Full wwPDB X-ray Structure Validation Report (i)

Feb 6, 2024 – 06:09 pm GMT

PDB ID	:	1QIU
Title	:	A triple beta-spiral in the adenovirus fibre shaft reveals a new structural motif for biological fibres
Authors	:	van Raaij, M.J.; Lavigne, G.; Mitraki, A.; Cusack, S.
Deposited on	:	1999-06-16
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
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) 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 (1)) were used in the production of this report:

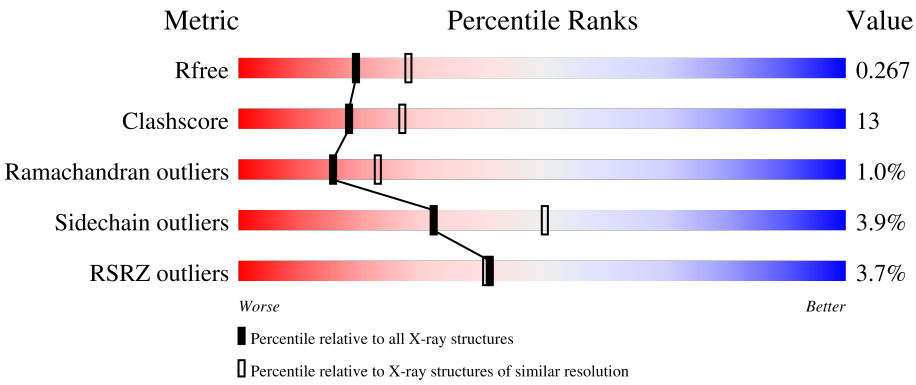
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 2.40 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}	130704	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (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 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	A	264	4% 74% 24% •
1	B	264	4% 69% 30% •
1	C	264	4% 76% 22% •
1	D	264	4% 75% 21% •
1	E	264	4% 71% 26% •

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Mol	Chain	Length	Quality of chain
1	F	264	2% 74% 23% •

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 12621 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 ADENOVIRUS FIBRE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	264	2007	1254	331	413	9	0	0	0
1	B	264	2007	1254	331	413	9	0	0	0
1	C	264	2007	1254	331	413	9	0	0	0
1	D	264	2007	1254	331	413	9	0	0	0
1	E	264	2007	1254	331	413	9	0	0	0
1	F	264	2007	1254	331	413	9	0	0	0

- Molecule 2 is water.

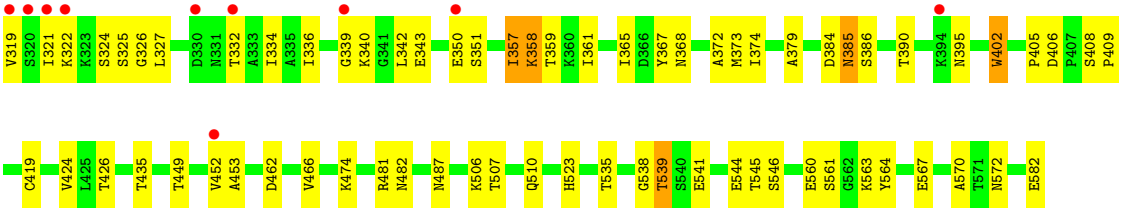
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	99	99	99	0	0
2	B	96	96	96	0	0
2	C	85	85	85	0	0
2	D	95	95	95	0	0
2	E	92	92	92	0	0
2	F	112	112	112	0	0

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.

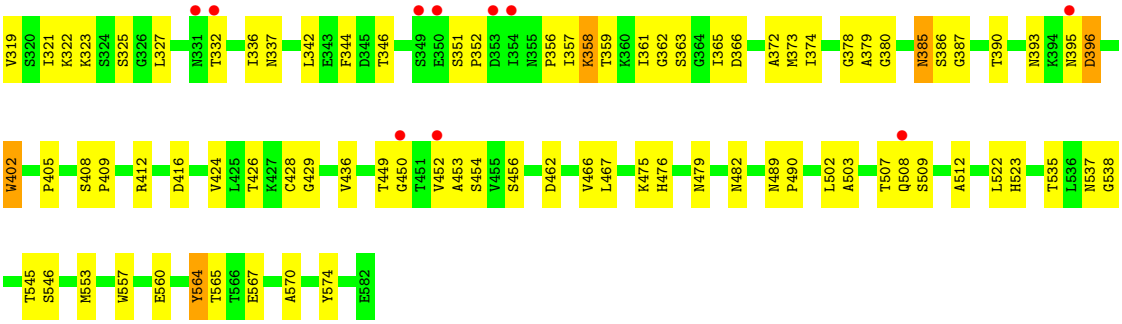
• Molecule 1: ADENOVIRUS FIBRE

Chain A:



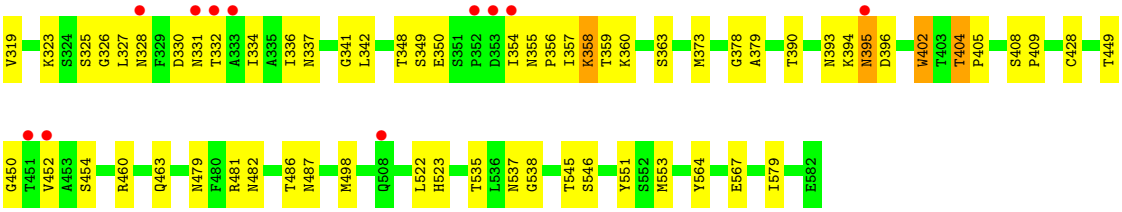
• Molecule 1: ADENOVIRUS FIBRE

Chain B:



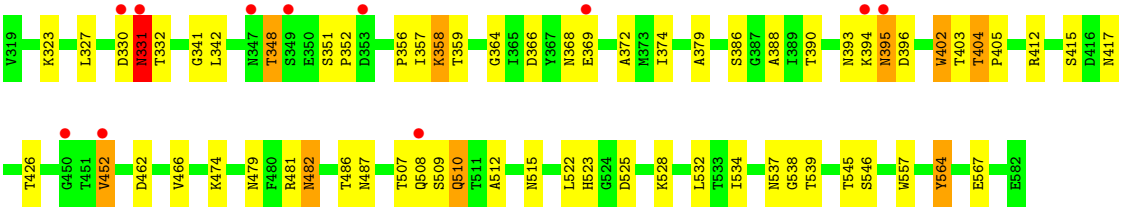
• Molecule 1: ADENOVIRUS FIBRE

Chain C:



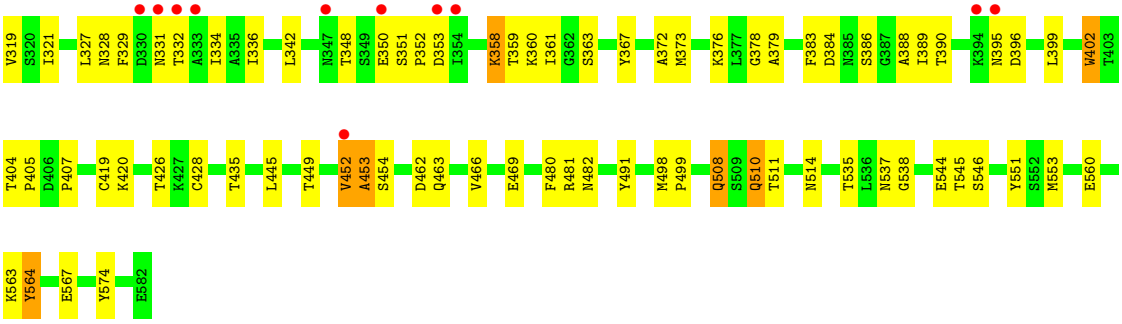
• Molecule 1: ADENOVIRUS FIBRE

Chain D:



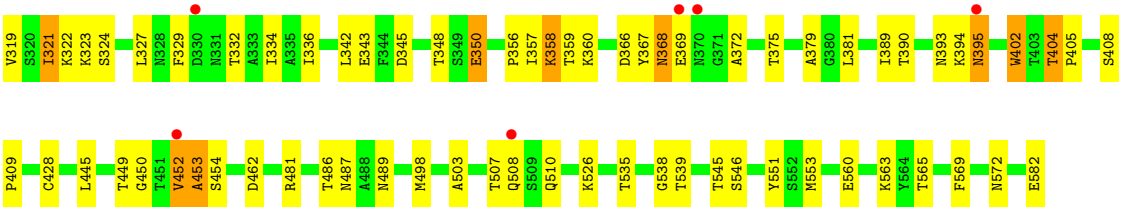
• Molecule 1: ADENOVIRUS FIBRE

Chain E:



• Molecule 1: ADENOVIRUS FIBRE

Chain F:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	165.51Å 95.87Å 211.77Å 90.00° 106.83° 90.00°	Depositor
Resolution (Å)	25.00 – 2.40 25.06 – 2.41	Depositor EDS
% Data completeness (in resolution range)	84.0 (25.00-2.40) 78.0 (25.06-2.41)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.13	Depositor
$< I/\sigma(I) >$	-	Xtriage
Refinement program	CNS 0.4	Depositor
R, R_{free}	0.232 , 0.265 0.232 , 0.267	Depositor DCC
R_{free} test set	827 reflections (0.80%)	wwPDB-VP
Wilson B-factor (Å^2)	(Not available)	Xtriage
Anisotropy	(Not available)	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 22.1	EDS
L-test for twinning[1]	$< \|L\| >$ = (Not available), $< L^2 >$ = (Not available)	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	12621	wwPDB-VP
Average B, all atoms (Å^2)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *(Not available)*

[1]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 (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	0.37	0/2046	0.65	0/2780
1	B	0.37	0/2046	0.64	0/2780
1	C	0.36	0/2046	0.65	0/2780
1	D	0.37	0/2046	0.66	0/2780
1	E	0.37	0/2046	0.65	0/2780
1	F	0.37	0/2046	0.64	0/2780
All	All	0.37	0/12276	0.65	0/16680

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 (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	2007	0	1949	63	0
1	B	2007	0	1949	73	0
1	C	2007	0	1949	60	0
1	D	2007	0	1949	59	0
1	E	2007	0	1949	71	0
1	F	2007	0	1949	68	0
2	A	99	0	0	4	0
2	B	96	0	0	3	0
2	C	85	0	0	0	0
2	D	95	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	92	0	0	1	0
2	F	112	0	0	7	0
All	All	12621	0	11694	317	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 13.

All (317) 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:321:ILE:HD11	1:B:327:LEU:HB3	1.45	0.98
1:A:321:ILE:HD11	1:A:334:ILE:HG12	1.51	0.92
1:B:336:ILE:HD11	1:C:326:GLY:HA3	1.47	0.92
1:A:336:ILE:HD11	1:B:336:ILE:HG22	1.51	0.91
1:A:560:GLU:HB2	1:A:563:LYS:HD2	1.56	0.88
1:C:350:GLU:HB2	1:C:354:ILE:HB	1.56	0.86
1:A:510:GLN:HE22	1:A:539:THR:HG21	1.41	0.84
1:F:560:GLU:HB2	1:F:563:LYS:HD2	1.62	0.81
1:A:326:GLY:HA3	1:C:336:ILE:HD11	1.63	0.80
1:D:417:ASN:HD21	1:D:474:LYS:HD2	1.52	0.74
1:E:462:ASP:OD2	1:E:466:VAL:HB	1.87	0.74
1:C:452:VAL:HG12	1:C:454:SER:H	1.51	0.74
1:B:452:VAL:HA	2:B:2031:HOH:O	1.88	0.73
1:A:368:ASN:HD21	1:A:372:ALA:HB3	1.54	0.73
1:F:449:THR:HB	1:F:452:VAL:HB	1.72	0.72
1:C:319:VAL:HG11	1:C:334:ILE:HD11	1.73	0.71
1:A:321:ILE:HD11	1:A:334:ILE:CG1	2.21	0.70
1:D:510:GLN:HA	1:D:510:GLN:HE21	1.57	0.70
1:B:359:THR:HG23	1:C:342:LEU:HD13	1.73	0.69
1:F:445:LEU:HD21	1:F:569:PHE:HB2	1.74	0.68
1:D:379:ALA:N	1:F:390:THR:HG22	2.08	0.68
1:F:342:LEU:O	1:F:343:GLU:HG3	1.93	0.67
1:E:390:THR:HG22	1:F:379:ALA:N	2.10	0.67
1:A:402:TRP:CZ2	1:A:481:ARG:HG3	2.31	0.66
1:F:322:LYS:HE3	1:F:324:SER:OG	1.95	0.66
1:F:321:ILE:HD11	1:F:334:ILE:HG23	1.77	0.66
1:B:356:PRO:HB3	1:C:337:ASN:ND2	2.10	0.66
1:E:402:TRP:CZ2	1:E:481:ARG:HG3	2.30	0.66
1:A:379:ALA:N	1:C:390:THR:HG22	2.12	0.65
1:A:545:THR:O	1:A:546:SER:HB2	1.95	0.65
1:C:348:THR:HG22	1:C:350:GLU:H	1.60	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:449:THR:HB	1:E:452:VAL:HB	1.78	0.64
1:D:393:ASN:HB2	2:D:2015:HOH:O	1.96	0.64
1:E:321:ILE:HD11	1:E:334:ILE:HG12	1.78	0.64
1:E:564:TYR:HB3	1:E:567:GLU:HG3	1.78	0.64
1:E:361:ILE:HD12	1:E:367:TYR:CZ	2.33	0.64
1:D:394:LYS:O	1:D:395:ASN:HB2	1.97	0.64
1:B:374:ILE:HG22	1:C:363:SER:N	2.14	0.62
1:F:332:THR:O	1:F:332:THR:HG22	2.00	0.62
1:B:361:ILE:HD11	1:B:373:MET:SD	2.40	0.61
1:D:390:THR:HG22	1:E:379:ALA:N	2.14	0.61
1:D:402:TRP:CH2	1:D:405:PRO:HD3	2.35	0.61
1:D:404:THR:HG22	1:D:486:THR:HG23	1.82	0.61
1:A:449:THR:HB	1:A:452:VAL:O	2.00	0.61
1:F:368:ASN:ND2	1:F:369:GLU:H	1.97	0.61
1:E:348:THR:HG22	1:E:350:GLU:H	1.66	0.61
1:F:452:VAL:HG12	1:F:454:SER:H	1.64	0.61
1:E:560:GLU:HB3	1:E:563:LYS:HD2	1.83	0.61
1:B:449:THR:HB	1:B:452:VAL:HB	1.83	0.60
1:C:332:THR:O	1:C:332:THR:HG22	2.01	0.60
1:F:368:ASN:HD22	1:F:369:GLU:H	1.48	0.60
1:B:358:LYS:HD2	1:B:359:THR:O	2.02	0.60
1:D:357:ILE:O	1:E:342:LEU:HD12	2.02	0.60
1:F:452:VAL:HG12	1:F:454:SER:N	2.15	0.60
1:F:452:VAL:HA	2:F:2045:HOH:O	1.99	0.60
1:B:336:ILE:HG12	1:C:325:SER:O	2.01	0.60
1:A:342:LEU:HD12	1:C:358:LYS:HA	1.82	0.60
1:D:368:ASN:OD1	1:D:369:GLU:N	2.35	0.59
1:E:510:GLN:HA	1:E:510:GLN:HE21	1.67	0.59
1:A:332:THR:O	1:A:332:THR:HG22	2.02	0.59
1:B:361:ILE:HG23	1:B:365:ILE:HB	1.83	0.59
1:A:510:GLN:HE22	1:A:539:THR:CG2	2.15	0.59
1:A:319:VAL:HG13	1:A:334:ILE:HD11	1.84	0.58
1:A:507:THR:H	1:A:538:GLY:HA2	1.68	0.58
1:C:393:ASN:HB3	1:C:396:ASP:O	2.03	0.58
1:C:449:THR:HB	1:C:452:VAL:O	2.03	0.58
1:F:359:THR:HG23	1:F:367:TYR:OH	2.03	0.58
1:A:390:THR:HG22	1:B:378:GLY:HA3	1.85	0.58
1:A:426:THR:HG21	1:B:428:CYS:O	2.04	0.58
1:A:324:SER:HA	1:C:355:ASN:HB2	1.85	0.57
1:A:336:ILE:HD11	1:B:336:ILE:CG2	2.29	0.57
1:F:452:VAL:O	1:F:453:ALA:HB3	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:336:ILE:HG23	1:C:327:LEU:HD11	1.86	0.57
1:F:375:THR:HG21	1:F:389:ILE:HD11	1.86	0.57
1:A:510:GLN:NE2	1:A:539:THR:OG1	2.38	0.57
1:E:545:THR:O	1:E:546:SER:HB2	2.05	0.57
1:E:327:LEU:HD23	1:E:336:ILE:HA	1.86	0.56
1:E:342:LEU:HD23	1:E:359:THR:HA	1.87	0.56
1:A:523:HIS:CD2	1:A:570:ALA:HB3	2.40	0.56
1:B:545:THR:O	1:B:546:SER:HB2	2.06	0.56
1:E:510:GLN:HE21	1:E:510:GLN:CA	2.16	0.56
1:A:358:LYS:HD3	1:A:359:THR:O	2.06	0.56
1:D:372:ALA:HA	1:E:360:LYS:HG3	1.88	0.56
1:D:379:ALA:H	1:F:390:THR:HG22	1.71	0.55
1:D:510:GLN:HE21	1:D:510:GLN:CA	2.19	0.55
1:A:357:ILE:O	1:B:342:LEU:HD12	2.06	0.55
1:E:452:VAL:O	1:E:453:ALA:HB3	2.05	0.55
1:B:462:ASP:OD2	1:B:466:VAL:HB	2.07	0.55
1:C:402:TRP:CH2	1:C:405:PRO:HD3	2.41	0.55
1:E:508:GLN:CD	1:E:508:GLN:H	2.09	0.55
1:D:374:ILE:HG22	1:E:363:SER:N	2.21	0.54
1:E:361:ILE:HD11	1:E:373:MET:HG2	1.90	0.54
1:D:510:GLN:HG2	1:D:539:THR:HG21	1.89	0.54
1:D:374:ILE:HG22	1:E:363:SER:H	1.71	0.54
1:A:321:ILE:CD1	1:A:334:ILE:HG12	2.33	0.54
1:A:506:LYS:HG3	1:A:541:GLU:HG3	1.90	0.54
1:D:390:THR:HG22	1:E:379:ALA:H	1.72	0.54
1:A:340:LYS:O	1:C:358:LYS:HB2	2.08	0.54
1:F:319:VAL:HA	2:F:2002:HOH:O	2.08	0.54
1:C:545:THR:O	1:C:546:SER:HB2	2.09	0.53
1:A:405:PRO:HB2	1:B:503:ALA:HB2	1.90	0.53
1:B:467:LEU:O	1:B:475:LYS:HE3	2.08	0.53
1:A:385:ASN:HD22	1:A:386:SER:N	2.07	0.53
1:C:522:LEU:HG	1:C:523:HIS:CD2	2.44	0.53
1:C:349:SER:HB3	1:C:350:GLU:OE2	2.08	0.53
1:E:358:LYS:HA	1:F:342:LEU:HD12	1.90	0.53
1:B:323:LYS:HA	1:B:327:LEU:O	2.09	0.52
1:A:325:SER:O	1:C:336:ILE:HG12	2.09	0.52
1:A:384:ASP:OD2	1:A:386:SER:HB3	2.08	0.52
1:B:390:THR:HG22	1:C:379:ALA:N	2.25	0.52
1:D:388:ALA:HB2	1:E:376:LYS:HG2	1.92	0.52
1:A:343:GLU:OE2	1:A:358:LYS:HD2	2.09	0.52
1:B:452:VAL:O	1:B:453:ALA:HB3	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:321:ILE:HD11	1:E:334:ILE:CG1	2.39	0.52
1:D:452:VAL:HG13	2:D:2025:HOH:O	2.09	0.52
1:D:357:ILE:O	1:D:357:ILE:HG23	2.09	0.51
1:F:560:GLU:HB2	1:F:563:LYS:CD	2.35	0.51
1:B:327:LEU:HD23	1:B:336:ILE:HG22	1.92	0.51
1:C:330:ASP:O	1:C:331:ASN:HB3	2.10	0.51
1:F:366:ASP:OD1	1:F:367:TYR:N	2.44	0.51
1:A:535:THR:HG21	1:A:539:THR:HG22	1.90	0.51
1:F:327:LEU:HD23	1:F:336:ILE:HA	1.93	0.51
1:A:368:ASN:ND2	1:A:372:ALA:HB3	2.22	0.51
1:A:453:ALA:HB2	1:A:561:SER:HA	1.92	0.51
1:A:462:ASP:OD2	1:A:466:VAL:HB	2.11	0.51
1:C:498:MET:HG3	1:C:551:TYR:CD2	2.46	0.51
1:A:406:ASP:HB2	1:B:502:LEU:HB3	1.92	0.51
1:C:450:GLY:C	1:C:452:VAL:H	2.14	0.51
1:B:402:TRP:CH2	1:B:405:PRO:HD3	2.46	0.51
1:D:479:ASN:ND2	1:D:487:ASN:HB3	2.26	0.51
1:A:359:THR:HG21	1:A:373:MET:SD	2.50	0.50
1:E:449:THR:O	1:E:452:VAL:HG23	2.11	0.50
1:E:560:GLU:O	1:E:563:LYS:HB2	2.11	0.50
1:F:404:THR:HG22	1:F:486:THR:HG23	1.93	0.50
1:A:582:GLU:HB3	2:A:2097:HOH:O	2.10	0.50
1:D:358:LYS:HD2	1:D:359:THR:O	2.10	0.50
1:A:390:THR:HG22	1:B:379:ALA:N	2.26	0.50
1:A:564:TYR:HB3	1:A:567:GLU:HG3	1.92	0.50
1:B:436:VAL:HG12	1:B:574:TYR:HB3	1.93	0.50
1:C:402:TRP:CZ2	1:C:481:ARG:HG3	2.46	0.50
1:D:342:LEU:HD12	1:F:357:ILE:O	2.12	0.50
1:B:351:SER:HB2	1:B:352:PRO:HA	1.94	0.49
1:D:417:ASN:ND2	1:D:474:LYS:HD2	2.25	0.49
1:D:512:ALA:O	1:F:572:ASN:HA	2.12	0.49
1:F:445:LEU:CD2	1:F:569:PHE:HB2	2.41	0.49
1:A:327:LEU:HD21	1:C:336:ILE:HG23	1.94	0.49
1:B:450:GLY:O	1:B:452:VAL:HG23	2.12	0.49
1:C:404:THR:HG22	1:C:486:THR:HG23	1.93	0.49
1:D:402:TRP:CZ2	1:D:481:ARG:HG3	2.48	0.49
1:D:545:THR:O	1:D:546:SER:HB2	2.13	0.49
1:E:445:LEU:HD12	1:E:564:TYR:O	2.12	0.49
1:B:321:ILE:HG22	1:F:489:ASN:OD1	2.11	0.49
1:B:380:GLY:H	1:B:393:ASN:ND2	2.11	0.49
1:D:394:LYS:O	1:D:395:ASN:CB	2.58	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:328:ASN:OD1	1:E:329:PHE:N	2.39	0.49
1:D:532:LEU:HD13	1:D:557:TRP:CE2	2.48	0.49
1:D:364:GLY:HA3	1:F:375:THR:HG22	1.93	0.48
1:E:481:ARG:HG2	1:E:482:ASN:N	2.27	0.48
1:D:323:LYS:HA	1:D:327:LEU:O	2.14	0.48
1:F:348:THR:HG22	1:F:350:GLU:H	1.78	0.48
1:B:535:THR:O	1:B:553:MET:HA	2.13	0.48
1:B:332:THR:HG22	1:B:332:THR:O	2.14	0.48
1:D:341:GLY:HA3	1:F:359:THR:HG22	1.95	0.48
1:D:372:ALA:CB	1:E:360:LYS:HG3	2.43	0.48
1:D:348:THR:HG21	1:D:356:PRO:CD	2.43	0.47
1:D:426:THR:HG21	1:E:428:CYS:O	2.13	0.47
1:A:507:THR:N	1:A:538:GLY:HA2	2.28	0.47
1:D:330:ASP:O	1:D:331:ASN:CB	2.61	0.47
1:A:374:ILE:HG22	1:B:363:SER:N	2.29	0.47
1:D:341:GLY:HA3	1:F:359:THR:CG2	2.45	0.47
1:F:323:LYS:HA	1:F:327:LEU:O	2.13	0.47
1:F:453:ALA:HB3	2:F:2044:HOH:O	2.14	0.47
1:A:395:ASN:HB2	2:A:2015:HOH:O	2.14	0.47
1:F:402:TRP:CZ2	1:F:481:ARG:HG3	2.50	0.47
1:E:402:TRP:CH2	1:E:405:PRO:HD3	2.50	0.47
1:F:394:LYS:HB3	2:F:2060:HOH:O	2.14	0.47
1:D:386:SER:HB3	1:E:376:LYS:NZ	2.30	0.46
1:B:319:VAL:HG21	1:C:319:VAL:HG22	1.97	0.46
1:B:372:ALA:HA	1:C:360:LYS:HG2	1.97	0.46
1:C:479:ASN:HB3	1:C:486:THR:HB	1.98	0.46
1:D:462:ASP:OD2	1:D:466:VAL:HB	2.16	0.46
1:B:564:TYR:HB3	1:B:567:GLU:HG3	1.97	0.46
1:A:334:ILE:HG22	1:B:327:LEU:CD1	2.45	0.46
1:B:322:LYS:HD3	2:B:2001:HOH:O	2.15	0.46
1:C:408:SER:O	1:C:409:PRO:C	2.53	0.46
1:B:374:ILE:HG22	1:C:363:SER:H	1.81	0.46
1:C:323:LYS:HA	1:C:327:LEU:O	2.14	0.46
1:B:408:SER:O	1:B:409:PRO:C	2.52	0.46
1:F:510:GLN:HG2	1:F:539:THR:HG21	1.98	0.46
1:B:359:THR:HG23	1:C:342:LEU:CD1	2.42	0.45
1:D:390:THR:HG22	1:E:378:GLY:HA3	1.98	0.45
1:D:510:GLN:HA	1:D:510:GLN:NE2	2.28	0.45
1:F:545:THR:O	1:F:546:SER:HB2	2.16	0.45
1:D:507:THR:HG23	1:D:509:SER:O	2.17	0.45
1:D:393:ASN:HB3	1:D:396:ASP:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:452:VAL:O	1:E:453:ALA:CB	2.64	0.45
1:F:394:LYS:O	1:F:395:ASN:O	2.35	0.45
1:A:334:ILE:O	1:B:327:LEU:HD12	2.17	0.45
1:D:327:LEU:HG	1:F:336:ILE:HG13	1.98	0.45
1:F:402:TRP:CH2	1:F:405:PRO:HD3	2.51	0.45
1:A:374:ILE:HG22	1:B:363:SER:H	1.81	0.45
1:E:480:PHE:CE2	1:E:491:TYR:HB2	2.52	0.45
1:A:321:ILE:HD11	1:A:334:ILE:CD1	2.47	0.45
1:F:357:ILE:O	1:F:357:ILE:HG23	2.17	0.45
1:F:408:SER:O	1:F:409:PRO:C	2.55	0.45
1:A:339:GLY:N	1:C:357:ILE:O	2.46	0.44
1:A:474:LYS:HE2	2:A:2043:HOH:O	2.17	0.44
1:E:384:ASP:OD2	1:E:388:ALA:HB3	2.17	0.44
1:F:321:ILE:HB	1:F:329:PHE:CZ	2.52	0.44
1:F:452:VAL:O	1:F:453:ALA:CB	2.65	0.44
1:E:361:ILE:HD12	1:E:367:TYR:CE1	2.52	0.44
1:E:351:SER:HB3	1:E:352:PRO:HA	1.99	0.44
1:B:522:LEU:O	1:B:523:HIS:HB2	2.18	0.44
1:E:390:THR:HG22	1:F:379:ALA:H	1.81	0.44
1:E:348:THR:HG21	2:E:2008:HOH:O	2.18	0.44
1:F:498:MET:HG3	1:F:551:TYR:CD2	2.53	0.44
1:A:322:LYS:HG3	1:C:330:ASP:OD2	2.18	0.44
1:B:523:HIS:CE1	1:B:570:ALA:HB3	2.53	0.44
1:D:372:ALA:CA	1:E:360:LYS:HG3	2.48	0.44
1:F:582:GLU:OXT	1:F:582:GLU:HG3	2.17	0.44
1:E:321:ILE:HD12	1:E:329:PHE:CZ	2.52	0.44
1:E:404:THR:O	1:E:407:PRO:HD3	2.18	0.44
1:E:498:MET:HG3	1:E:551:TYR:CD2	2.53	0.43
1:A:408:SER:O	1:A:409:PRO:C	2.56	0.43
1:B:412:ARG:HD2	1:B:416:ASP:OD1	2.18	0.43
1:D:481:ARG:HG2	1:D:482:ASN:N	2.33	0.43
1:E:510:GLN:HA	1:E:510:GLN:NE2	2.30	0.43
1:C:537:ASN:O	1:C:538:GLY:C	2.55	0.43
1:D:330:ASP:O	1:D:331:ASN:HB2	2.18	0.43
1:D:522:LEU:HG	1:D:523:HIS:CD2	2.54	0.43
1:E:407:PRO:O	1:E:420:LYS:HD3	2.18	0.43
1:F:321:ILE:HD12	1:F:334:ILE:HG12	1.99	0.43
1:A:572:ASN:HA	1:B:512:ALA:O	2.19	0.43
1:C:482:ASN:HB2	1:C:487:ASN:HD22	1.82	0.43
1:C:358:LYS:HD2	1:C:359:THR:O	2.18	0.43
1:D:545:THR:O	1:D:546:SER:CB	2.66	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:450:GLY:C	1:F:452:VAL:H	2.21	0.43
1:A:321:ILE:CG2	1:A:327:LEU:HB2	2.49	0.43
1:B:560:GLU:HB3	2:B:2089:HOH:O	2.19	0.43
1:E:399:LEU:HD23	1:E:399:LEU:HA	1.84	0.43
1:C:463:GLN:CD	1:C:463:GLN:H	2.22	0.43
1:E:426:THR:HG21	1:F:428:CYS:O	2.18	0.43
1:E:511:THR:O	1:E:514:ASN:HB2	2.18	0.43
1:F:507:THR:HG22	1:F:538:GLY:HA2	2.00	0.43
1:C:394:LYS:O	1:C:395:ASN:HB2	2.19	0.43
1:D:412:ARG:HG3	1:D:415:SER:O	2.18	0.43
1:E:498:MET:HA	1:E:499:PRO:HD3	1.88	0.43
1:F:345:ASP:HB3	1:F:356:PRO:HD2	2.01	0.43
1:A:357:ILE:H	1:A:357:ILE:HG13	1.65	0.43
1:C:564:TYR:HB3	1:C:567:GLU:HG3	2.00	0.43
1:E:405:PRO:HB2	1:F:503:ALA:HB2	2.01	0.43
1:B:426:THR:HG21	1:C:428:CYS:O	2.19	0.42
1:F:535:THR:O	1:F:553:MET:HA	2.19	0.42
1:E:319:VAL:HG21	1:F:319:VAL:HG22	2.01	0.42
1:E:359:THR:HG23	1:F:342:LEU:HG	2.01	0.42
1:E:372:ALA:HA	1:F:360:LYS:HG3	2.01	0.42
1:D:390:THR:CG2	1:E:378:GLY:HA3	2.49	0.42
1:A:322:LYS:HD3	1:A:322:LYS:HA	1.84	0.42
1:A:424:VAL:HB	1:A:435:THR:HG22	2.01	0.42
1:F:449:THR:O	1:F:452:VAL:HG23	2.19	0.42
1:F:462:ASP:HB2	2:F:2049:HOH:O	2.19	0.42
2:A:2018:HOH:O	1:B:429:GLY:HA3	2.19	0.42
1:C:341:GLY:O	1:C:342:LEU:HD12	2.19	0.42
1:E:351:SER:HA	1:E:353:ASP:N	2.34	0.42
1:E:389:ILE:O	1:F:381:LEU:HD12	2.19	0.42
1:E:537:ASN:O	1:E:538:GLY:C	2.57	0.42
1:B:319:VAL:CG2	1:C:319:VAL:HG22	2.50	0.42
1:B:424:VAL:HG21	1:C:579:ILE:CD1	2.50	0.42
1:F:368:ASN:N	1:F:372:ALA:O	2.51	0.42
1:A:334:ILE:HG22	1:B:327:LEU:HD11	2.01	0.42
1:D:403:THR:O	1:D:404:THR:OG1	2.33	0.42
1:F:368:ASN:ND2	1:F:369:GLU:N	2.65	0.42
1:B:507:THR:OG1	1:B:509:SER:HB3	2.20	0.42
1:E:435:THR:HA	1:E:574:TYR:O	2.19	0.42
1:B:454:SER:HA	1:B:557:TRP:O	2.19	0.42
1:D:525:ASP:OD2	1:D:528:LYS:HE2	2.19	0.42
1:D:564:TYR:HB3	1:D:567:GLU:HG3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:350:GLU:HB3	2:F:2013:HOH:O	2.19	0.42
1:B:344:PHE:O	1:B:346:THR:HG23	2.19	0.42
1:B:390:THR:HG22	1:C:378:GLY:HA3	2.02	0.42
1:B:395:ASN:O	1:B:396:ASP:HB2	2.20	0.42
1:D:372:ALA:HB2	1:E:360:LYS:HG3	2.02	0.41
1:F:368:ASN:HD22	1:F:369:GLU:N	2.14	0.41
1:B:385:ASN:C	1:B:387:GLY:H	2.24	0.41
1:F:358:LYS:HD2	1:F:358:LYS:C	2.40	0.41
1:B:336:ILE:HG23	1:C:327:LEU:CD1	2.50	0.41
1:B:537:ASN:O	1:B:538:GLY:C	2.57	0.41
1:B:424:VAL:HG21	1:C:579:ILE:HD11	2.01	0.41
1:D:515:ASN:HA	1:D:534:ILE:O	2.20	0.41
1:E:383:PHE:HA	1:E:388:ALA:O	2.20	0.41
1:D:351:SER:HB3	1:D:352:PRO:HA	2.02	0.41
1:C:394:LYS:H	1:C:394:LYS:HG2	1.46	0.41
1:C:450:GLY:C	1:C:452:VAL:N	2.74	0.41
1:D:332:THR:O	1:D:332:THR:HG22	2.20	0.41
1:C:326:GLY:O	1:C:327:LEU:HD12	2.20	0.41
1:C:535:THR:O	1:C:553:MET:HA	2.20	0.41
1:A:379:ALA:H	1:C:390:THR:HG22	1.85	0.41
1:D:537:ASN:O	1:D:538:GLY:C	2.59	0.41
1:A:361:ILE:HD11	1:A:367:TYR:CE2	2.56	0.41
1:B:321:ILE:CD1	1:B:327:LEU:HB3	2.33	0.41
1:B:336:ILE:HD12	1:B:357:ILE:HD12	2.03	0.41
1:B:452:VAL:HG12	1:B:454:SER:N	2.35	0.41
1:B:545:THR:O	1:B:546:SER:CB	2.67	0.41
1:E:321:ILE:HD11	1:E:334:ILE:CD1	2.51	0.41
1:E:452:VAL:HG12	1:E:454:SER:N	2.35	0.41
1:F:526:LYS:HG2	2:F:2095:HOH:O	2.20	0.41
1:B:323:LYS:C	1:B:325:SER:H	2.24	0.41
1:C:354:ILE:HG22	1:C:356:PRO:HD3	2.02	0.41
1:A:365:ILE:HD12	1:C:373:MET:O	2.21	0.40
1:B:361:ILE:HG22	1:B:362:GLY:N	2.36	0.40
1:B:479:ASN:HD21	1:B:490:PRO:HA	1.86	0.40
1:E:463:GLN:CD	1:E:463:GLN:H	2.25	0.40
1:A:351:SER:O	1:E:386:SER:CB	2.70	0.40
1:E:331:ASN:O	1:E:332:THR:HB	2.21	0.40
1:E:372:ALA:HA	1:F:360:LYS:CG	2.52	0.40
1:E:535:THR:O	1:E:553:MET:HA	2.21	0.40
1:B:336:ILE:CG1	1:C:325:SER:O	2.67	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	262/264 (99%)	241 (92%)	21 (8%)	0	100	100
1	B	262/264 (99%)	230 (88%)	30 (12%)	2 (1%)	19	29
1	C	262/264 (99%)	234 (89%)	26 (10%)	2 (1%)	19	29
1	D	262/264 (99%)	235 (90%)	23 (9%)	4 (2%)	10	14
1	E	262/264 (99%)	238 (91%)	21 (8%)	3 (1%)	14	20
1	F	262/264 (99%)	242 (92%)	16 (6%)	4 (2%)	10	14
All	All	1572/1584 (99%)	1420 (90%)	137 (9%)	15 (1%)	15	23

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	395	ASN
1	E	452	VAL
1	F	395	ASN
1	D	331	ASN
1	D	395	ASN
1	E	453	ALA
1	F	452	VAL
1	B	396	ASP
1	F	453	ALA
1	B	386	SER
1	C	395	ASN
1	D	348	THR
1	D	404	THR
1	C	404	THR
1	F	404	THR

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	228/228 (100%)	218 (96%)	10 (4%)	28	45
1	B	228/228 (100%)	216 (95%)	12 (5%)	22	37
1	C	228/228 (100%)	224 (98%)	4 (2%)	59	76
1	D	228/228 (100%)	219 (96%)	9 (4%)	32	50
1	E	228/228 (100%)	219 (96%)	9 (4%)	32	50
1	F	228/228 (100%)	219 (96%)	9 (4%)	32	50
All	All	1368/1368 (100%)	1315 (96%)	53 (4%)	32	50

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	350	GLU
1	A	357	ILE
1	A	358	LYS
1	A	385	ASN
1	A	402	TRP
1	A	419	CYS
1	A	482	ASN
1	A	487	ASN
1	A	539	THR
1	A	544	GLU
1	B	337	ASN
1	B	358	LYS
1	B	366	ASP
1	B	385	ASN
1	B	402	TRP
1	B	456	SER
1	B	476	HIS
1	B	482	ASN
1	B	489	ASN
1	B	508	GLN
1	B	564	TYR
1	B	565	THR

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Mol	Chain	Res	Type
1	C	328	ASN
1	C	358	LYS
1	C	402	TRP
1	C	460	ARG
1	D	331	ASN
1	D	358	LYS
1	D	366	ASP
1	D	402	TRP
1	D	452	VAL
1	D	482	ASN
1	D	508	GLN
1	D	510	GLN
1	D	564	TYR
1	E	358	LYS
1	E	396	ASP
1	E	402	TRP
1	E	419	CYS
1	E	469	GLU
1	E	508	GLN
1	E	510	GLN
1	E	544	GLU
1	E	564	TYR
1	F	321	ILE
1	F	350	GLU
1	F	358	LYS
1	F	368	ASN
1	F	393	ASN
1	F	402	TRP
1	F	487	ASN
1	F	508	GLN
1	F	565	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (35) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	385	ASN
1	A	482	ASN
1	A	510	GLN
1	A	523	HIS
1	B	331	ASN
1	B	337	ASN
1	B	370	ASN

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Mol	Chain	Res	Type
1	B	393	ASN
1	B	464	ASN
1	B	482	ASN
1	C	328	ASN
1	C	337	ASN
1	C	355	ASN
1	C	370	ASN
1	C	464	ASN
1	C	476	HIS
1	C	487	ASN
1	C	515	ASN
1	D	337	ASN
1	D	393	ASN
1	D	417	ASN
1	D	482	ASN
1	D	508	GLN
1	D	510	GLN
1	D	515	ASN
1	E	370	ASN
1	E	395	ASN
1	E	464	ASN
1	E	510	GLN
1	E	515	ASN
1	F	337	ASN
1	F	368	ASN
1	F	385	ASN
1	F	464	ASN
1	F	515	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 (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	264/264 (100%)	-0.08	10 (3%)	40	39	7, 29, 68, 89	0
1	B	264/264 (100%)	-0.06	10 (3%)	40	39	10, 31, 68, 87	0
1	C	264/264 (100%)	-0.00	11 (4%)	36	35	11, 30, 75, 89	0
1	D	264/264 (100%)	-0.06	11 (4%)	36	35	9, 31, 68, 90	0
1	E	264/264 (100%)	-0.02	11 (4%)	36	35	10, 31, 75, 89	0
1	F	264/264 (100%)	-0.12	6 (2%)	60	58	10, 30, 66, 76	0
All	All	1584/1584 (100%)	-0.06	59 (3%)	41	41	7, 30, 72, 90	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	395	ASN	4.4
1	D	349	SER	4.2
1	A	321	ILE	4.0
1	C	352	PRO	3.8
1	D	394	LYS	3.8
1	A	394	LYS	3.6
1	B	508	GLN	3.6
1	C	452	VAL	3.5
1	C	332	THR	3.5
1	A	452	VAL	3.5
1	E	332	THR	3.4
1	A	350	GLU	3.4
1	D	508	GLN	3.4
1	C	331	ASN	3.4
1	C	353	ASP	3.4
1	D	452	VAL	3.4
1	B	450	GLY	3.3
1	A	322	LYS	3.3
1	E	331	ASN	3.2

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Mol	Chain	Res	Type	RSRZ
1	D	353	ASP	3.2
1	D	450	GLY	3.2
1	C	333	ALA	3.1
1	E	394	LYS	3.1
1	C	354	ILE	3.1
1	B	353	ASP	3.1
1	E	395	ASN	3.1
1	F	508	GLN	3.1
1	E	353	ASP	3.0
1	F	395	ASN	2.9
1	E	354	ILE	2.8
1	A	332	THR	2.8
1	F	369	GLU	2.7
1	C	451	THR	2.7
1	E	350	GLU	2.6
1	B	349	SER	2.6
1	F	370	ASN	2.6
1	F	452	VAL	2.6
1	A	320	SER	2.5
1	B	452	VAL	2.5
1	E	333	ALA	2.5
1	C	328	ASN	2.5
1	C	508	GLN	2.5
1	E	452	VAL	2.4
1	B	331	ASN	2.4
1	D	330	ASP	2.4
1	B	354	ILE	2.4
1	E	330	ASP	2.4
1	C	395	ASN	2.3
1	F	330	ASP	2.3
1	D	347	ASN	2.2
1	D	331	ASN	2.1
1	B	350	GLU	2.1
1	A	339	GLY	2.1
1	D	369	GLU	2.1
1	D	395	ASN	2.1
1	E	347	ASN	2.0
1	A	330	ASP	2.0
1	B	332	THR	2.0
1	A	319	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 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.