



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:31 AM GMT

PDB ID : 2HIG
Title : Crystal Structure of Phosphofructokinase apoenzyme from Trypanosoma brucei.
Authors : Martinez-Oyanedel, J.; McNae, I.W.; Fothergill-Gilmore, L.A.; Walkinshaw, M.D.
Deposited on : 2006-06-29
Resolution : 2.40 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

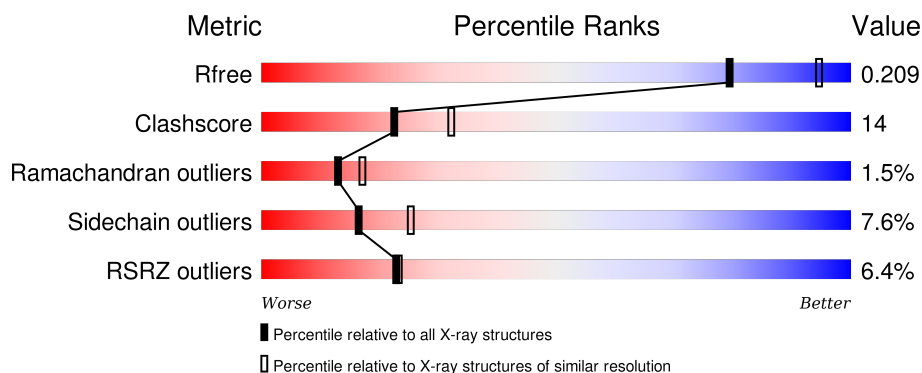
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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	487	
1	B	487	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	NA	A	488	-	-	-	X
2	NA	B	488	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6807 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 6-phospho-1-fructokinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	440	Total	C	N	O	S	0	0	0
			3374	2121	614	622	17			
1	B	422	Total	C	N	O	S	0	0	0
			3229	2034	583	595	17			

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Na	0	0
			1	1		
2	A	1	Total	Na	0	0
			1	1		

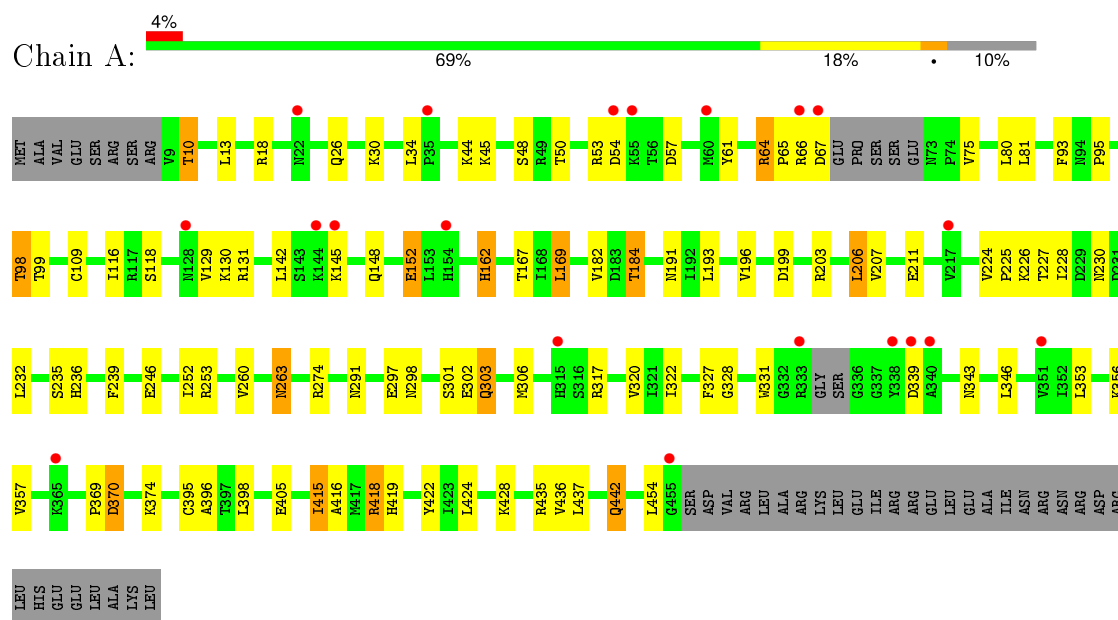
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	114	Total	O	0	0
			114	114		
3	B	88	Total	O	0	0
			88	88		

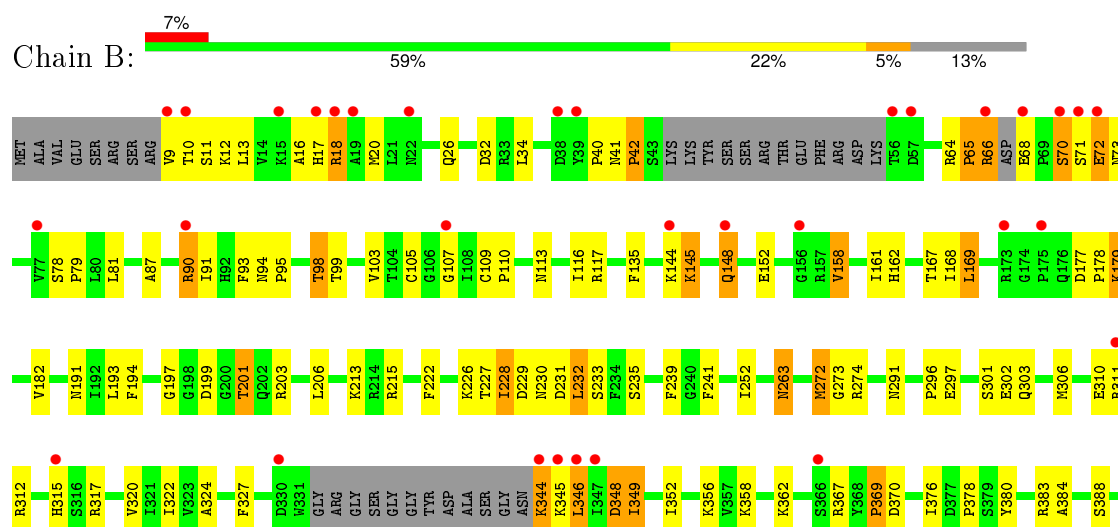
3 Residue-property plots [i](#)

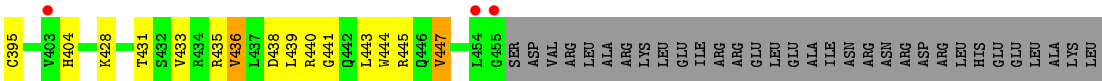
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: 6-phospho-1-fructokinase



• Molecule 1: 6-phospho-1-fructokinase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	118.77Å 92.50Å 79.53Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	21.45 – 2.40 21.45 – 2.40	Depositor EDS
% Data completeness (in resolution range)	98.2 (21.45-2.40) 98.3 (21.45-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.07 (at 2.41Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.201 , 0.238 0.203 , 0.209	Depositor DCC
R_{free} test set	1717 reflections (5.02%)	DCC
Wilson B-factor (Å ²)	43.8	Xtriage
Anisotropy	0.200	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 57.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 34225 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6807	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.27% 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, 0.375 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: NA

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	0.36	0/3433	0.62	0/4645
1	B	0.36	0/3285	0.62	0/4450
All	All	0.36	0/6718	0.62	0/9095

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	3374	0	3426	75	0
1	B	3229	0	3285	110	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	114	0	0	3	0
3	B	88	0	0	4	0
All	All	6807	0	6711	183	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 14.

All (183) 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:B:291:ASN:HD22	1:B:320:VAL:H	1.08	1.02
1:A:291:ASN:HD22	1:A:320:VAL:H	1.17	0.93
1:A:99:THR:H	1:A:191:ASN:HD22	1.19	0.91
1:B:263:ASN:HD21	1:B:317:ARG:HH11	1.14	0.90
1:B:99:THR:H	1:B:191:ASN:HD22	1.17	0.87
1:A:26:GLN:HE22	1:A:81:LEU:H	1.25	0.84
1:B:103:VAL:HG12	1:B:135:PHE:HB2	1.60	0.83
1:B:26:GLN:HE22	1:B:81:LEU:H	1.27	0.80
1:A:297:GLU:HG2	1:A:436:VAL:HG13	1.65	0.79
1:B:199:ASP:HB2	1:B:431:THR:HG22	1.66	0.76
1:B:103:VAL:CG1	1:B:135:PHE:HB2	2.16	0.75
1:B:291:ASN:ND2	1:B:320:VAL:H	1.85	0.74
1:B:263:ASN:HD21	1:B:317:ARG:NH1	1.85	0.74
1:B:179:LYS:HD3	1:B:215:ARG:HH21	1.54	0.72
1:B:233:SER:CB	1:B:433:VAL:HG23	2.20	0.71
1:B:144:LYS:O	1:B:145:LYS:HB2	1.91	0.69
1:A:26:GLN:HE22	1:A:81:LEU:N	1.90	0.69
1:A:302:GLU:OE2	1:A:356:LYS:HE2	1.93	0.68
1:A:339:ASP:OD2	1:A:343:ASN:HB2	1.95	0.66
1:B:144:LYS:O	1:B:145:LYS:CB	2.44	0.66
1:A:297:GLU:CG	1:A:436:VAL:HG13	2.25	0.66
1:A:303:GLN:NE2	1:A:303:GLN:H	1.93	0.65
1:B:346:LEU:HD13	1:B:346:LEU:H	1.63	0.64
1:B:274:ARG:NH1	1:B:346:LEU:HD11	2.13	0.64
1:A:13:LEU:HD13	1:A:61:TYR:OH	1.98	0.63
1:A:236:HIS:CE1	1:A:419:HIS:H	2.17	0.62
1:B:291:ASN:HD22	1:B:320:VAL:N	1.89	0.62
1:B:345:LYS:NZ	1:B:345:LYS:HB2	2.15	0.62
1:B:362:LYS:NZ	1:B:362:LYS:HB3	2.15	0.61
1:B:263:ASN:ND2	1:B:317:ARG:HH11	1.92	0.61
1:A:405:GLU:HG3	1:A:424:LEU:HD21	1.82	0.60
1:A:10:THR:HB	1:A:442:GLN:OE1	2.01	0.60
1:B:116:ILE:HD13	1:B:169:LEU:HD22	1.82	0.60
1:A:246:GLU:OE2	1:B:388:SER:HB2	2.02	0.59
1:A:418:ARG:HG2	1:A:419:HIS:ND1	2.18	0.58
1:B:66:ARG:HB3	3:B:540:HOH:O	2.02	0.58
1:A:301:SER:HB2	1:A:303:GLN:HE22	1.68	0.58
1:B:162:HIS:CD2	1:B:162:HIS:H	2.20	0.58
1:B:302:GLU:HG2	1:B:303:GLN:NE2	2.19	0.58
1:B:274:ARG:HH11	1:B:346:LEU:HD11	1.67	0.58
1:B:228:ILE:HD12	1:B:229:ASP:H	1.68	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:252:ILE:HD13	1:A:322:ILE:HD11	1.87	0.57
1:B:117:ARG:HA	1:B:158:VAL:HG21	1.85	0.57
1:B:94:ASN:O	1:B:98:THR:HG22	2.03	0.57
1:B:109:CYS:HB2	1:B:228:ILE:HG23	1.86	0.57
1:B:12:LYS:HE2	3:B:489:HOH:O	2.05	0.57
1:B:344:LYS:HD3	1:B:344:LYS:N	2.19	0.57
1:B:95:PRO:O	1:B:98:THR:HG23	2.06	0.56
1:B:91:ILE:HD11	1:B:95:PRO:HG3	1.88	0.55
1:B:68:GLU:C	1:B:70:SER:H	2.10	0.55
1:B:20:MET:SD	1:B:428:LYS:HG2	2.47	0.54
1:A:199:ASP:H	1:A:230:ASN:ND2	2.05	0.54
1:A:93:PHE:HB3	1:A:98:THR:CG2	2.38	0.54
1:A:109:CYS:HB2	1:A:228:ILE:HD11	1.90	0.54
1:B:226:LYS:HA	3:B:492:HOH:O	2.07	0.54
1:B:232:LEU:HD13	1:B:235:SER:HB2	1.89	0.54
1:B:199:ASP:CB	1:B:431:THR:HG22	2.37	0.54
1:B:239:PHE:HB2	1:B:395:CYS:SG	2.48	0.54
1:B:233:SER:OG	1:B:433:VAL:HG23	2.08	0.53
1:A:199:ASP:OD2	1:A:203:ARG:NH1	2.42	0.53
1:A:291:ASN:ND2	1:A:320:VAL:H	1.97	0.53
1:A:291:ASN:HD22	1:A:320:VAL:N	1.98	0.53
1:A:301:SER:HB2	1:A:303:GLN:NE2	2.24	0.53
1:B:116:ILE:CD1	1:B:169:LEU:HD22	2.39	0.53
1:B:310:GLU:CD	1:B:367:ARG:HH12	2.12	0.53
1:B:348:ASP:O	1:B:352:ILE:HG12	2.09	0.53
1:B:206:LEU:HD11	1:B:428:LYS:HG3	1.91	0.52
1:B:230:ASN:HB2	1:B:241:PHE:CD2	2.44	0.52
1:A:66:ARG:HH21	1:A:75:VAL:HG21	1.74	0.52
1:A:184:THR:HG22	3:A:545:HOH:O	2.09	0.52
1:B:199:ASP:HB2	1:B:431:THR:CG2	2.39	0.52
1:B:358:LYS:O	1:B:362:LYS:HG2	2.09	0.52
1:B:362:LYS:HZ2	1:B:362:LYS:HB3	1.74	0.52
1:B:228:ILE:HD12	1:B:229:ASP:N	2.24	0.52
1:A:239:PHE:HB2	1:A:395:CYS:SG	2.49	0.52
1:B:199:ASP:O	1:B:203:ARG:HG3	2.10	0.52
1:B:302:GLU:O	1:B:306:MET:HG2	2.10	0.52
1:A:109:CYS:SG	1:A:226:LYS:HD2	2.50	0.51
1:B:297:GLU:CG	1:B:436:VAL:HG22	2.40	0.51
1:B:9:VAL:HG23	1:B:11:SER:H	1.73	0.51
1:B:105:CYS:HB2	1:B:201:THR:HG21	1.94	0.50
1:A:162:HIS:CD2	1:A:162:HIS:H	2.27	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:196:VAL:HA	1:A:224:VAL:O	2.12	0.50
1:B:296:PRO:HD3	1:B:324:ALA:HB3	1.92	0.50
1:A:57:ASP:HB3	1:B:64:ARG:NH2	2.27	0.50
1:B:235:SER:O	1:B:435:ARG:HD3	2.11	0.50
1:A:93:PHE:HB3	1:A:98:THR:HG21	1.94	0.50
1:B:93:PHE:O	1:B:95:PRO:HD3	2.12	0.50
1:B:310:GLU:HG3	1:B:367:ARG:NH1	2.26	0.50
1:B:252:ILE:HD13	1:B:322:ILE:HD11	1.94	0.50
1:A:235:SER:OG	1:A:416:ALA:HB1	2.12	0.49
1:A:263:ASN:HD21	1:A:317:ARG:HH11	1.59	0.49
1:B:110:PRO:HD3	1:B:383:ARG:HA	1.93	0.49
1:B:327:PHE:O	1:B:349:ILE:HG13	2.13	0.48
1:B:274:ARG:HH12	1:B:346:LEU:HD21	1.77	0.48
1:B:177:ASP:OD1	1:B:179:LYS:CB	2.61	0.48
1:A:303:GLN:HE21	1:A:303:GLN:H	1.59	0.48
1:A:252:ILE:CD1	1:A:322:ILE:HD11	2.43	0.48
1:B:117:ARG:HA	1:B:158:VAL:CG2	2.43	0.48
1:B:178:PRO:O	1:B:182:VAL:HG23	2.13	0.48
1:B:227:THR:CG2	1:B:229:ASP:HB2	2.44	0.48
1:B:199:ASP:CG	1:B:203:ARG:HH12	2.17	0.48
1:B:302:GLU:OE2	1:B:356:LYS:HD3	2.14	0.47
1:B:232:LEU:CD1	1:B:235:SER:HB2	2.44	0.47
1:B:444:TRP:HA	1:B:447:VAL:HG13	1.95	0.47
1:A:203:ARG:O	1:A:207:VAL:HG23	2.14	0.47
1:A:142:LEU:O	1:A:184:THR:HG21	2.13	0.47
1:A:116:ILE:HD13	1:A:169:LEU:HD22	1.95	0.47
1:B:441:GLY:O	1:B:445:ARG:HD3	2.15	0.47
1:A:66:ARG:HE	1:A:75:VAL:CG2	2.28	0.46
1:A:369:PRO:HG2	1:A:370:ASP:H	1.79	0.46
1:A:398:LEU:HD23	1:A:422:TYR:CE1	2.51	0.46
1:B:158:VAL:HG23	1:B:161:ILE:HD13	1.98	0.46
1:B:32:ASP:OD2	1:B:213:LYS:NZ	2.45	0.46
1:A:230:ASN:ND2	1:A:232:LEU:H	2.14	0.46
1:A:53:ARG:NH1	1:A:57:ASP:HB2	2.31	0.46
1:B:105:CYS:HB2	1:B:201:THR:CG2	2.46	0.46
1:A:99:THR:N	1:A:191:ASN:HD22	2.00	0.46
1:B:345:LYS:HB3	1:B:346:LEU:HD13	1.98	0.46
1:B:98:THR:HA	1:B:191:ASN:ND2	2.31	0.46
1:B:99:THR:H	1:B:191:ASN:ND2	1.98	0.46
1:B:194:PHE:HA	1:B:222:PHE:O	2.15	0.46
1:B:91:ILE:HD11	1:B:95:PRO:CG	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:199:ASP:HB2	1:B:203:ARG:HH12	1.80	0.45
1:B:168:ILE:HG13	1:B:169:LEU:N	2.29	0.45
1:B:26:GLN:HE22	1:B:81:LEU:N	2.06	0.45
1:B:145:LYS:O	1:B:148:GLN:HB3	2.16	0.45
1:B:301:SER:HB2	1:B:303:GLN:OE1	2.17	0.45
1:A:226:LYS:HB3	1:A:239:PHE:CZ	2.51	0.45
1:A:145:LYS:O	1:A:148:GLN:HB3	2.17	0.45
1:A:99:THR:HG23	1:A:131:ARG:HG2	1.98	0.45
1:B:274:ARG:HG2	1:B:274:ARG:NH1	2.32	0.45
1:B:274:ARG:HG2	1:B:274:ARG:HH11	1.81	0.45
1:A:327:PHE:CD1	1:A:328:GLY:N	2.85	0.44
1:B:369:PRO:HG2	1:B:370:ASP:H	1.82	0.44
1:B:227:THR:HG22	1:B:229:ASP:N	2.33	0.44
1:B:312:ARG:HA	1:B:315:HIS:CE1	2.52	0.44
1:A:274:ARG:HD2	1:A:346:LEU:HD11	1.99	0.44
1:B:438:ASP:C	1:B:440:ARG:H	2.20	0.44
1:B:113:ASN:OD1	1:B:168:ILE:HG23	2.18	0.44
1:A:353:LEU:O	1:A:357:VAL:HG23	2.18	0.44
1:A:302:GLU:O	1:A:306:MET:HG2	2.17	0.44
1:A:418:ARG:HD3	3:A:559:HOH:O	2.16	0.44
1:B:13:LEU:HD21	3:B:545:HOH:O	2.18	0.44
1:A:182:VAL:HG21	1:A:211:GLU:CG	2.47	0.44
1:A:44:LYS:HB3	1:A:44:LYS:NZ	2.33	0.43
1:A:26:GLN:NE2	1:A:80:LEU:HA	2.34	0.43
1:B:345:LYS:HB3	1:B:346:LEU:H	1.47	0.43
1:B:380:TYR:O	1:B:384:ALA:HB3	2.17	0.43
1:A:454:LEU:HD23	1:A:454:LEU:HA	1.88	0.43
1:B:197:GLY:HA3	1:B:201:THR:HG21	2.00	0.43
1:A:328:GLY:O	1:A:331:TRP:HB2	2.19	0.43
1:A:230:ASN:HD21	1:A:232:LEU:HB3	1.84	0.43
1:A:298:ASN:HB3	3:A:494:HOH:O	2.18	0.43
1:A:206:LEU:HD11	1:A:428:LYS:HG3	2.00	0.43
1:A:65:PRO:HA	1:A:75:VAL:O	2.19	0.42
1:A:227:THR:HG22	1:A:232:LEU:HD23	2.02	0.42
1:A:327:PHE:CG	1:A:328:GLY:N	2.86	0.42
1:A:130:LYS:O	1:A:131:ARG:HB3	2.18	0.42
1:B:438:ASP:O	1:B:440:ARG:N	2.52	0.42
1:A:45:LYS:HG2	1:A:48:SER:OG	2.20	0.42
1:B:376:ILE:HG22	1:B:378:PRO:HD3	2.01	0.42
1:A:236:HIS:CE1	1:A:418:ARG:HG3	2.55	0.41
1:B:9:VAL:HG23	1:B:10:THR:N	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:87:ALA:HA	1:B:404:HIS:HB3	2.01	0.41
1:A:274:ARG:CD	1:A:346:LEU:HD11	2.50	0.41
1:B:18:ARG:O	1:B:203:ARG:NH2	2.42	0.41
1:B:78:SER:HA	1:B:79:PRO:HD3	1.86	0.41
1:A:131:ARG:NH2	1:A:152:GLU:OE2	2.41	0.41
1:B:177:ASP:OD1	1:B:179:LYS:HB2	2.21	0.41
1:A:162:HIS:N	1:A:162:HIS:CD2	2.89	0.41
1:B:40:PRO:HB3	1:B:90:ARG:HH21	1.85	0.41
1:B:41:ASN:HA	1:B:42:PRO:HD3	1.89	0.41
1:A:98:THR:HA	1:A:191:ASN:ND2	2.36	0.41
1:B:16:ALA:O	1:B:17:HIS:C	2.59	0.41
1:A:398:LEU:HD22	1:A:415:ILE:HG12	2.02	0.40
1:B:272:MET:HG2	1:B:272:MET:H	1.64	0.40
1:A:118:SER:OG	1:A:396:ALA:HA	2.21	0.40
1:A:436:VAL:CG1	1:A:437:LEU:N	2.85	0.40
1:B:203:ARG:HG3	1:B:203:ARG:HH11	1.86	0.40
1:B:346:LEU:HD13	1:B:346:LEU:N	2.35	0.40
1:A:64:ARG:HH11	1:A:64:ARG:CG	2.33	0.40
1:B:107:GLY:O	1:B:226:LYS:NZ	2.36	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	434/487 (89%)	412 (95%)	20 (5%)	2 (0%)	34	48
1	B	414/487 (85%)	387 (94%)	16 (4%)	11 (3%)	6	6
All	All	848/974 (87%)	799 (94%)	36 (4%)	13 (2%)	13	17

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	18	ARG
1	B	70	SER
1	B	71	SER
1	B	73	ASN
1	B	145	LYS
1	B	273	GLY
1	B	439	LEU
1	B	348	ASP
1	A	95	PRO
1	B	65	PRO
1	B	72	GLU
1	B	42	PRO
1	B	369	PRO

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	363/405 (90%)	336 (93%)	27 (7%)	17	26
1	B	350/405 (86%)	323 (92%)	27 (8%)	16	24
All	All	713/810 (88%)	659 (92%)	54 (8%)	16	25

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	THR
1	A	30	LYS
1	A	34	LEU
1	A	50	THR
1	A	54	ASP
1	A	64	ARG
1	A	67	ASP
1	A	98	THR
1	A	129	VAL
1	A	152	GLU
1	A	162	HIS

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Mol	Chain	Res	Type
1	A	167	THR
1	A	169	LEU
1	A	184	THR
1	A	193	LEU
1	A	206	LEU
1	A	225	PRO
1	A	253	ARG
1	A	260	VAL
1	A	263	ASN
1	A	303	GLN
1	A	370	ASP
1	A	374	LYS
1	A	415	ILE
1	A	418	ARG
1	A	435	ARG
1	A	442	GLN
1	B	18	ARG
1	B	34	LEU
1	B	65	PRO
1	B	66	ARG
1	B	72	GLU
1	B	90	ARG
1	B	98	THR
1	B	148	GLN
1	B	152	GLU
1	B	158	VAL
1	B	167	THR
1	B	169	LEU
1	B	179	LYS
1	B	193	LEU
1	B	201	THR
1	B	228	ILE
1	B	231	ASP
1	B	232	LEU
1	B	263	ASN
1	B	272	MET
1	B	311	ARG
1	B	344	LYS
1	B	346	LEU
1	B	349	ILE
1	B	436	VAL
1	B	443	LEU

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Mol	Chain	Res	Type
1	B	447	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	26	GLN
1	A	73	ASN
1	A	162	HIS
1	A	191	ASN
1	A	230	ASN
1	A	242	GLN
1	A	263	ASN
1	A	282	GLN
1	A	289	GLN
1	A	291	ASN
1	A	298	ASN
1	A	303	GLN
1	A	421	ASN
1	A	442	GLN
1	A	446	GLN
1	B	26	GLN
1	B	125	ASN
1	B	148	GLN
1	B	162	HIS
1	B	163	HIS
1	B	191	ASN
1	B	263	ASN
1	B	282	GLN
1	B	289	GLN
1	B	291	ASN
1	B	442	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 [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	440/487 (90%)	0.03	20 (4%) 37 38	25, 41, 66, 88	0
1	B	422/487 (86%)	0.20	35 (8%) 14 14	28, 44, 71, 105	0
All	All	862/974 (88%)	0.11	55 (6%) 23 23	25, 43, 68, 105	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	344	LYS	5.8
1	B	346	LEU	5.6
1	B	71	SER	5.4
1	B	9	VAL	5.2
1	B	56	THR	4.5
1	B	345	LYS	4.3
1	B	17	HIS	4.2
1	B	18	ARG	3.6
1	B	455	GLY	3.6
1	B	330	ASP	3.4
1	B	22	ASN	3.4
1	B	175	PRO	3.4
1	B	72	GLU	3.4
1	A	338	TYR	3.2
1	B	10	THR	3.2
1	A	54	ASP	3.1
1	A	315	HIS	3.1
1	B	366	SER	3.1
1	B	70	SER	3.1
1	A	333	ARG	3.0
1	B	454	LEU	3.0
1	A	340	ALA	2.9
1	A	145	LYS	2.7
1	B	403	VAL	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	144	LYS	2.7
1	B	107	GLY	2.6
1	B	38	ASP	2.6
1	B	311	ARG	2.6
1	A	217	VAL	2.5
1	B	77	VAL	2.5
1	A	22	ASN	2.5
1	A	55	LYS	2.5
1	B	90	ARG	2.5
1	A	66	ARG	2.5
1	B	39	TYR	2.5
1	A	60	MET	2.4
1	B	315	HIS	2.4
1	A	144	LYS	2.4
1	B	148	GLN	2.4
1	B	66	ARG	2.3
1	A	339	ASP	2.3
1	B	57	ASP	2.3
1	B	19	ALA	2.3
1	B	68	GLU	2.3
1	A	351	VAL	2.3
1	A	455	GLY	2.2
1	A	154	HIS	2.2
1	B	173	ARG	2.2
1	A	67	ASP	2.2
1	B	15	LYS	2.2
1	A	35	PRO	2.1
1	B	156	GLY	2.1
1	A	128	ASN	2.1
1	B	347	ILE	2.1
1	A	365	LYS	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 ⓘ

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
2	NA	B	488	1/1	0.95	0.21	6.55	44,44,44,44	0
2	NA	A	488	1/1	0.91	0.27	4.34	46,46,46,46	0

6.5 Other polymers [i](#)

There are no such residues in this entry.