Full wwPDB X-ray Structure Validation Report ⓘ

Aug 8, 2023 – 06:30 AM EDT

PDB ID	:	1OY3
Title	:	CRYSTAL STRUCTURE OF AN IKBBETA/NF-KB P65 HOMODIMER COMPLEX
Authors	:	Malek, S.; Huang, D.B.; Huxford, T.; Ghosh, S.; Ghosh, G.
Deposited on	:	2003-04-03
Resolution	:	2.05 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 types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references ⓘ) were used in the production of this report:

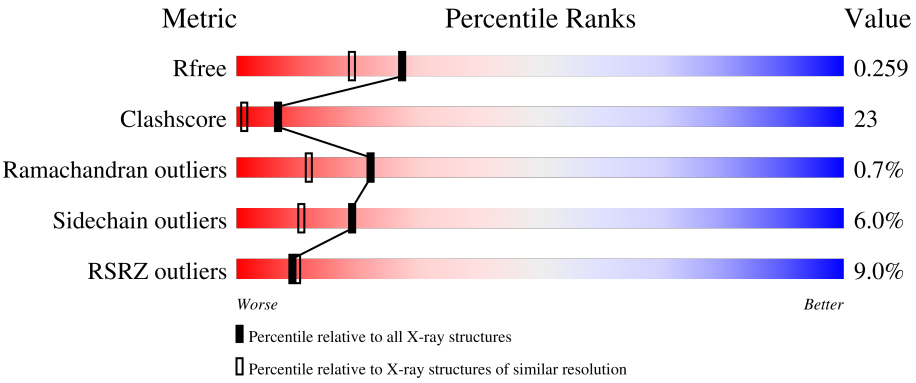
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

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 2.05 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	1692 (2.04-2.04)
Clashscore	141614	1773 (2.04-2.04)
Ramachandran outliers	138981	1752 (2.04-2.04)
Sidechain outliers	138945	1752 (2.04-2.04)
RSRZ outliers	127900	1672 (2.04-2.04)

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 of 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 <=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	B	136	9% 46% 33% 18%
1	C	136	11% 61% 35%
2	D	282	5% 49% 27% 22%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 3888 atoms, of which 4 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 Transcription factor p65.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	135	Total	C	N	O	S	0	0	0
			1093	683	199	205	6			
1	B	112	Total	C	N	O	S	0	0	0
			891	553	160	173	5			

- Molecule 2 is a protein called transcription factor inhibitor I-kappa-B-beta.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	D	220	Total	C	H	N	O	S	0	0	0
			1656	1040	4	303	303	6			

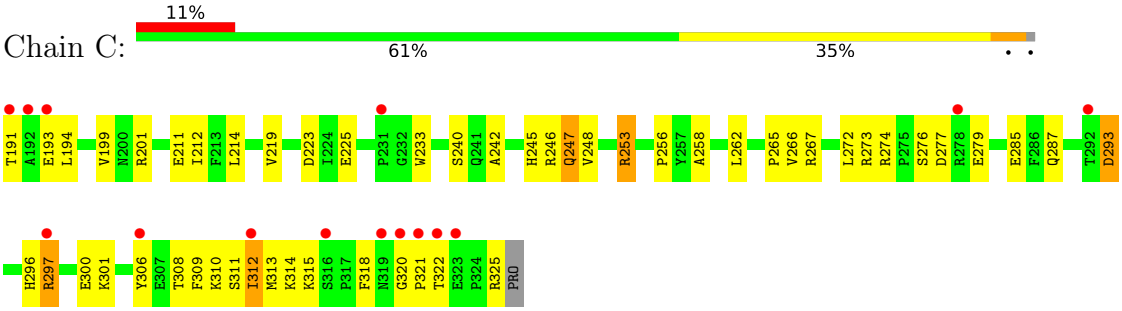
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	63	Total	O	0	0
			63	63		
3	B	62	Total	O	0	0
			62	62		
3	D	123	Total	O	0	0
			123	123		

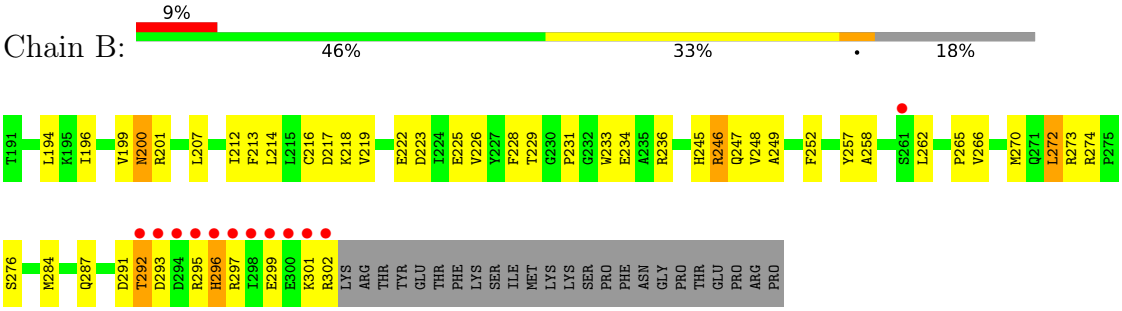
3 Residue-property plots (i)

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

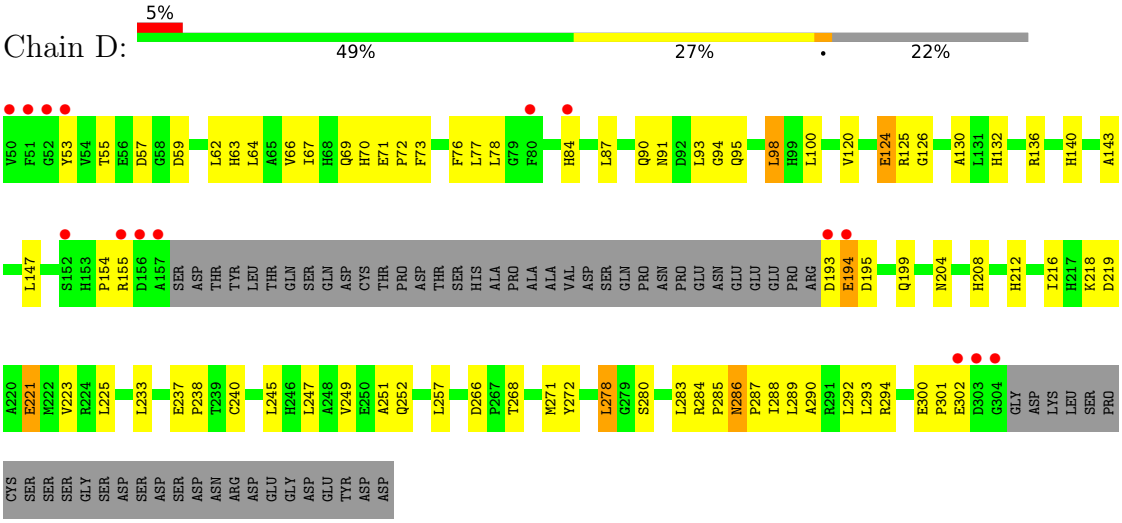
• Molecule 1: Transcription factor p65



• Molecule 1: Transcription factor p65



• Molecule 2: transcription factor inhibitor I-kappa-B-beta



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	48.62Å 54.79Å 59.57Å 102.56° 96.07° 91.90°	Depositor
Resolution (Å)	30.00 – 2.05 35.47 – 2.04	Depositor EDS
% Data completeness (in resolution range)	78.1 (30.00-2.05) 84.5 (35.47-2.04)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.47 (at 2.05Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.218 , 0.247 0.234 , 0.259	Depositor DCC
R_{free} test set	1589 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å^2)	20.2	Xtriage
Anisotropy	0.531	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 54.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	3888	wwPDB-VP
Average B, all atoms (Å^2)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.61% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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	B	0.38	0/909	0.68	1/1229 (0.1%)
1	C	0.35	0/1118	0.62	0/1508
2	D	0.34	0/1688	0.63	3/2300 (0.1%)
All	All	0.35	0/3715	0.64	4/5037 (0.1%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	D	84	HIS	CB-CA-C	10.01	130.43	110.40
2	D	238	PRO	N-CA-C	5.79	127.15	112.10
1	B	296	HIS	CB-CA-C	5.33	121.06	110.40
2	D	71	GLU	CB-CA-C	5.06	120.51	110.40

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	B	891	0	854	50	0
1	C	1093	0	1073	49	0
2	D	1652	4	1631	77	0
3	B	62	0	0	2	0
3	C	63	0	0	2	0
3	D	123	0	0	10	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	3884	4	3558	167	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 23.

All (167) 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:246:ARG:HG3	1:B:248:VAL:HG21	1.27	1.15
2:D:70:HIS:HD2	2:D:73:PHE:H	1.17	0.93
2:D:286:ASN:HD21	2:D:288:ILE:HG12	1.32	0.93
2:D:55:THR:HG23	2:D:57:ASP:H	1.33	0.92
1:B:291:ASP:OD1	3:B:653:HOH:O	1.97	0.82
1:C:223:ASP:OD1	1:C:276:SER:HB2	1.81	0.81
1:B:296:HIS:O	1:B:301:LYS:HB2	1.80	0.81
1:C:262:LEU:HD11	1:C:266:VAL:HG21	1.62	0.80
2:D:55:THR:HG22	2:D:59:ASP:H	1.48	0.79
1:C:248:VAL:HG22	1:B:245:HIS:NE2	1.97	0.78
1:B:296:HIS:HA	1:B:299:GLU:HB2	1.65	0.78
1:B:297:ARG:HE	1:B:302:ARG:HD2	1.49	0.77
2:D:212:HIS:O	2:D:216:ILE:HG12	1.86	0.76
1:B:200:ASN:H	1:B:200:ASN:HD22	1.32	0.76
2:D:93:LEU:HB3	2:D:125:ARG:NH1	2.02	0.75
2:D:208:HIS:HA	3:D:826:HOH:O	1.88	0.73
1:B:200:ASN:ND2	1:B:213:PHE:H	1.87	0.73
1:B:199:VAL:HG12	1:B:201:ARG:H	1.53	0.73
1:C:274:ARG:HG2	1:C:276:SER:HB3	1.71	0.73
2:D:266:ASP:OD2	2:D:268:THR:HG22	1.88	0.72
1:C:308:THR:O	1:C:312:ILE:HG12	1.89	0.72
2:D:125:ARG:HG2	2:D:125:ARG:HH11	1.55	0.72
2:D:252:GLN:HA	2:D:289:LEU:HD11	1.69	0.72
1:C:265:PRO:HB3	1:C:287:GLN:OE1	1.91	0.70
2:D:62:LEU:HD11	2:D:87:LEU:HD23	1.73	0.70
1:B:262:LEU:HD11	1:B:266:VAL:CG2	2.20	0.70
1:C:199:VAL:HG12	1:C:201:ARG:H	1.56	0.69
1:B:293:ASP:HA	1:B:297:ARG:HD2	1.73	0.69
2:D:70:HIS:CD2	2:D:73:PHE:H	2.06	0.69
1:B:223:ASP:OD1	1:B:276:SER:HB2	1.93	0.68
2:D:290:ALA:HB1	2:D:294:ARG:HH12	1.58	0.68
1:C:293:ASP:O	1:C:297:ARG:HB2	1.94	0.68
2:D:290:ALA:O	2:D:294:ARG:HG2	1.94	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:208:HIS:HD2	3:D:826:HOH:O	1.76	0.67
1:C:262:LEU:HD11	1:C:266:VAL:CG2	2.25	0.67
1:B:246:ARG:O	1:B:247:GLN:HG2	1.94	0.67
2:D:288:ILE:O	2:D:292:LEU:HD13	1.95	0.66
2:D:73:PHE:CE2	2:D:77:LEU:HD11	2.30	0.66
1:C:301:LYS:HD3	2:D:59:ASP:OD2	1.94	0.66
2:D:55:THR:HG23	2:D:57:ASP:N	2.08	0.66
2:D:125:ARG:NH1	2:D:125:ARG:HG2	2.10	0.66
1:B:270:MET:CE	1:B:284:MET:HE1	2.26	0.65
2:D:140:HIS:HE1	2:D:219:ASP:OD2	1.79	0.65
1:C:225:GLU:OE1	1:C:273:ARG:HD3	1.97	0.65
1:C:191:THR:HB	1:C:274:ARG:NH2	2.13	0.64
2:D:286:ASN:ND2	2:D:288:ILE:HG12	2.09	0.63
2:D:73:PHE:O	2:D:77:LEU:HD13	1.99	0.63
2:D:140:HIS:CE1	2:D:221:GLU:HB3	2.34	0.62
1:B:225:GLU:OE1	1:B:273:ARG:HD3	2.00	0.62
1:C:242:ALA:HB2	2:D:283:LEU:HD12	1.81	0.61
1:B:262:LEU:HD11	1:B:266:VAL:HG21	1.82	0.60
2:D:266:ASP:CG	2:D:268:THR:HG22	2.22	0.60
1:C:297:ARG:HG2	1:C:297:ARG:HH11	1.66	0.60
1:C:318:PHE:HZ	3:D:704:HOH:O	1.83	0.60
2:D:278:LEU:HD13	2:D:301:PRO:HD3	1.84	0.59
2:D:55:THR:HG21	3:D:621:HOH:O	2.00	0.59
2:D:280:SER:O	2:D:284:ARG:HG2	2.01	0.59
1:B:262:LEU:HD11	1:B:266:VAL:HG23	1.84	0.59
1:B:200:ASN:HD22	1:B:200:ASN:N	2.01	0.59
2:D:73:PHE:CZ	2:D:77:LEU:HD11	2.38	0.58
2:D:290:ALA:HB1	2:D:294:ARG:NH1	2.18	0.58
1:B:194:LEU:HD21	1:B:274:ARG:HG3	1.86	0.58
1:C:219:VAL:O	1:C:247:GLN:HB2	2.04	0.57
2:D:194:GLU:O	2:D:194:GLU:HG2	2.04	0.57
1:B:199:VAL:HG21	1:B:270:MET:HE1	1.85	0.57
2:D:72:PRO:HB2	3:D:704:HOH:O	2.03	0.57
2:D:240:CYS:O	2:D:271:MET:HB2	2.04	0.56
1:C:308:THR:HG21	2:D:53:TYR:OH	2.06	0.56
1:C:313:MET:HE1	2:D:73:PHE:HB2	1.87	0.55
2:D:91:ASN:HD21	2:D:95:GLN:HE21	1.54	0.55
2:D:212:HIS:ND1	3:D:826:HOH:O	2.33	0.55
1:C:321:PRO:HG3	3:D:704:HOH:O	2.04	0.55
2:D:237:GLU:HB3	3:D:826:HOH:O	2.06	0.55
1:B:233:TRP:CD1	1:B:258:ALA:HB2	2.41	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:219:VAL:O	1:B:247:GLN:HB2	2.06	0.55
1:B:246:ARG:HH11	1:B:246:ARG:HG2	1.72	0.54
2:D:132:HIS:O	2:D:136:ARG:HG3	2.07	0.54
1:B:207:LEU:HD21	1:B:291:ASP:HB2	1.88	0.54
1:C:191:THR:HB	1:C:274:ARG:CZ	2.37	0.54
1:B:218:LYS:HG3	1:B:247:GLN:HG3	1.88	0.54
1:C:193:GLU:HG2	1:C:194:LEU:N	2.22	0.53
2:D:55:THR:HG22	2:D:59:ASP:N	2.20	0.53
1:C:225:GLU:CG	1:C:273:ARG:HB3	2.38	0.53
1:C:245:HIS:NE2	1:B:248:VAL:HG22	2.24	0.53
1:B:297:ARG:HG3	1:B:302:ARG:HB2	1.90	0.52
2:D:204:ASN:HD21	2:D:208:HIS:HB2	1.74	0.52
2:D:300:GLU:HG3	2:D:301:PRO:HD2	1.91	0.52
1:B:214:LEU:C	1:B:214:LEU:HD23	2.29	0.52
1:C:233:TRP:CD1	1:C:258:ALA:HB2	2.46	0.51
1:B:293:ASP:O	1:B:295:ARG:N	2.43	0.51
2:D:120:VAL:HG21	2:D:199:GLN:OE1	2.11	0.51
2:D:66:VAL:HG21	2:D:100:LEU:HB2	1.93	0.51
1:B:297:ARG:NE	1:B:302:ARG:HD2	2.23	0.50
1:C:256:PRO:HG3	3:C:808:HOH:O	2.10	0.50
2:D:286:ASN:HD21	2:D:288:ILE:CG1	2.15	0.50
2:D:237:GLU:CB	3:D:826:HOH:O	2.59	0.49
1:B:231:PRO:HD2	3:B:738:HOH:O	2.12	0.49
2:D:278:LEU:HA	2:D:293:LEU:HD13	1.93	0.49
1:C:253:ARG:HH11	1:C:253:ARG:HG3	1.76	0.49
2:D:240:CYS:HA	2:D:272:TYR:CD1	2.48	0.49
2:D:125:ARG:HH11	2:D:125:ARG:CG	2.24	0.49
2:D:193:ASP:C	2:D:195:ASP:H	2.15	0.49
1:C:225:GLU:HG3	1:C:273:ARG:HB3	1.95	0.48
1:C:306:TYR:HE1	1:C:310:LYS:HE3	1.78	0.48
1:C:246:ARG:HH11	1:C:246:ARG:HG2	1.77	0.48
2:D:219:ASP:O	2:D:223:VAL:HG23	2.14	0.48
2:D:245:LEU:O	2:D:249:VAL:HG23	2.14	0.48
1:C:267:ARG:HD2	1:C:285:GLU:OE1	2.14	0.47
1:B:265:PRO:HB3	1:B:287:GLN:HE21	1.78	0.47
2:D:120:VAL:CG2	2:D:199:GLN:OE1	2.63	0.47
1:B:295:ARG:C	1:B:297:ARG:H	2.18	0.47
1:B:228:PHE:CE2	1:B:270:MET:HG3	2.50	0.47
1:C:277:ASP:OD1	1:C:279:GLU:HB2	2.15	0.47
1:B:199:VAL:HG12	1:B:201:ARG:N	2.25	0.46
2:D:286:ASN:HD22	2:D:287:PRO:CD	2.28	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:216:CYS:O	1:B:248:VAL:HG22	2.15	0.46
1:C:199:VAL:HG13	1:C:212:ILE:HG23	1.98	0.46
1:B:270:MET:HE2	1:B:284:MET:HE1	1.97	0.46
2:D:120:VAL:HG21	2:D:199:GLN:CD	2.36	0.46
1:C:309:PHE:HA	1:C:312:ILE:HD11	1.96	0.46
1:B:293:ASP:HA	1:B:297:ARG:CD	2.45	0.46
2:D:64:LEU:HD23	2:D:67:ILE:HD12	1.98	0.45
2:D:90:GLN:HA	2:D:95:GLN:O	2.16	0.45
1:C:311:SER:OG	1:C:315:LYS:HE2	2.17	0.45
1:C:199:VAL:HG13	1:C:212:ILE:CG2	2.47	0.45
2:D:216:ILE:HD11	2:D:247:LEU:HD13	1.99	0.45
1:C:318:PHE:CE1	1:C:320:GLY:HA2	2.51	0.45
1:B:293:ASP:CG	1:B:297:ARG:HH11	2.19	0.45
1:B:226:VAL:HG11	1:B:252:PHE:CZ	2.51	0.45
1:C:191:THR:HG21	1:C:277:ASP:OD2	2.17	0.45
2:D:143:ALA:O	2:D:147:LEU:HG	2.17	0.45
1:C:297:ARG:HG2	1:C:297:ARG:NH1	2.29	0.45
2:D:249:VAL:O	2:D:252:GLN:NE2	2.49	0.44
2:D:155:ARG:NE	2:D:155:ARG:HA	2.32	0.44
1:C:297:ARG:O	1:C:300:GLU:HG3	2.17	0.44
2:D:124:GLU:OE1	2:D:126:GLY:N	2.51	0.44
1:C:248:VAL:HG22	1:C:248:VAL:O	2.17	0.44
1:B:223:ASP:CG	1:B:276:SER:HB2	2.37	0.44
1:B:270:MET:HE1	1:B:284:MET:HE1	2.00	0.44
2:D:63:HIS:O	2:D:66:VAL:HG22	2.16	0.44
2:D:98:LEU:HD13	2:D:130:ALA:HB2	2.00	0.44
2:D:69:GLN:HA	3:D:715:HOH:O	2.17	0.43
1:C:242:ALA:HB2	2:D:283:LEU:CD1	2.47	0.43
2:D:62:LEU:HD23	2:D:62:LEU:HA	1.87	0.43
2:D:98:LEU:HD13	2:D:130:ALA:CB	2.49	0.43
1:C:211:GLU:HB2	1:C:253:ARG:NH1	2.34	0.43
1:C:314:LYS:HE2	1:C:314:LYS:HB2	1.85	0.43
1:B:257:TYR:CG	1:B:258:ALA:N	2.87	0.42
1:B:229:THR:HG22	1:B:234:GLU:HG2	2.01	0.42
2:D:302:GLU:N	2:D:302:GLU:OE1	2.52	0.42
1:B:199:VAL:HG13	1:B:212:ILE:CG2	2.49	0.42
2:D:257:LEU:HD22	2:D:289:LEU:HD22	2.02	0.42
1:C:293:ASP:CG	1:C:296:HIS:HD2	2.23	0.42
1:C:246:ARG:C	1:C:247:GLN:HG2	2.40	0.41
1:C:214:LEU:C	1:C:214:LEU:HD23	2.41	0.41
1:C:318:PHE:HB2	2:D:76:PHE:CD2	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:286:ASN:HD22	2:D:286:ASN:C	2.22	0.41
1:B:196:ILE:HG13	1:B:272:LEU:HD22	2.02	0.41
1:B:245:HIS:HB3	1:B:249:ALA:HB3	2.03	0.41
1:B:291:ASP:HB3	1:B:295:ARG:CB	2.51	0.41
2:D:286:ASN:HA	2:D:287:PRO:HD3	1.85	0.41
2:D:94:GLY:O	2:D:124:GLU:HA	2.21	0.41
1:B:216:CYS:O	1:B:248:VAL:CG2	2.69	0.41
1:B:292:THR:O	1:B:293:ASP:HB2	2.20	0.41
2:D:218:LYS:HG2	2:D:251:ALA:HB1	2.03	0.41
2:D:286:ASN:HD22	2:D:287:PRO:N	2.18	0.41
1:C:322:THR:HG23	3:C:773:HOH:O	2.21	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	B	110/136 (81%)	99 (90%)	11 (10%)	0	100 100
1	C	133/136 (98%)	123 (92%)	9 (7%)	1 (1%)	19 10
2	D	216/282 (77%)	200 (93%)	14 (6%)	2 (1%)	17 8
All	All	459/554 (83%)	422 (92%)	34 (7%)	3 (1%)	22 12

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	154	PRO
1	C	247	GLN
2	D	285	PRO

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	B	97/123 (79%)	90 (93%)	7 (7%)	14 7
1	C	121/123 (98%)	114 (94%)	7 (6%)	20 11
2	D	165/225 (73%)	156 (94%)	9 (6%)	21 13
All	All	383/471 (81%)	360 (94%)	23 (6%)	19 11

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	240	SER
1	C	253	ARG
1	C	272	LEU
1	C	293	ASP
1	C	297	ARG
1	C	312	ILE
1	C	325	ARG
1	B	200	ASN
1	B	217	ASP
1	B	222	GLU
1	B	236	ARG
1	B	246	ARG
1	B	272	LEU
1	B	292	THR
2	D	78	LEU
2	D	98	LEU
2	D	124	GLU
2	D	194	GLU
2	D	221	GLU
2	D	225	LEU
2	D	233	LEU
2	D	278	LEU
2	D	286	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	247	GLN
1	C	296	HIS
1	B	200	ASN
1	B	287	GLN
2	D	68	HIS
2	D	70	HIS
2	D	95	GLN
2	D	140	HIS
2	D	286	ASN

5.3.3 RNA (i)

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 monosaccharides 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

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^{th} 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	B	112/136 (82%)	0.64	12 (10%) 6 6	11, 23, 87, 95	0
1	C	135/136 (99%)	0.50	15 (11%) 5 5	13, 31, 64, 89	0
2	D	220/282 (78%)	0.37	15 (6%) 17 18	9, 24, 69, 98	0
All	All	467/554 (84%)	0.47	42 (8%) 9 10	9, 26, 76, 98	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	193	ASP	8.8
1	C	191	THR	8.3
1	B	293	ASP	7.9
1	B	299	GLU	7.2
1	B	295	ARG	6.7
2	D	52	GLY	6.1
2	D	50	VAL	6.1
1	B	296	HIS	6.0
2	D	157	ALA	5.9
1	C	321	PRO	5.8
2	D	156	ASP	5.7
1	B	300	GLU	5.5
1	B	297	ARG	5.5
1	B	294	ASP	5.2
2	D	84	HIS	5.2
1	C	319	ASN	5.1
1	C	192	ALA	4.9
2	D	303	ASP	4.7
1	B	302	ARG	4.5
2	D	304	GLY	4.3
1	C	292	THR	4.3
1	B	301	LYS	4.2
2	D	155	ARG	3.8

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Mol	Chain	Res	Type	RSRZ
2	D	51	PHE	3.8
1	B	298	ILE	3.6
1	B	261	SER	3.3
1	C	306	TYR	3.3
1	C	323	GLU	3.1
2	D	80	PHE	3.1
1	C	312	ILE	2.8
1	B	292	THR	2.8
1	C	320	GLY	2.8
2	D	194	GLU	2.6
1	C	322	THR	2.5
1	C	231	PRO	2.5
2	D	152	SER	2.4
1	C	297	ARG	2.3
1	C	316	SER	2.2
2	D	302	GLU	2.1
2	D	53	TYR	2.1
1	C	193	GLU	2.0
1	C	278	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 monosaccharides 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.