:THR:O	2.30	0.40
1:F:11:LYS:O	1:F:15:VAL:HG13	2.21	0.40
1:C:76:TYR:CE1	1:C:82:MET:HG3	2.57	0.40
1:D:11:LYS:O	1:D:15:VAL:HG23	2.21	0.40
1:B:138:CYS:HB3	1:B:162:PHE:CE2	2.56	0.40
1:A:177:VAL:HG21	1:A:183:ALA:HA	2.04	0.40
1:C:49:ILE:HG12	1:C:49:ILE:H	1.58	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	205/215 (95%)	195 (95%)	9 (4%)	1 (0%)	34 72
1	B	204/215 (95%)	188 (92%)	15 (7%)	1 (0%)	34 72

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	204/215 (95%)	174 (85%)	23 (11%)	7 (3%)	5	25
1	D	204/215 (95%)	190 (93%)	13 (6%)	1 (0%)	34	72
1	E	205/215 (95%)	188 (92%)	14 (7%)	3 (2%)	13	46
1	F	203/215 (94%)	180 (89%)	22 (11%)	1 (0%)	34	72
All	All	1225/1290 (95%)	1115 (91%)	96 (8%)	14 (1%)	17	55

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	9	ASN
1	B	66	ARG
1	C	143	ASN
1	E	166	PRO
1	C	97	ARG
1	C	186	MET
1	D	98	SER
1	C	17	PHE
1	C	166	PRO
1	C	188	GLY
1	F	98	SER
1	A	170	VAL
1	C	180	PRO
1	E	194	ILE

### 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	182/190 (96%)	168 (92%)	14 (8%)	16	50
1	B	180/190 (95%)	164 (91%)	16 (9%)	12	42
1	C	178/190 (94%)	153 (86%)	25 (14%)	4	18
1	D	181/190 (95%)	166 (92%)	15 (8%)	14	46
1	E	180/190 (95%)	161 (89%)	19 (11%)	8	31

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	F	175/190 (92%)	146 (83%)	29 (17%)	3	12
All	All	1076/1140 (94%)	958 (89%)	118 (11%)	8	30

All (118) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	LEU
1	A	43	LEU
1	A	72	GLN
1	A	73	SER
1	A	95	GLU
1	A	103	SER
1	A	106	ARG
1	A	146	LEU
1	A	152	TRP
1	A	164	THR
1	A	172	LEU
1	A	173	GLU
1	A	185	GLN
1	A	186	MET
1	B	10	LEU
1	B	13	VAL
1	B	14	LEU
1	B	23	GLU
1	B	43	LEU
1	B	53	ILE
1	B	62	ARG
1	B	68	ARG
1	B	97	ARG
1	B	106	ARG
1	B	115	LEU
1	B	120	LEU
1	B	134	THR
1	B	164	THR
1	B	181	GLU
1	B	182	GLN
1	C	14	LEU
1	C	23	GLU
1	C	24	LYS
1	C	26	GLU
1	C	30	ASP
1	C	43	LEU

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Mol	Chain	Res	Type
1	C	49	ILE
1	C	72	GLN
1	C	78	SER
1	C	106	ARG
1	C	115	LEU
1	C	120	LEU
1	C	134	THR
1	C	137	LEU
1	C	138	CYS
1	C	146	LEU
1	C	153	VAL
1	C	164	THR
1	C	167	THR
1	C	171	PHE
1	C	182	GLN
1	C	191	LEU
1	C	195	GLN
1	C	197	VAL
1	C	200	VAL
1	D	45	SER
1	D	48	THR
1	D	73	SER
1	D	95	GLU
1	D	97	ARG
1	D	103	SER
1	D	106	ARG
1	D	128	SER
1	D	131	ASP
1	D	146	LEU
1	D	159	THR
1	D	170	VAL
1	D	171	PHE
1	D	189	GLU
1	D	203	LYS
1	E	10	LEU
1	E	15	VAL
1	E	16	SER
1	E	17	PHE
1	E	21	LEU
1	E	23	GLU
1	E	29	LEU
1	E	51	SER

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Mol	Chain	Res	Type
1	E	53	ILE
1	E	78	SER
1	E	95	GLU
1	E	100	HIS
1	E	106	ARG
1	E	119	GLN
1	E	128	SER
1	E	156	ARG
1	E	164	THR
1	E	186	MET
1	E	191	LEU
1	F	12	VAL
1	F	13	VAL
1	F	15	VAL
1	F	23	GLU
1	F	27	VAL
1	F	39	LEU
1	F	46	LEU
1	F	49	ILE
1	F	53	ILE
1	F	72	GLN
1	F	74	GLU
1	F	89	ASN
1	F	90	ARG
1	F	95	GLU
1	F	97	ARG
1	F	98	SER
1	F	103	SER
1	F	106	ARG
1	F	110	ARG
1	F	138	CYS
1	F	141	SER
1	F	153	VAL
1	F	154	VAL
1	F	175	MET
1	F	184	VAL
1	F	187	LEU
1	F	196	ARG
1	F	201	SER
1	F	205	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	112	HIS
1	C	185	GLN
1	D	89	ASN
1	E	89	ASN
1	F	85	HIS
1	F	116	HIS
1	F	195	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 carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

6 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	1Q0	A	301	-	42,43,43	0.82	2 (4%)	43,48,48	0.81	1 (2%)
2	1Q0	B	301	-	25,26,43	1.07	2 (8%)	25,31,48	1.09	2 (8%)
2	1Q0	C	301	-	42,43,43	0.82	2 (4%)	43,48,48	0.80	1 (2%)
2	1Q0	D	301	-	42,43,43	0.97	2 (4%)	43,48,48	0.88	1 (2%)
2	1Q0	E	301	-	42,43,43	0.83	2 (4%)	43,48,48	0.80	1 (2%)
2	1Q0	F	301	-	14,15,43	1.04	2 (14%)	16,20,48	1.33	3 (18%)

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	1Q0	A	301	-	-	0/46/46/46	0/0/0/0
2	1Q0	B	301	-	-	0/29/29/46	0/0/0/0
2	1Q0	C	301	-	-	0/46/46/46	0/0/0/0
2	1Q0	D	301	-	-	0/46/46/46	0/0/0/0
2	1Q0	E	301	-	-	0/46/46/46	0/0/0/0
2	1Q0	F	301	-	-	0/17/17/46	0/0/0/0

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	301	1Q0	CBQ-NBM	-3.25	1.40	1.46
2	E	301	1Q0	CBQ-NBM	-2.17	1.42	1.46
2	A	301	1Q0	CBQ-NBM	-2.17	1.42	1.46
2	B	301	1Q0	CBQ-NBM	-2.15	1.42	1.46
2	C	301	1Q0	CBQ-NBM	-2.11	1.42	1.46
2	F	301	1Q0	CBQ-NBM	-2.09	1.42	1.46
2	F	301	1Q0	CBO-NBM	2.08	1.42	1.34
2	D	301	1Q0	CBO-NBM	3.27	1.40	1.34
2	C	301	1Q0	CBO-NBM	4.12	1.42	1.34
2	A	301	1Q0	CBO-NBM	4.14	1.42	1.34
2	E	301	1Q0	CBO-NBM	4.15	1.42	1.34
2	B	301	1Q0	CBO-NBM	4.15	1.42	1.34

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	301	1Q0	CBP-CAK-CAJ	-3.33	119.99	124.85
2	F	301	1Q0	CBQ-NBM-CBO	-2.38	119.99	123.04
2	F	301	1Q0	CBK-CBO-NBM	2.03	119.99	116.11
2	B	301	1Q0	OBK-CBL-CBQ	2.20	111.92	108.33
2	D	301	1Q0	CBK-CBO-NBM	2.24	119.49	115.83
2	B	301	1Q0	CBK-CBO-NBM	2.54	119.98	115.83
2	E	301	1Q0	CBK-CBO-NBM	2.55	119.99	115.83
2	C	301	1Q0	CBK-CBO-NBM	2.55	119.99	115.83
2	A	301	1Q0	CBK-CBO-NBM	2.57	120.02	115.83

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 38 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	1Q0	9	0
2	B	301	1Q0	5	0
2	C	301	1Q0	1	0
2	D	301	1Q0	9	0
2	E	301	1Q0	11	0
2	F	301	1Q0	3	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	207/215 (96%)	-0.48	0 100 100	37, 63, 107, 132	0
1	B	206/215 (95%)	-0.34	1 (0%) 91 83	44, 79, 118, 142	0
1	C	206/215 (95%)	-0.27	0 100 100	49, 95, 140, 191	0
1	D	206/215 (95%)	-0.30	1 (0%) 91 83	47, 86, 130, 207	0
1	E	207/215 (96%)	-0.40	0 100 100	44, 84, 129, 174	0
1	F	205/215 (95%)	0.34	12 (5%) 26 11	89, 149, 195, 226	0
All	All	1237/1290 (95%)	-0.24	14 (1%) 82 66	37, 89, 165, 226	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	185	GLN	3.6
1	F	26	GLU	3.1
1	F	71	PRO	2.8
1	F	91	LEU	2.7
1	F	132	ALA	2.6
1	B	62	ARG	2.5
1	F	97	ARG	2.3
1	F	177	VAL	2.2
1	F	170	VAL	2.2
1	F	129	PRO	2.2
1	D	214	PRO	2.2
1	F	209	SER	2.1
1	F	95	GLU	2.1
1	F	102	GLU	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 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, 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	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	1Q0	A	301	44/44	0.93	0.24	1.68	66,66,68,69	0
2	1Q0	C	301	44/44	0.91	0.25	1.64	71,77,85,88	0
2	1Q0	E	301	44/44	0.93	0.24	1.29	69,72,79,81	0
2	1Q0	D	301	44/44	0.93	0.23	0.60	82,85,88,89	0
2	1Q0	B	301	27/44	0.94	0.22	0.45	49,57,61,62	27
2	1Q0	F	301	16/44	0.92	0.22	-0.14	94,96,98,100	16

### 6.5 Other polymers [i](#)

There are no such residues in this entry.