



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 03:40 am BST

PDB ID : 4QGU
Title : protein domain complex with ssDNA
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Deposited on : 2014-05-25
Resolution : 2.54 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
Xtriage (Phenix)	:	1.13
EDS	:	2.11
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)	:	2.11

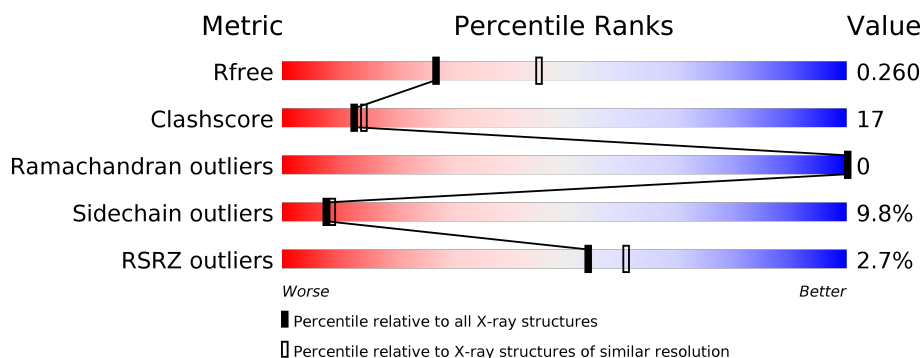
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.54 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	1284 (2.56-2.52)
Clashscore	141614	1332 (2.56-2.52)
Ramachandran outliers	138981	1315 (2.56-2.52)
Sidechain outliers	138945	1315 (2.56-2.52)
RSRZ outliers	127900	1272 (2.56-2.52)

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	206	<div> <div>76%</div> <div>14%</div> <div>7%</div> </div>
1	B	206	<div>2%</div> <div>61%</div> <div>25%</div> <div>6%</div> <div>7%</div>
2	C	12	<div>33%</div> <div>8%</div> <div>33%</div> <div>58%</div>
2	D	12	<div>8%</div> <div>33%</div> <div>50%</div> <div>17%</div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3608 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 Gamma-interferon-inducible protein 16.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	191	Total	C	N	O	S	4	0	0
			1547	989	262	287	9			
1	B	191	Total	C	N	O	S	5	0	0
			1547	989	262	287	9			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	188	GLY	-	EXPRESSION TAG	UNP Q16666
A	189	SER	-	EXPRESSION TAG	UNP Q16666
A	190	HIS	-	EXPRESSION TAG	UNP Q16666
A	191	MET	-	EXPRESSION TAG	UNP Q16666
B	188	GLY	-	EXPRESSION TAG	UNP Q16666
B	189	SER	-	EXPRESSION TAG	UNP Q16666
B	190	HIS	-	EXPRESSION TAG	UNP Q16666
B	191	MET	-	EXPRESSION TAG	UNP Q16666

- Molecule 2 is a DNA chain called DNA (5'-D(P*AP*GP*GP*CP*CP*GP*GP*CP*GP*TP*GP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	12	Total	C	N	O	P	0	0	0
			251	117	51	71	12			
2	D	10	Total	C	N	O	P	0	0	0
			204	97	41	56	10			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	33	Total	O	0	0
			33	33		

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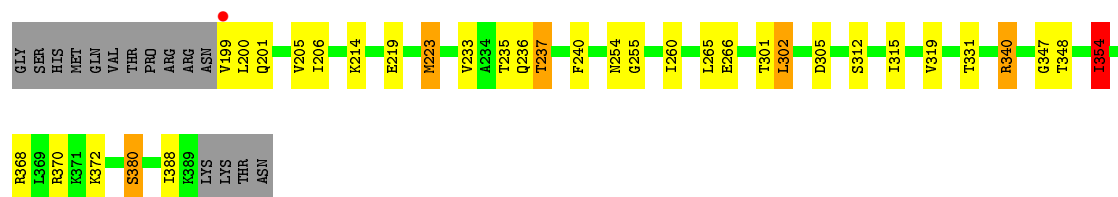
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	18	Total 18	O 18	0	0
3	C	4	Total 4	O 4	0	0
3	D	4	Total 4	O 4	0	0

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

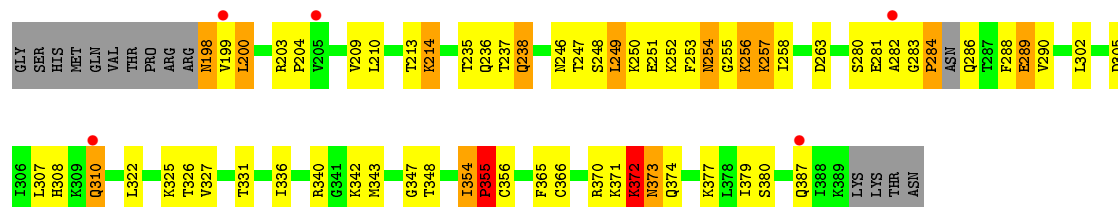
• Molecule 1: Gamma-interferon-inducible protein 16

Chain A: 



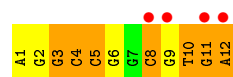
• Molecule 1: Gamma-interferon-inducible protein 16

Chain B: 



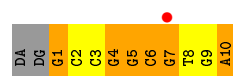
• Molecule 2: DNA (5'-D(P*AP*GP*GP*CP*CP*GP*GP*CP*GP*TP*GP*A)-3')

Chain C: 



• Molecule 2: DNA (5'-D(P*AP*GP*GP*CP*CP*GP*GP*CP*GP*TP*GP*A)-3')

Chain D: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	85.72Å 91.86Å 69.98Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.40 – 2.54 38.40 – 2.55	Depositor EDS
% Data completeness (in resolution range)	96.5 (38.40-2.54) 92.6 (38.40-2.55)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.14 (at 2.54Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.238 , 0.258 0.240 , 0.260	Depositor DCC
R_{free} test set	898 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	55.3	Xtriage
Anisotropy	0.485	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 39.0	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	0.92	EDS
Total number of atoms	3608	wwPDB-VP
Average B, all atoms (Å ²)	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.69% 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

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	0.51	0/1573	0.66	1/2113 (0.0%)
1	B	0.51	1/1572 (0.1%)	0.62	2/2110 (0.1%)
2	C	0.65	0/282	2.17	14/434 (3.2%)
2	D	0.63	0/229	1.60	8/347 (2.3%)
All	All	0.53	1/3656 (0.0%)	0.97	25/5004 (0.5%)

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

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	355	PRO	N-CD	5.34	1.55	1.47

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	11	DG	O4'-C1'-N9	16.12	119.29	108.00
2	C	3	DG	O4'-C1'-N9	14.08	117.86	108.00
2	C	6	DG	O4'-C1'-N9	9.93	114.95	108.00
2	C	5	DC	O4'-C1'-N1	9.84	114.89	108.00
2	C	10	DT	O4'-C1'-N1	9.47	114.63	108.00
2	D	6	DC	O4'-C1'-N1	9.05	114.33	108.00
2	C	1	DA	O4'-C1'-N9	8.80	114.16	108.00
2	D	4	DG	O4'-C1'-N9	-8.78	101.86	108.00
2	C	12	DA	C1'-O4'-C4'	-7.54	102.56	110.10
2	D	5	DG	O4'-C1'-N9	7.00	112.90	108.00
2	D	4	DG	O4'-C4'-C3'	-6.93	101.73	104.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	4	DC	C1'-O4'-C4'	-6.72	103.38	110.10
2	C	8	DC	O4'-C1'-N1	-6.69	103.31	108.00
2	D	1	DG	O4'-C1'-N9	-6.22	103.64	108.00
1	A	354	ILE	C-N-CD	5.96	140.92	128.40
2	D	10	DA	C1'-O4'-C4'	-5.79	104.31	110.10
1	B	354	ILE	C-N-CD	5.78	140.53	128.40
2	C	3	DG	C3'-C2'-C1'	-5.77	95.58	102.50
2	D	7	DG	O4'-C1'-N9	-5.54	104.12	108.00
2	C	1	DA	N9-C1'-C2'	5.48	123.01	112.60
2	C	3	DG	C1'-O4'-C4'	-5.46	104.64	110.10
1	B	284	PRO	N-CA-C	5.38	126.08	112.10
2	C	6	DG	C3'-C2'-C1'	-5.26	96.19	102.50
2	D	4	DG	C4'-C3'-C2'	-5.23	98.39	103.10
2	C	10	DT	C5-C4-O4	-5.11	121.33	124.90

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	372	LYS	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	1547	0	1596	23	0
1	B	1547	0	1595	70	1
2	C	251	0	134	13	1
2	D	204	0	112	17	0
3	A	33	0	0	0	0
3	B	18	0	0	1	0
3	C	4	0	0	0	0
3	D	4	0	0	2	0
All	All	3608	0	3437	114	1

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 17.

All (114) 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:A:305:ASP:CG	1:A:340:ARG:NH1	1.86	1.29
2:C:5:DC:N4	2:D:8:DT:O4	1.66	1.27
1:A:305:ASP:OD2	1:A:340:ARG:NH1	1.70	1.24
1:A:305:ASP:CG	1:A:340:ARG:HH12	1.43	1.21
1:A:305:ASP:OD1	1:A:340:ARG:NH1	1.79	1.14
1:B:256:LYS:NZ	1:B:283:GLY:O	1.88	1.07
1:B:284:PRO:C	1:B:286:GLN:N	2.13	1.01
1:B:256:LYS:HD3	1:B:281:GLU:HB3	1.42	1.00
1:B:256:LYS:CE	1:B:283:GLY:O	2.11	0.98
2:C:3:DG:O6	2:D:10:DA:N6	1.99	0.95
2:C:4:DC:N4	2:D:9:DG:O6	1.98	0.94
1:B:373:ASN:ND2	1:B:373:ASN:O	2.03	0.91
1:B:236:GLN:NE2	1:B:289:GLU:O	2.10	0.84
1:A:305:ASP:CG	1:A:340:ARG:HH11	1.79	0.82
1:B:256:LYS:HE3	1:B:258:ILE:CG2	2.10	0.81
1:B:200:LEU:O	1:B:200:LEU:HD23	1.81	0.79
1:B:235:THR:HG23	1:B:237:THR:H	1.53	0.74
1:B:198:ASN:O	1:B:198:ASN:ND2	2.22	0.73
1:A:201:GLN:NE2	1:A:266:GLU:OE1	2.14	0.73
1:B:256:LYS:HE2	1:B:283:GLY:O	1.89	0.71
1:B:235:THR:HG22	1:B:238:GLN:H	1.56	0.70
2:C:8:DC:N4	2:D:5:DG:O6	2.18	0.69
1:B:235:THR:OG1	3:B:417:HOH:O	2.10	0.68
1:B:377:LYS:HD3	1:B:379:ILE:HD11	1.76	0.67
2:D:9:DG:H2'	2:D:10:DA:H5'	1.77	0.66
1:B:213:THR:O	1:B:250:LYS:NZ	2.29	0.64
1:B:373:ASN:O	1:B:374:GLN:HB2	1.98	0.64
2:D:9:DG:H5''	2:D:9:DG:H8	1.63	0.63
1:B:256:LYS:CD	1:B:281:GLU:HB3	2.24	0.63
2:D:3:DC:H2'	2:D:4:DG:C8	2.33	0.63
2:D:8:DT:H2''	2:D:9:DG:OP1	1.98	0.63
1:A:235:THR:OG1	1:A:236:GLN:N	2.32	0.62
1:B:256:LYS:CE	1:B:283:GLY:H	2.12	0.62
1:B:282:ALA:O	1:B:286:GLN:NE2	2.33	0.61
2:C:5:DC:C4	2:D:8:DT:O4	2.51	0.60
1:A:331:THR:HG22	1:A:348:THR:HG22	1.81	0.60
1:B:256:LYS:HE3	1:B:258:ILE:HG22	1.82	0.60
1:B:235:THR:CG2	1:B:238:GLN:H	2.15	0.60
1:B:373:ASN:H	1:B:373:ASN:ND2	1.98	0.60
1:B:331:THR:HG22	1:B:348:THR:HG22	1.83	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:322:LEU:HD22	1:B:356:CYS:SG	2.42	0.59
1:B:254:ASN:HD22	1:B:254:ASN:H	1.48	0.59
1:B:310:GLN:O	1:B:371:LYS:HE2	2.03	0.59
2:D:9:DG:H2'	2:D:10:DA:C5'	2.33	0.58
1:B:246:ASN:O	1:B:249:LEU:HD12	2.04	0.58
1:B:256:LYS:HZ2	1:B:256:LYS:HB2	1.68	0.57
2:C:10:DT:H2''	2:C:11:DG:O5'	2.05	0.57
1:B:256:LYS:CE	1:B:258:ILE:HG22	2.35	0.57
1:A:206:ILE:HG12	1:A:260:ILE:HG12	1.87	0.57
1:A:340:ARG:HB3	1:A:340:ARG:HH11	1.70	0.56
1:B:214:LYS:HE3	1:B:387:GLN:HE22	1.71	0.56
1:B:235:THR:HG23	1:B:237:THR:N	2.19	0.56
1:B:355:PRO:HG2	1:B:355:PRO:O	2.06	0.56
1:B:209:VAL:HG11	1:B:253:PHE:O	2.08	0.54
1:B:254:ASN:H	1:B:254:ASN:ND2	2.05	0.54
2:C:12:DA:H1'	2:D:2:DC:C5	2.43	0.54
1:A:347:GLY:HA2	1:A:380:SER:HB3	1.90	0.53
1:B:256:LYS:NZ	1:B:256:LYS:HB2	2.23	0.53
1:B:200:LEU:C	1:B:200:LEU:HD23	2.26	0.53
1:B:256:LYS:HE3	1:B:283:GLY:H	1.73	0.52
1:B:256:LYS:NZ	1:B:256:LYS:CB	2.73	0.52
1:B:256:LYS:HE2	1:B:283:GLY:H	1.73	0.52
2:D:3:DC:H2''	2:D:4:DG:O5'	2.09	0.52
1:B:210:LEU:HD21	1:B:290:VAL:HG11	1.92	0.52
1:B:365:PHE:CD1	1:B:365:PHE:N	2.77	0.51
2:D:8:DT:P	3:D:101:HOH:O	2.68	0.51
1:B:258:ILE:HD11	1:B:288:PHE:CE2	2.46	0.51
1:B:373:ASN:C	1:B:373:ASN:HD22	2.03	0.51
1:B:256:LYS:HE3	1:B:258:ILE:HG23	1.92	0.50
1:B:209:VAL:O	1:B:257:LYS:HG3	2.11	0.50
1:B:254:ASN:HD22	1:B:254:ASN:N	2.08	0.50
1:B:254:ASN:N	1:B:254:ASN:ND2	2.60	0.50
1:B:325:LYS:HD3	1:B:327:VAL:HG23	1.95	0.49
1:B:256:LYS:NZ	1:B:258:ILE:HG22	2.27	0.49
1:A:199:VAL:HG13	1:A:200:LEU:H	1.78	0.48
2:D:2:DC:H2''	2:D:3:DC:O4'	2.14	0.48
1:B:258:ILE:O	1:B:258:ILE:HG23	2.14	0.48
1:B:254:ASN:HD22	1:B:255:GLY:H	1.61	0.47
1:B:284:PRO:CA	1:B:286:GLN:N	2.78	0.47
1:B:256:LYS:HZ2	1:B:258:ILE:HG22	1.80	0.46
2:D:1:DG:H5'	3:D:104:HOH:O	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:235:THR:HG23	1:A:237:THR:H	1.81	0.46
1:A:302:LEU:H	1:A:302:LEU:HD12	1.80	0.46
1:A:301:THR:HG21	1:A:319:VAL:O	2.14	0.46
1:B:336:ILE:HD11	1:B:343:MET:HE3	1.96	0.46
1:B:325:LYS:HD3	1:B:327:VAL:CG2	2.46	0.46
1:B:256:LYS:HE2	1:B:283:GLY:C	2.36	0.46
2:D:6:DC:H2''	2:D:7:DG:C8	2.51	0.45
1:B:256:LYS:HG3	1:B:258:ILE:O	2.16	0.45
2:C:9:DG:H2'	2:C:10:DT:H71	1.98	0.44
1:A:331:THR:HG22	1:A:348:THR:CG2	2.47	0.44
1:B:256:LYS:HE2	1:B:283:GLY:N	2.32	0.44
2:C:3:DG:O6	2:D:10:DA:C6	2.68	0.43
1:B:280:SER:OG	1:B:281:GLU:N	2.52	0.43
2:C:10:DT:H2''	2:C:11:DG:C5'	2.48	0.43
1:A:219:GLU:OE2	1:A:223:MET:HG3	2.18	0.43
1:B:373:ASN:ND2	1:B:373:ASN:N	2.60	0.43
1:A:233:VAL:HG22	1:A:240:PHE:HB2	1.99	0.43
1:A:199:VAL:HG22	1:A:265:LEU:HD22	2.01	0.42
1:B:203:ARG:HD2	1:B:204:PRO:HD2	2.00	0.42
1:A:235:THR:HG23	1:A:237:THR:N	2.35	0.42
1:A:354:ILE:CG2	1:A:388:ILE:HD11	2.49	0.42
1:B:347:GLY:HA2	1:B:380:SER:HB3	2.02	0.42
1:A:315:ILE:HD11	1:A:368:ARG:CZ	2.50	0.41
1:B:305:ASP:OD2	1:B:340:ARG:NH1	2.54	0.41
2:C:2:DG:H2''	2:C:3:DG:OP1	2.20	0.41
1:B:307:LEU:HA	1:B:307:LEU:HD23	1.93	0.41
1:B:308:HIS:O	1:B:371:LYS:NZ	2.54	0.41
1:A:254:ASN:OD1	1:A:255:GLY:N	2.53	0.41
1:B:237:THR:OG1	1:B:238:GLN:HG3	2.21	0.41
1:B:373:ASN:HA	2:C:2:DG:OP2	2.21	0.41
1:B:354:ILE:HG23	1:B:355:PRO:HD2	2.02	0.40
1:B:374:GLN:N	2:C:2:DG:OP2	2.48	0.40
1:B:282:ALA:C	1:B:286:GLN:HE21	2.24	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:372:LYS:NZ	2:C:2:DG:O6[2_655]	1.36	0.84

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	189/206 (92%)	187 (99%)	2 (1%)	0	100	100
1	B	187/206 (91%)	185 (99%)	2 (1%)	0	100	100
All	All	376/412 (91%)	372 (99%)	4 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	178/192 (93%)	167 (94%)	11 (6%)	18	24
1	B	178/192 (93%)	154 (86%)	24 (14%)	4	3
All	All	356/384 (93%)	321 (90%)	35 (10%)	8	9

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	205	VAL
1	A	214	LYS
1	A	223	MET
1	A	237	THR
1	A	302	LEU
1	A	312	SER
1	A	340	ARG
1	A	354	ILE

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Mol	Chain	Res	Type
1	A	370	ARG
1	A	372	LYS
1	A	380	SER
1	B	198	ASN
1	B	199	VAL
1	B	200	LEU
1	B	214	LYS
1	B	238	GLN
1	B	247	THR
1	B	248	SER
1	B	249	LEU
1	B	251	GLU
1	B	252	LYS
1	B	254	ASN
1	B	256	LYS
1	B	257	LYS
1	B	263	ASP
1	B	289	GLU
1	B	302	LEU
1	B	310	GLN
1	B	326	THR
1	B	342	LYS
1	B	355	PRO
1	B	366	CYS
1	B	370	ARG
1	B	372	LYS
1	B	373	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	373	ASN
1	B	254	ASN
1	B	286	GLN
1	B	373	ASN

5.3.3 RNA ⓘ

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	191/206 (92%)	-0.01	1 (0%) 91 94	34, 61, 99, 115	4 (2%)
1	B	191/206 (92%)	0.23	5 (2%) 56 62	42, 79, 116, 130	3 (1%)
2	C	12/12 (100%)	1.76	4 (33%) 0 0	76, 140, 173, 175	0
2	D	10/12 (83%)	1.12	1 (10%) 7 9	112, 122, 147, 157	0
All	All	404/436 (92%)	0.19	11 (2%) 54 61	34, 72, 119, 175	7 (1%)

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	205	VAL	3.8
2	C	9	DG	3.6
2	C	8	DC	3.3
1	B	282	ALA	3.3
2	C	11	DG	2.9
2	C	12	DA	2.8
1	B	310	GLN	2.3
2	D	7	DG	2.2
1	B	387	GLN	2.2
1	A	199	VAL	2.1
1	B	199	VAL	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 carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.