



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 01:20 am BST

PDB ID : 5E1T
Title : Crystal structure of TRAF1 TRAF domain
Authors : Park, H.H.; Kim, C.M.
Deposited on : 2015-09-30
Resolution : 2.80 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

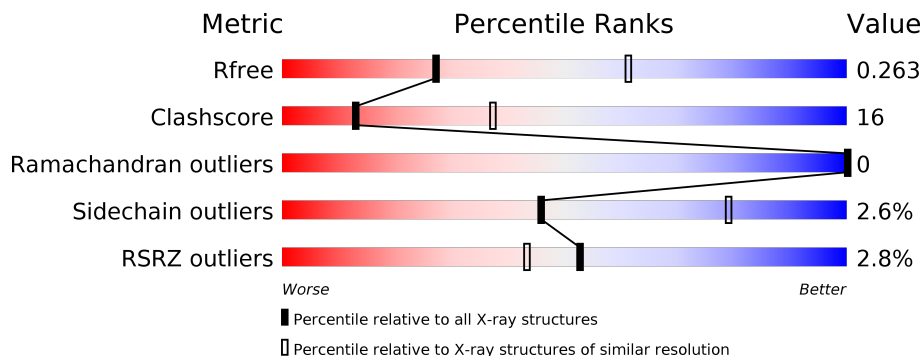
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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	197	<div style="display: flex; align-items: center;"> <div style="width: 2%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 66%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 30%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: grey;"></div> </div> <p style="font-size: small; margin-top: 5px;">2% 66% 30% ••</p>
1	B	197	<div style="display: flex; align-items: center;"> <div style="width: 4%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 61%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 32%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: grey;"></div> </div> <p style="font-size: small; margin-top: 5px;">4% 61% 32% ••</p>
1	C	197	<div style="display: flex; align-items: center;"> <div style="width: 3%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 74%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 18%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: grey;"></div> </div> <p style="font-size: small; margin-top: 5px;">3% 74% 18% •• 6%</p>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4612 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 TNF receptor-associated factor 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	192	1549	986	271	283	9	0	0	0
1	B	191	1545	984	269	283	9	0	0	0
1	C	186	1505	961	262	273	9	0	0	0

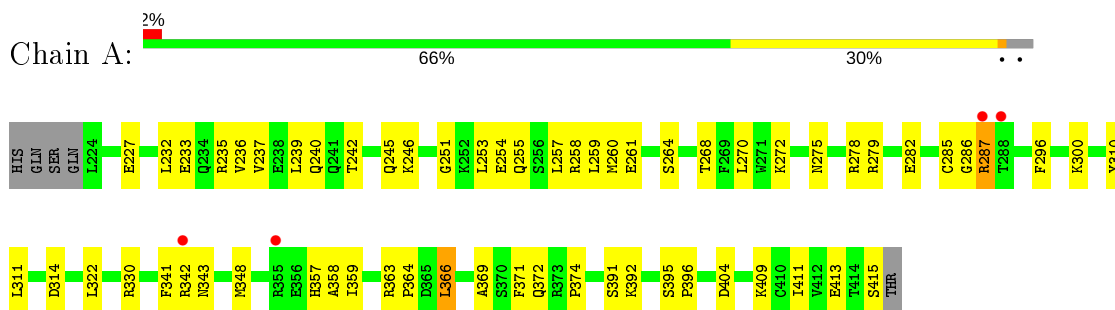
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	6	Total 6	O 6	0	0
2	B	3	Total 3	O 3	0	0
2	C	4	Total 4	O 4	0	0

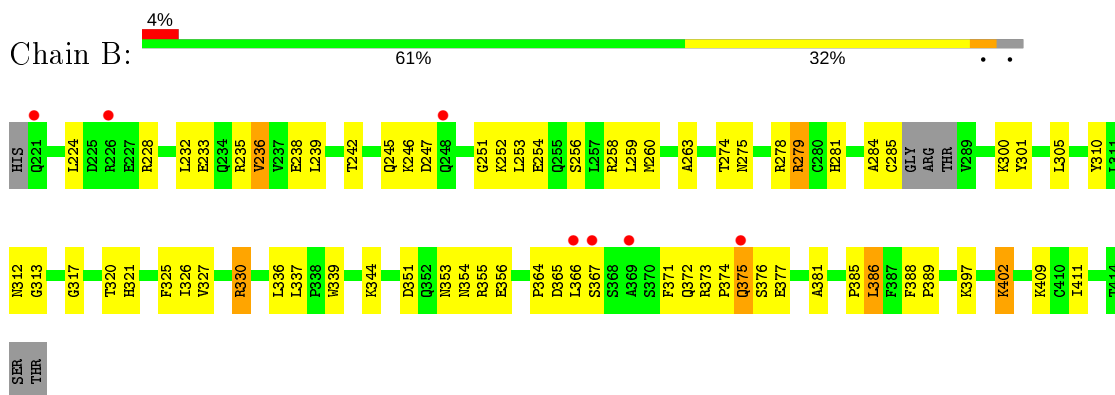
3 Residue-property plots [i](#)

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.

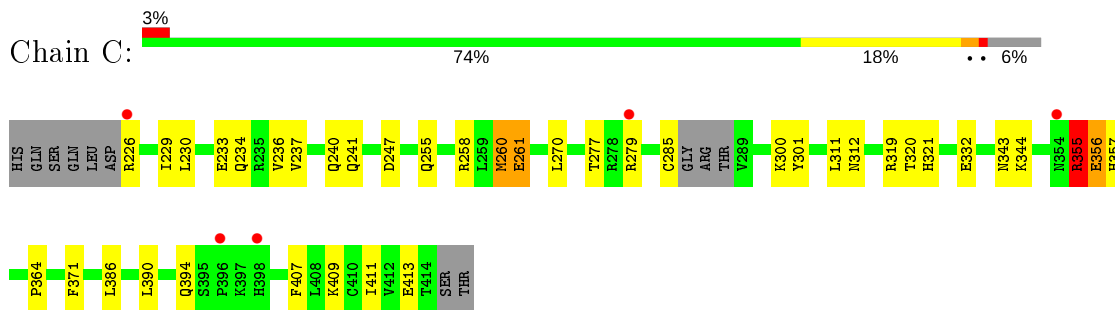
- Molecule 1: TNF receptor-associated factor 1



- Molecule 1: TNF receptor-associated factor 1



- Molecule 1: TNF receptor-associated factor 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	75.35Å 79.40Å 108.98Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.52 – 2.80 29.52 – 2.38	Depositor EDS
% Data completeness (in resolution range)	94.9 (29.52-2.80) 91.0 (29.52-2.38)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.00 (at 2.39Å)	Xtrriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.211 , 0.262 0.210 , 0.263	Depositor DCC
R_{free} test set	2019 reflections (8.26%)	wwPDB-VP
Wilson B-factor (Å ²)	43.7	Xtrriage
Anisotropy	0.455	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 53.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	0.016 for k,h,-l	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4612	wwPDB-VP
Average B, all atoms (Å ²)	76.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.30% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.55	0/1582	0.86	5/2135 (0.2%)
1	B	0.53	0/1577	0.89	5/2127 (0.2%)
1	C	0.44	0/1537	0.65	1/2073 (0.0%)
All	All	0.51	0/4696	0.81	11/6335 (0.2%)

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 (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	386	LEU	CB-CG-CD2	-12.38	89.95	111.00
1	A	287	ARG	NE-CZ-NH1	8.48	124.54	120.30
1	B	236	VAL	CG1-CB-CG2	-8.35	97.54	110.90
1	A	287	ARG	NE-CZ-NH2	-7.88	116.36	120.30
1	A	227	GLU	CA-CB-CG	7.02	128.85	113.40
1	A	366	LEU	CB-CG-CD1	6.39	121.87	111.00
1	C	260	MET	CB-CG-SD	-6.35	93.36	112.40
1	B	402	LYS	CD-CE-NZ	-6.21	97.43	111.70
1	B	344	LYS	CD-CE-NZ	-5.97	97.97	111.70
1	B	375	GLN	C-N-CA	5.19	134.68	121.70
1	A	227	GLU	N-CA-CB	5.02	119.64	110.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	355	ARG	Peptide

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1549	0	1552	68	0
1	B	1545	0	1544	66	0
1	C	1505	0	1508	43	0
2	A	6	0	0	0	0
2	B	3	0	0	0	0
2	C	4	0	0	0	0
All	All	4612	0	4604	151	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 16.

All (151) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:255:GLN:OE1	1:C:258:ARG:NH1	1.78	1.17
1:B:330:ARG:NH1	1:B:376:SER:C	2.17	0.98
1:A:342:ARG:NH2	1:A:364:PRO:O	2.00	0.94
1:B:330:ARG:NH2	1:B:377:GLU:HG2	1.84	0.93
1:B:239:LEU:HD21	1:C:240:GLN:HG2	1.52	0.92
1:A:348:MET:HG2	1:A:359:ILE:HD12	1.58	0.84
1:A:341:PHE:O	1:A:342:ARG:NH1	2.12	0.83
1:C:277:THR:HG22	1:C:394:GLN:HE22	1.43	0.81
1:B:330:ARG:HH12	1:B:376:SER:C	1.81	0.80
1:A:357:HIS:HB2	1:A:359:ILE:HD11	1.63	0.80
1:B:354:ASN:OD1	1:B:355:ARG:N	2.15	0.79
1:A:239:LEU:HD11	1:B:236:VAL:HG13	1.65	0.78
1:B:330:ARG:NH1	1:B:376:SER:O	2.20	0.75
1:C:355:ARG:HE	1:C:356:GLU:HA	1.51	0.74
1:B:232:LEU:HD21	1:C:233:GLU:HG2	1.71	0.73
1:C:355:ARG:HH21	1:C:357:HIS:H	1.37	0.73
1:B:274:THR:O	1:B:279:ARG:NH1	2.22	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:342:ARG:HH22	1:A:364:PRO:HB2	1.53	0.72
1:B:353:ASN:OD1	1:B:402:LYS:NZ	2.23	0.71
1:A:268:THR:HG22	1:A:411:ILE:HG23	1.72	0.70
1:C:279:ARG:HG3	1:C:311:LEU:HD13	1.74	0.69
1:A:366:LEU:N	1:A:366:LEU:HD23	2.06	0.69
1:B:330:ARG:HH12	1:B:376:SER:CA	2.07	0.68
1:A:364:PRO:HB3	1:A:371:PHE:CZ	2.29	0.68
1:B:375:GLN:HA	1:B:375:GLN:OE1	1.92	0.68
1:B:372:GLN:HG3	1:B:373:ARG:H	1.58	0.68
1:A:255:GLN:OE1	1:A:258:ARG:NH1	2.27	0.67
1:A:282:GLU:HG2	1:A:287:ARG:NH2	2.09	0.67
1:B:330:ARG:HB3	1:B:374:PRO:HB2	1.76	0.67
1:A:260:MET:HB3	1:C:260:MET:HE1	1.77	0.66
1:B:330:ARG:CZ	1:B:377:GLU:HG2	2.26	0.65
1:A:342:ARG:NE	1:A:342:ARG:HA	2.13	0.62
1:B:300:LYS:NZ	1:C:261:GLU:O	2.32	0.62
1:B:281:HIS:CE1	1:B:285:CYS:SG	2.92	0.62
1:B:235:ARG:NH2	1:C:240:GLN:OE1	2.30	0.62
1:B:228:ARG:NH2	1:C:233:GLU:OE2	2.22	0.62
1:B:305:LEU:HB3	1:B:326:ILE:HD11	1.80	0.62
1:B:235:ARG:O	1:B:239:LEU:HG	2.00	0.61
1:C:344:LYS:HB3	1:C:413:GLU:HG2	1.81	0.60
1:B:321:HIS:ND1	1:B:386:LEU:HB3	2.17	0.60
1:B:321:HIS:CE1	1:B:386:LEU:HB3	2.38	0.59
1:C:312:ASN:OD1	1:C:319:ARG:HA	2.03	0.59
1:A:300:LYS:NZ	1:B:263:ALA:O	2.36	0.58
1:B:351:ASP:OD1	1:B:354:ASN:HB3	2.04	0.57
1:A:242:THR:O	1:A:246:LYS:HG2	2.03	0.57
1:A:233:GLU:O	1:A:237:VAL:HG12	2.05	0.57
1:A:348:MET:CG	1:A:359:ILE:HD12	2.32	0.57
1:A:254:GLU:OE1	1:A:258:ARG:NH2	2.38	0.56
1:A:342:ARG:NE	1:A:366:LEU:CD2	2.69	0.55
1:B:330:ARG:NE	1:B:377:GLU:OE2	2.40	0.55
1:C:321:HIS:CA	1:C:390:LEU:HD23	2.37	0.55
1:B:246:LYS:NZ	1:C:247:ASP:OD2	2.40	0.54
1:A:260:MET:SD	1:C:260:MET:CE	2.95	0.54
1:B:258:ARG:HG3	1:B:259:LEU:N	2.21	0.54
1:B:242:THR:O	1:B:246:LYS:HG2	2.07	0.54
1:C:321:HIS:C	1:C:390:LEU:HD23	2.27	0.54
1:A:342:ARG:NH2	1:A:364:PRO:C	2.61	0.53
1:B:245:GLN:HG3	1:B:245:GLN:O	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:284:ALA:HA	1:B:312:ASN:HD22	1.73	0.53
1:A:246:LYS:NZ	1:B:247:ASP:OD1	2.37	0.52
1:B:325:PHE:HE1	1:B:385:PRO:HG3	1.75	0.52
1:A:343:ASN:HB3	1:A:413:GLU:O	2.09	0.52
1:B:251:GLY:O	1:B:254:GLU:HB2	2.10	0.52
1:A:341:PHE:O	1:A:342:ARG:CZ	2.57	0.52
1:A:369:ALA:HA	1:A:372:GLN:HG2	1.92	0.52
1:A:260:MET:SD	1:C:260:MET:HE2	2.50	0.51
1:C:229:ILE:O	1:C:233:GLU:HG3	2.10	0.51
1:A:282:GLU:HG2	1:A:287:ARG:CZ	2.41	0.51
1:A:357:HIS:HB2	1:A:359:ILE:CD1	2.39	0.51
1:C:321:HIS:CG	1:C:386:LEU:HA	2.46	0.51
1:B:238:GLU:O	1:B:242:THR:HG23	2.11	0.50
1:A:253:LEU:HD21	1:B:253:LEU:HD22	1.93	0.50
1:C:409:LYS:HE2	1:C:411:ILE:HD11	1.93	0.50
1:B:317:GLY:CA	1:B:386:LEU:HD21	2.41	0.50
1:C:364:PRO:HB3	1:C:371:PHE:CE2	2.47	0.50
1:A:342:ARG:NE	1:A:366:LEU:HD22	2.27	0.50
1:A:409:LYS:HE2	1:A:411:ILE:HD11	1.94	0.50
1:A:239:LEU:HD11	1:B:236:VAL:CG1	2.41	0.49
1:A:236:VAL:O	1:A:240:GLN:HG3	2.11	0.49
1:A:272:LYS:NZ	1:C:332:GLU:OE1	2.34	0.48
1:A:358:ALA:C	1:A:359:ILE:HD13	2.33	0.48
1:A:261:GLU:O	1:C:300:LYS:NZ	2.47	0.48
1:A:391:SER:OG	1:A:392:LYS:HE2	2.13	0.48
1:A:330:ARG:HB2	1:A:374:PRO:HB2	1.95	0.47
1:A:343:ASN:OD1	1:A:343:ASN:N	2.37	0.47
1:B:321:HIS:CD2	1:B:389:PRO:HB3	2.50	0.47
1:B:310:TYR:CD1	1:B:325:PHE:HE2	2.34	0.46
1:B:327:VAL:HG22	1:B:381:ALA:HB2	1.97	0.46
1:A:260:MET:CE	1:C:260:MET:HE3	2.46	0.46
1:A:342:ARG:NE	1:A:342:ARG:CA	2.75	0.46
1:B:372:GLN:CG	1:B:373:ARG:H	2.27	0.46
1:B:364:PRO:HB3	1:B:371:PHE:CZ	2.50	0.46
1:A:257:LEU:HD11	1:C:260:MET:SD	2.56	0.46
1:A:342:ARG:HH12	1:A:371:PHE:CB	2.28	0.46
1:B:339:TRP:CE2	1:B:373:ARG:HB2	2.51	0.46
1:B:336:LEU:HD11	1:C:407:PHE:CD2	2.51	0.46
1:A:311:LEU:HA	1:A:322:LEU:HD23	1.96	0.46
1:A:342:ARG:HH22	1:A:364:PRO:CB	2.27	0.45
1:B:313:GLY:HA3	1:B:317:GLY:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:342:ARG:HH22	1:A:364:PRO:C	2.20	0.45
1:C:226:ARG:O	1:C:229:ILE:HB	2.16	0.45
1:B:321:HIS:HA	1:B:389:PRO:HA	1.99	0.45
1:A:260:MET:CE	1:C:260:MET:CE	2.95	0.44
1:A:342:ARG:CZ	1:A:342:ARG:HA	2.46	0.44
1:C:355:ARG:HE	1:C:356:GLU:CA	2.24	0.44
1:A:285:CYS:SG	1:A:286:GLY:N	2.91	0.44
1:C:411:ILE:N	1:C:411:ILE:HD12	2.33	0.44
1:B:364:PRO:HB3	1:B:371:PHE:CE1	2.52	0.44
1:A:270:LEU:HB3	1:C:301:TYR:CE1	2.52	0.44
1:A:232:LEU:HD21	1:B:233:GLU:HG2	2.00	0.44
1:A:251:GLY:O	1:A:254:GLU:HB3	2.17	0.44
1:A:264:SER:HB3	1:A:296:PHE:HB2	2.00	0.44
1:B:373:ARG:HA	1:B:374:PRO:HD3	1.85	0.44
1:A:363:ARG:HG2	1:A:363:ARG:NH1	2.33	0.43
1:A:279:ARG:HA	1:A:282:GLU:HG3	2.00	0.43
1:C:230:LEU:CD1	1:C:234:GLN:HE21	2.31	0.43
1:B:301:TYR:CZ	1:C:270:LEU:HB2	2.54	0.43
1:A:278:ARG:HG2	1:A:282:GLU:OE1	2.19	0.43
1:A:235:ARG:O	1:A:239:LEU:HG	2.19	0.43
1:A:341:PHE:C	1:A:342:ARG:HD2	2.39	0.43
1:B:330:ARG:CB	1:B:374:PRO:HB2	2.47	0.43
1:C:237:VAL:O	1:C:241:GLN:HG2	2.18	0.43
1:B:366:LEU:HA	1:B:366:LEU:HD23	1.55	0.43
1:B:224:LEU:O	1:B:228:ARG:HG3	2.18	0.43
1:A:311:LEU:HA	1:A:322:LEU:CD2	2.49	0.42
1:B:235:ARG:HH21	1:C:240:GLN:CD	2.17	0.42
1:C:321:HIS:CD2	1:C:386:LEU:HA	2.54	0.42
1:B:235:ARG:HE	1:C:240:GLN:HE22	1.68	0.42
1:A:310:TYR:CD2	1:A:314:ASP:HB2	2.55	0.42
1:B:397:LYS:HG2	1:B:397:LYS:H	1.65	0.42
1:C:236:VAL:O	1:C:240:GLN:HG3	2.20	0.42
1:A:342:ARG:NH2	1:A:364:PRO:HB2	2.29	0.41
1:A:255:GLN:O	1:A:259:LEU:HG	2.21	0.41
1:A:275:ASN:O	1:A:279:ARG:HG3	2.20	0.41
1:C:320:THR:OG1	1:C:321:HIS:ND1	2.34	0.41
1:B:365:ASP:N	1:B:365:ASP:OD1	2.53	0.41
1:B:388:PHE:HA	1:B:389:PRO:HD3	1.98	0.41
1:B:409:LYS:HE2	1:B:411:ILE:HD11	2.02	0.41
1:B:275:ASN:HB3	1:B:278:ARG:HD2	2.02	0.41
1:A:395:SER:HA	1:A:396:PRO:HD2	1.91	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:363:ARG:HG2	1:A:363:ARG:HH11	1.86	0.40
1:B:320:THR:CB	1:B:321:HIS:CD2	3.04	0.40
1:C:364:PRO:HB3	1:C:371:PHE:CD2	2.56	0.40
1:B:354:ASN:OD1	1:B:356:GLU:N	2.53	0.40
1:A:260:MET:HB3	1:C:260:MET:CE	2.49	0.40
1:A:275:ASN:HA	1:A:404:ASP:OD2	2.22	0.40
1:A:342:ARG:NE	1:A:366:LEU:HD21	2.37	0.40
1:B:320:THR:CB	1:B:321:HIS:HD2	2.35	0.40
1:C:364:PRO:HB3	1:C:371:PHE:CZ	2.56	0.40
1:B:260:MET:HE2	1:B:260:MET:HB3	1.77	0.40
1:B:372:GLN:HG3	1:B:373:ARG:N	2.32	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	190/197 (96%)	178 (94%)	12 (6%)	0	100	100
1	B	187/197 (95%)	179 (96%)	8 (4%)	0	100	100
1	C	182/197 (92%)	176 (97%)	6 (3%)	0	100	100
All	All	559/591 (95%)	533 (95%)	26 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	171/176 (97%)	169 (99%)	2 (1%)	71	92
1	B	171/176 (97%)	165 (96%)	6 (4%)	36	70
1	C	166/176 (94%)	161 (97%)	5 (3%)	41	75
All	All	508/528 (96%)	495 (97%)	13 (3%)	46	79

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	245	GLN
1	A	415	SER
1	B	252	LYS
1	B	256	SER
1	B	279	ARG
1	B	330	ARG
1	B	337	LEU
1	B	367	SER
1	C	261	GLU
1	C	285	CYS
1	C	343	ASN
1	C	355	ARG
1	C	356	GLU

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

Mol	Chain	Res	Type
1	A	240	GLN
1	A	241	GLN
1	B	248	GLN
1	B	281	HIS
1	B	321	HIS
1	C	234	GLN
1	C	394	GLN

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	192/197 (97%)	-0.25	4 (2%) 63 54	43, 64, 97, 132	0
1	B	191/197 (96%)	-0.05	7 (3%) 41 31	46, 81, 123, 163	0
1	C	186/197 (94%)	-0.22	5 (2%) 54 44	49, 73, 124, 144	0
All	All	569/591 (96%)	-0.18	16 (2%) 53 43	43, 72, 121, 163	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	367	SER	4.9
1	A	287	ARG	3.7
1	A	288	THR	3.6
1	C	396	PRO	3.2
1	C	354	ASN	3.1
1	B	221	GLN	3.1
1	B	366	LEU	2.8
1	B	226	ARG	2.7
1	C	279	ARG	2.7
1	B	369	ALA	2.6
1	A	342	ARG	2.4
1	C	398	HIS	2.2
1	C	226	ARG	2.1
1	B	248	GLN	2.1
1	B	375	GLN	2.0
1	A	355	ARG	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 [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.