Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 01:22 pm BST

PDB ID	:	1FLK
Title	:	MOLECULAR BASIS FOR CD40 SIGNALING MEDIATED BY TRAF3
Authors	:	Ni, C.-Z.; Welsh, K.; Leo, E.; Chiou, C.-K.; Wu, H.; Reed, J.C.; Ely, K.R.
Deposited on	:	2000-08-14
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 ①) were used in the production of this report:

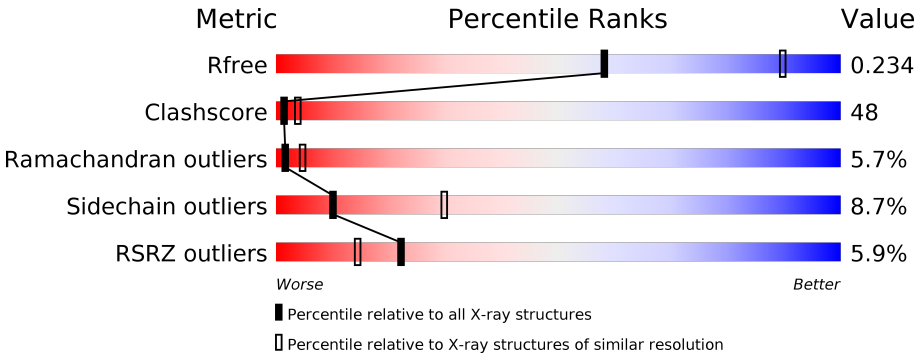
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation 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 <=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	228	5%, 36%, 43%, 9%, •, 10%
1	B	228	5%, 32%, 49%, 9%, 10%

2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 3252 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 3.

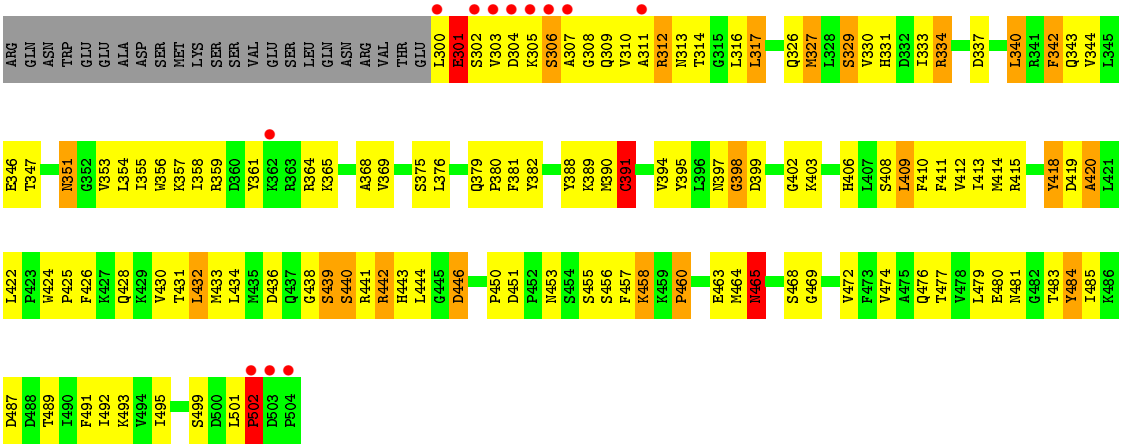
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	205	Total	C	N	O	S	0	0	0
			1626	1035	277	303	11			
1	B	205	Total	C	N	O	S	0	0	0
			1626	1035	277	303	11			

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.

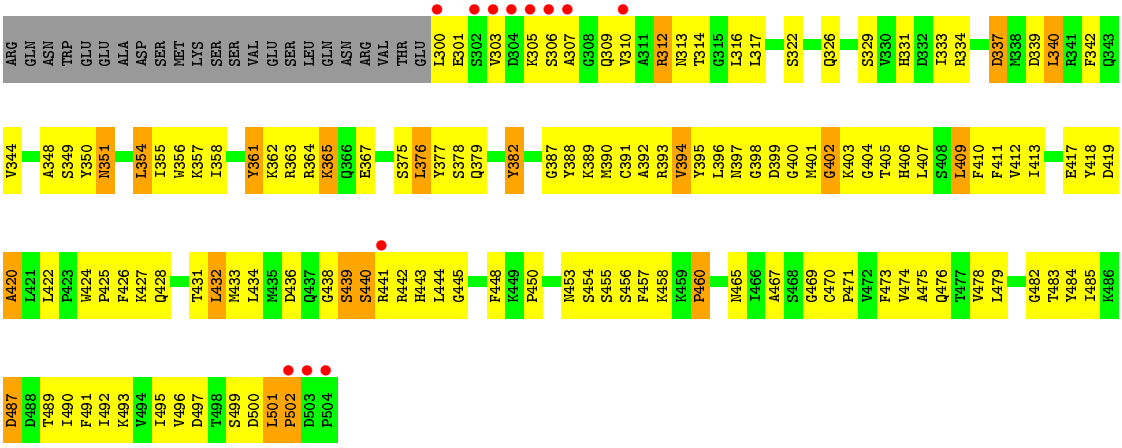
- Molecule 1: TNF RECEPTOR ASSOCIATED FACTOR 3

Chain A:



- Molecule 1: TNF RECEPTOR ASSOCIATED FACTOR 3

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α, β, γ	84.53Å 84.53Å 319.50Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	8.00 – 2.80 23.79 – 2.80	Depositor EDS
% Data completeness (in resolution range)	88.2 (8.00-2.80) 95.1 (23.79-2.80)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	4.77 (at 2.80Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.224 , 0.286 0.238 , 0.234	Depositor DCC
R_{free} test set	971 reflections (4.67%)	wwPDB-VP
Wilson B-factor (Å^2)	45.9	Xtriage
Anisotropy	0.636	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.42 , 43.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.487 for -h-k,k,-l	Xtriage
F_o,F_c correlation	0.90	EDS
Total number of atoms	3252	wwPDB-VP
Average B, all atoms (Å^2)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.84% 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	A	0.42	0/1661	0.67	1/2239 (0.0%)
1	B	0.41	0/1661	0.68	0/2239
All	All	0.41	0/3322	0.68	1/4478 (0.0%)

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	A	0	1
1	B	0	1
All	All	0	2

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	317	LEU	CA-CB-CG	5.50	127.94	115.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	418	TYR	Sidechain
1	B	350	TYR	Sidechain

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	1626	0	1622	156	0
1	B	1626	0	1622	153	0
All	All	3252	0	3244	309	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 48.

All (309) 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:436:ASP:OD2	1:B:441:ARG:HB3	1.54	1.08
1:A:441:ARG:HG2	1:A:442:ARG:HD3	1.39	1.03
1:B:314:THR:HA	1:B:317:LEU:HD23	1.40	1.01
1:A:410:PHE:HD2	1:A:469:GLY:H	1.08	0.96
1:A:434:LEU:HD21	1:A:479:LEU:HD12	1.45	0.96
1:B:390:MET:SD	1:B:496:VAL:HG21	2.09	0.93
1:A:455:SER:HA	1:A:458:LYS:HD2	1.50	0.92
1:A:413:ILE:O	1:A:465:ASN:HB2	1.70	0.92
1:B:441:ARG:NH1	1:B:483:THR:H	1.69	0.89
1:A:479:LEU:HD23	1:A:485:ILE:HD11	1.56	0.88
1:A:357:LYS:NZ	1:A:359:ARG:HG2	1.89	0.88
1:B:427:LYS:HA	1:B:450:PRO:HB2	1.56	0.87
1:A:305:LYS:O	1:A:309:GLN:HB2	1.76	0.86
1:A:441:ARG:HG2	1:A:442:ARG:H	1.42	0.84
1:B:441:ARG:HH12	1:B:483:THR:H	1.26	0.81
1:A:410:PHE:HD2	1:A:469:GLY:N	1.80	0.79
1:A:369:VAL:HG22	1:A:397:ASN:ND2	1.97	0.79
1:A:342:PHE:O	1:A:346:GLU:HG3	1.83	0.79
1:A:395:TYR:HB2	1:A:408:SER:HB2	1.65	0.78
1:A:436:ASP:OD2	1:A:441:ARG:HB3	1.84	0.77
1:B:312:ARG:HA	1:B:312:ARG:HE	1.49	0.77
1:B:441:ARG:NH2	1:B:483:THR:HG23	2.01	0.76
1:B:351:ASN:HD22	1:B:351:ASN:H	1.33	0.76
1:A:355:ILE:HD11	1:A:493:LYS:HE3	1.68	0.76
1:A:434:LEU:CD2	1:A:479:LEU:HD12	2.15	0.76
1:A:434:LEU:HD23	1:A:444:LEU:HD12	1.68	0.76
1:B:487:ASP:O	1:B:489:THR:HG23	1.85	0.75
1:A:442:ARG:H	1:A:442:ARG:HD3	1.50	0.75
1:B:300:LEU:HD23	1:B:303:VAL:HG21	1.68	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:410:PHE:HA	1:B:469:GLY:HA3	1.69	0.74
1:A:441:ARG:NH2	1:A:483:THR:HG23	2.02	0.74
1:A:357:LYS:HZ2	1:A:359:ARG:HG2	1.51	0.74
1:B:303:VAL:HG12	1:B:307:ALA:HB2	1.68	0.74
1:A:340:LEU:O	1:A:344:VAL:HG23	1.88	0.73
1:A:327:MET:HG3	1:A:331:HIS:CE1	2.23	0.73
1:A:355:ILE:HD13	1:A:493:LYS:HG3	1.72	0.72
1:A:472:VAL:HG23	1:A:472:VAL:O	1.90	0.72
1:A:344:VAL:CG1	1:A:380:PRO:HD2	2.18	0.72
1:A:433:MET:HG2	1:A:493:LYS:HB3	1.70	0.71
1:A:402:GLY:CA	1:A:472:VAL:HG12	2.20	0.71
1:B:439:SER:O	1:B:440:SER:HB2	1.90	0.70
1:B:426:PHE:O	1:B:457:PHE:HB3	1.91	0.70
1:B:363:ARG:O	1:B:367:GLU:HG3	1.91	0.70
1:A:438:GLY:HA3	1:A:483:THR:HG21	1.74	0.69
1:B:312:ARG:O	1:B:316:LEU:HB2	1.93	0.69
1:A:303:VAL:HA	1:A:306:SER:HB2	1.75	0.69
1:A:390:MET:O	1:A:391:CYS:HB3	1.93	0.69
1:B:441:ARG:HH12	1:B:483:THR:N	1.90	0.68
1:B:303:VAL:HA	1:B:306:SER:HB2	1.75	0.68
1:A:314:THR:HA	1:A:317:LEU:HD23	1.76	0.67
1:A:413:ILE:HD12	1:A:457:PHE:HA	1.74	0.67
1:B:493:LYS:HE2	1:B:495:ILE:HD11	1.77	0.67
1:A:303:VAL:HG12	1:A:307:ALA:HB3	1.77	0.66
1:B:441:ARG:HH22	1:B:483:THR:HG23	1.59	0.66
1:B:475:ALA:HB3	1:B:478:VAL:HG23	1.78	0.65
1:A:483:THR:O	1:A:485:ILE:N	2.30	0.65
1:B:473:PHE:CD2	1:B:474:VAL:HG23	2.31	0.65
1:A:410:PHE:HA	1:A:469:GLY:HA3	1.77	0.65
1:A:419:ASP:O	1:A:420:ALA:CB	2.44	0.65
1:A:480:GLU:HA	1:A:485:ILE:CD1	2.27	0.64
1:A:375:SER:HB3	1:A:395:TYR:CE2	2.33	0.64
1:B:361:TYR:CD2	1:B:485:ILE:HD11	2.32	0.64
1:B:455:SER:HA	1:B:458:LYS:HG3	1.80	0.64
1:A:357:LYS:HZ1	1:A:359:ARG:HG2	1.61	0.63
1:A:376:LEU:HB2	1:A:394:VAL:HG12	1.79	0.63
1:A:357:LYS:HG3	1:A:491:PHE:CZ	2.32	0.63
1:B:394:VAL:HG23	1:B:409:LEU:HD23	1.79	0.63
1:B:354:LEU:HD22	1:B:355:ILE:N	2.13	0.63
1:A:487:ASP:O	1:A:489:THR:HG23	1.99	0.62
1:A:451:ASP:OD2	1:A:453:ASN:HB2	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:364:ARG:HA	1:B:367:GLU:OE2	2.00	0.62
1:A:361:TYR:O	1:A:365:LYS:HB2	2.00	0.62
1:B:444:LEU:HD13	1:B:474:VAL:HG22	1.82	0.61
1:B:339:ASP:HA	1:B:342:PHE:HD1	1.64	0.61
1:B:448:PHE:O	1:B:448:PHE:CD1	2.54	0.61
1:B:493:LYS:CE	1:B:495:ILE:HD11	2.30	0.61
1:B:410:PHE:HD2	1:B:469:GLY:CA	2.14	0.61
1:B:460:PRO:HA	1:B:465:ASN:HD21	1.66	0.61
1:B:333:ILE:O	1:B:337:ASP:OD1	2.19	0.60
1:A:419:ASP:O	1:A:420:ALA:HB2	2.01	0.60
1:B:406:HIS:HA	1:B:476:GLN:HG3	1.82	0.60
1:A:415:ARG:HA	1:A:460:PRO:HB2	1.82	0.60
1:A:300:LEU:HD23	1:A:304:ASP:OD2	2.00	0.60
1:A:301:GLU:HG3	1:A:302:SER:H	1.67	0.60
1:B:358:ILE:HB	1:B:490:ILE:CG1	2.32	0.59
1:B:356:TRP:CD2	1:B:378:SER:HB3	2.38	0.59
1:A:376:LEU:HB2	1:A:394:VAL:CG1	2.32	0.59
1:A:303:VAL:HA	1:A:306:SER:CB	2.32	0.59
1:A:409:LEU:CD1	1:A:432:LEU:HB2	2.32	0.59
1:B:438:GLY:O	1:B:440:SER:N	2.36	0.59
1:A:357:LYS:HZ2	1:A:359:ARG:CG	2.16	0.59
1:B:426:PHE:N	1:B:457:PHE:O	2.36	0.59
1:B:413:ILE:HD11	1:B:426:PHE:CG	2.38	0.59
1:A:330:VAL:O	1:A:334:ARG:HG2	2.03	0.58
1:B:500:ASP:O	1:B:501:LEU:HB2	2.03	0.58
1:B:432:LEU:HD23	1:B:432:LEU:H	1.68	0.58
1:A:301:GLU:CG	1:A:302:SER:H	2.15	0.57
1:A:439:SER:O	1:A:440:SER:HB2	2.04	0.57
1:B:441:ARG:HH12	1:B:482:GLY:CA	2.16	0.57
1:A:354:LEU:HD22	1:A:355:ILE:N	2.20	0.57
1:A:412:VAL:HG11	1:A:464:MET:HB3	1.85	0.57
1:B:365:LYS:HD3	1:B:365:LYS:O	2.05	0.57
1:B:406:HIS:N	1:B:476:GLN:HE21	2.03	0.57
1:B:356:TRP:CZ3	1:B:394:VAL:HB	2.39	0.57
1:B:442:ARG:HH11	1:B:442:ARG:HG2	1.70	0.57
1:A:389:LYS:C	1:A:390:MET:HG3	2.24	0.57
1:B:361:TYR:CE2	1:B:485:ILE:HD11	2.40	0.57
1:A:415:ARG:HA	1:A:460:PRO:CB	2.35	0.56
1:A:436:ASP:HB2	1:A:484:TYR:HE1	1.69	0.56
1:A:480:GLU:HA	1:A:485:ILE:HD12	1.87	0.56
1:B:354:LEU:HD21	1:B:379:GLN:NE2	2.20	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:340:LEU:O	1:B:344:VAL:HG23	2.05	0.56
1:B:400:GLY:O	1:B:402:GLY:N	2.34	0.56
1:B:382:TYR:CE2	1:B:389:LYS:HG3	2.40	0.56
1:B:395:TYR:O	1:B:407:LEU:HA	2.05	0.56
1:A:344:VAL:HG11	1:A:380:PRO:HD2	1.88	0.56
1:A:326:GLN:O	1:A:329:SER:OG	2.24	0.56
1:B:493:LYS:NZ	1:B:495:ILE:HD11	2.20	0.55
1:B:351:ASN:HD22	1:B:351:ASN:N	1.96	0.55
1:A:398:GLY:HA2	1:A:408:SER:OG	2.07	0.55
1:B:358:ILE:HB	1:B:490:ILE:HG13	1.88	0.55
1:A:394:VAL:HG22	1:A:395:TYR:N	2.22	0.55
1:B:425:PRO:O	1:B:427:LYS:HE3	2.06	0.55
1:A:402:GLY:HA3	1:A:472:VAL:HG12	1.89	0.55
1:A:395:TYR:HB2	1:A:408:SER:CB	2.35	0.54
1:A:326:GLN:O	1:A:330:VAL:HG23	2.07	0.54
1:A:353:VAL:HG12	1:A:354:LEU:N	2.23	0.54
1:B:454:SER:OG	1:B:456:SER:HB2	2.08	0.54
1:A:431:THR:HB	1:A:495:ILE:HB	1.89	0.54
1:A:361:TYR:O	1:A:365:LYS:N	2.39	0.54
1:B:354:LEU:HD11	1:B:378:SER:OG	2.08	0.54
1:B:349:SER:OG	1:B:351:ASN:ND2	2.40	0.54
1:B:409:LEU:O	1:B:469:GLY:HA3	2.08	0.54
1:B:303:VAL:HA	1:B:306:SER:CB	2.37	0.54
1:B:307:ALA:O	1:B:310:VAL:HG13	2.08	0.54
1:B:356:TRP:HZ3	1:B:394:VAL:HB	1.72	0.54
1:A:351:ASN:HD22	1:A:351:ASN:H	1.55	0.53
1:A:432:LEU:HD23	1:A:432:LEU:O	2.08	0.53
1:B:400:GLY:C	1:B:402:GLY:H	2.12	0.53
1:B:410:PHE:CE2	1:B:471:PRO:HG3	2.43	0.53
1:A:300:LEU:HG	1:A:304:ASP:HA	1.90	0.53
1:A:313:ASN:HA	1:A:316:LEU:HD23	1.91	0.53
1:B:444:LEU:CD1	1:B:474:VAL:HG22	2.39	0.53
1:A:354:LEU:HD21	1:A:379:GLN:NE2	2.24	0.53
1:A:436:ASP:HB2	1:A:484:TYR:CE1	2.44	0.53
1:B:432:LEU:N	1:B:432:LEU:HD23	2.23	0.53
1:B:357:LYS:HD2	1:B:491:PHE:CZ	2.43	0.53
1:A:409:LEU:HD12	1:A:432:LEU:HB2	1.90	0.53
1:B:411:PHE:CD1	1:B:412:VAL:N	2.77	0.53
1:A:479:LEU:CD2	1:A:485:ILE:HD11	2.35	0.53
1:B:406:HIS:CA	1:B:476:GLN:HG3	2.37	0.53
1:A:455:SER:HA	1:A:458:LYS:CD	2.34	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:309:GLN:O	1:A:309:GLN:HG2	2.09	0.52
1:A:444:LEU:HD13	1:A:474:VAL:HG21	1.91	0.52
1:A:333:ILE:O	1:A:337:ASP:OD1	2.28	0.52
1:B:410:PHE:HD2	1:B:469:GLY:HA3	1.74	0.51
1:A:476:GLN:O	1:A:480:GLU:HG3	2.10	0.51
1:A:364:ARG:HG2	1:A:376:LEU:HD11	1.93	0.51
1:A:438:GLY:O	1:A:440:SER:N	2.44	0.51
1:A:343:GLN:O	1:A:347:THR:HG23	2.10	0.51
1:B:377:TYR:HA	1:B:392:ALA:O	2.10	0.51
1:A:331:HIS:HA	1:A:334:ARG:HG3	1.93	0.51
1:B:305:LYS:HB2	1:B:305:LYS:NZ	2.26	0.51
1:B:303:VAL:HG12	1:B:307:ALA:CB	2.40	0.50
1:A:436:ASP:CG	1:A:441:ARG:HB3	2.30	0.50
1:B:398:GLY:C	1:B:403:LYS:HA	2.32	0.50
1:B:410:PHE:CD2	1:B:469:GLY:HA3	2.47	0.50
1:B:448:PHE:O	1:B:448:PHE:HD1	1.94	0.50
1:A:441:ARG:CZ	1:A:483:THR:HG23	2.41	0.50
1:A:436:ASP:OD1	1:A:483:THR:OG1	2.23	0.50
1:A:430:VAL:HG12	1:A:431:THR:N	2.26	0.50
1:A:357:LYS:HG3	1:A:491:PHE:CE2	2.47	0.50
1:B:410:PHE:CZ	1:B:471:PRO:HG3	2.47	0.50
1:B:322:SER:O	1:B:326:GLN:HG3	2.12	0.49
1:A:414:MET:O	1:A:460:PRO:HG2	2.13	0.49
1:B:441:ARG:NH1	1:B:483:THR:N	2.49	0.49
1:A:433:MET:CG	1:A:493:LYS:HB3	2.41	0.49
1:A:300:LEU:O	1:A:304:ASP:HB2	2.13	0.49
1:A:359:ARG:HA	1:A:489:THR:HG22	1.95	0.49
1:A:432:LEU:CD2	1:A:446:ASP:O	2.60	0.49
1:A:480:GLU:HA	1:A:485:ILE:HD11	1.95	0.49
1:B:331:HIS:O	1:B:334:ARG:HB2	2.13	0.49
1:B:361:TYR:C	1:B:361:TYR:CD1	2.86	0.48
1:B:434:LEU:HB2	1:B:473:PHE:HE2	1.77	0.48
1:B:390:MET:CE	1:B:496:VAL:HG21	2.42	0.48
1:B:329:SER:O	1:B:333:ILE:HG13	2.14	0.48
1:B:357:LYS:HB2	1:B:491:PHE:CE2	2.48	0.48
1:A:436:ASP:HB3	1:A:443:HIS:ND1	2.28	0.48
1:B:354:LEU:CD2	1:B:379:GLN:NE2	2.76	0.48
1:A:382:TYR:N	1:A:382:TYR:CD1	2.80	0.48
1:A:402:GLY:O	1:A:406:HIS:HB2	2.13	0.48
1:A:409:LEU:HD21	1:A:492:ILE:HG21	1.95	0.48
1:B:351:ASN:N	1:B:351:ASN:ND2	2.62	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:478:VAL:O	1:B:478:VAL:HG12	2.14	0.48
1:A:441:ARG:HG2	1:A:442:ARG:N	2.21	0.48
1:B:387:GLY:HA2	1:B:422:LEU:HD11	1.96	0.48
1:A:430:VAL:CG1	1:A:431:THR:N	2.77	0.47
1:B:303:VAL:HG12	1:B:303:VAL:O	2.14	0.47
1:B:465:ASN:HD22	1:B:465:ASN:N	2.11	0.47
1:B:417:GLU:HB3	1:B:418:TYR:CE1	2.49	0.47
1:A:411:PHE:CE2	1:A:430:VAL:HG21	2.50	0.47
1:A:472:VAL:CG2	1:A:472:VAL:O	2.62	0.47
1:A:442:ARG:N	1:A:442:ARG:HD3	2.26	0.47
1:A:434:LEU:HB3	1:A:444:LEU:HB2	1.97	0.47
1:A:457:PHE:CZ	1:A:468:SER:OG	2.65	0.47
1:A:409:LEU:CD2	1:A:492:ILE:HD13	2.44	0.47
1:B:356:TRP:HH2	1:B:376:LEU:O	1.97	0.47
1:A:479:LEU:O	1:A:485:ILE:HG13	2.14	0.47
1:B:303:VAL:HG13	1:B:306:SER:HB2	1.96	0.47
1:A:303:VAL:O	1:A:307:ALA:N	2.48	0.47
1:B:375:SER:OG	1:B:393:ARG:NE	2.48	0.47
1:B:413:ILE:HD11	1:B:426:PHE:CB	2.44	0.47
1:A:375:SER:HB3	1:A:395:TYR:CD2	2.50	0.47
1:A:450:PRO:HB3	1:A:457:PHE:CD1	2.50	0.47
1:B:407:LEU:HD23	1:B:407:LEU:C	2.35	0.47
1:A:465:ASN:HA	1:A:465:ASN:HD22	1.52	0.46
1:B:382:TYR:CD1	1:B:382:TYR:N	2.83	0.46
1:B:433:MET:HB3	1:B:445:GLY:HA3	1.97	0.46
1:A:388:TYR:CD1	1:A:422:LEU:HD23	2.50	0.46
1:A:355:ILE:CD1	1:A:493:LYS:HG3	2.44	0.46
1:B:397:ASN:OD1	1:B:404:GLY:HA2	2.16	0.46
1:B:425:PRO:HD2	1:B:427:LYS:NZ	2.30	0.46
1:A:444:LEU:HD13	1:A:474:VAL:CG2	2.46	0.46
1:B:431:THR:O	1:B:431:THR:HG22	2.16	0.46
1:B:454:SER:HB3	1:B:457:PHE:CD2	2.51	0.46
1:A:483:THR:OG1	1:A:484:TYR:N	2.49	0.45
1:A:501:LEU:O	1:A:502:PRO:C	2.53	0.45
1:B:313:ASN:O	1:B:317:LEU:HD22	2.15	0.45
1:A:495:ILE:HG22	1:A:495:ILE:O	2.17	0.45
1:B:357:LYS:HD2	1:B:491:PHE:CE2	2.52	0.45
1:A:369:VAL:HA	1:A:397:ASN:HD22	1.81	0.45
1:A:501:LEU:HD22	1:A:501:LEU:HA	1.78	0.45
1:B:351:ASN:ND2	1:B:351:ASN:H	2.07	0.45
1:B:428:GLN:HB3	1:B:497:ASP:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:311:ALA:C	1:A:314:THR:HG22	2.37	0.44
1:B:395:TYR:CD1	1:B:399:ASP:HB2	2.51	0.44
1:B:410:PHE:HD2	1:B:469:GLY:N	2.16	0.44
1:B:441:ARG:CZ	1:B:483:THR:H	2.29	0.44
1:A:394:VAL:HG23	1:A:408:SER:O	2.18	0.44
1:A:356:TRP:CD1	1:A:357:LYS:N	2.85	0.44
1:B:419:ASP:O	1:B:420:ALA:HB2	2.18	0.44
1:B:358:ILE:HB	1:B:490:ILE:HD11	1.99	0.44
1:A:456:SER:O	1:A:465:ASN:OD1	2.36	0.44
1:B:453:ASN:O	1:B:454:SER:C	2.55	0.44
1:B:455:SER:HA	1:B:458:LYS:CG	2.46	0.44
1:A:477:THR:O	1:A:481:ASN:HB2	2.17	0.43
1:B:405:THR:C	1:B:476:GLN:HG3	2.39	0.43
1:B:361:TYR:HD1	1:B:361:TYR:C	2.21	0.43
1:B:436:ASP:HB2	1:B:484:TYR:CE1	2.53	0.43
1:A:303:VAL:CA	1:A:306:SER:HB2	2.45	0.43
1:A:426:PHE:CE1	1:A:428:GLN:HB2	2.54	0.43
1:B:313:ASN:O	1:B:317:LEU:CD2	2.66	0.43
1:A:309:GLN:HG2	1:A:312:ARG:HB2	2.01	0.43
1:A:313:ASN:O	1:A:317:LEU:HD22	2.18	0.43
1:B:394:VAL:HG23	1:B:409:LEU:CD2	2.48	0.43
1:A:301:GLU:HG3	1:A:302:SER:N	2.33	0.43
1:B:312:ARG:NE	1:B:312:ARG:HA	2.25	0.43
1:B:382:TYR:HE2	1:B:389:LYS:HG3	1.83	0.43
1:B:454:SER:HB3	1:B:457:PHE:HD2	1.84	0.43
1:A:303:VAL:HA	1:A:306:SER:OG	2.19	0.43
1:B:405:THR:OG1	1:B:406:HIS:HD2	2.01	0.43
1:A:303:VAL:HG12	1:A:303:VAL:O	2.20	0.42
1:B:391:CYS:SG	1:B:412:VAL:HB	2.59	0.42
1:B:484:TYR:O	1:B:490:ILE:HG22	2.19	0.42
1:B:305:LYS:O	1:B:309:GLN:HB2	2.19	0.42
1:A:351:ASN:N	1:A:351:ASN:HD22	2.17	0.42
1:A:409:LEU:HD23	1:A:492:ILE:HD13	2.01	0.42
1:B:396:LEU:HD23	1:B:407:LEU:HD12	2.01	0.42
1:A:351:ASN:H	1:A:351:ASN:ND2	2.16	0.42
1:A:395:TYR:CD1	1:A:408:SER:HB2	2.55	0.42
1:B:303:VAL:CG1	1:B:307:ALA:HB2	2.44	0.42
1:B:300:LEU:HB3	1:B:303:VAL:HG23	2.02	0.42
1:B:400:GLY:C	1:B:402:GLY:N	2.73	0.42
1:B:455:SER:HA	1:B:458:LYS:CD	2.50	0.42
1:A:308:GLY:C	1:A:310:VAL:H	2.23	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:441:ARG:HG2	1:A:442:ARG:CD	2.28	0.41
1:B:306:SER:O	1:B:309:GLN:HB3	2.20	0.41
1:B:395:TYR:CD1	1:B:410:PHE:HE1	2.38	0.41
1:A:356:TRP:CZ3	1:A:358:ILE:HD11	2.55	0.41
1:B:460:PRO:CA	1:B:465:ASN:HD21	2.32	0.41
1:A:499:SER:C	1:A:501:LEU:H	2.24	0.41
1:B:300:LEU:HD12	1:B:300:LEU:N	2.35	0.41
1:B:436:ASP:H	1:B:443:HIS:CE1	2.37	0.41
1:A:354:LEU:CD2	1:A:379:GLN:NE2	2.84	0.41
1:A:483:THR:O	1:A:484:TYR:C	2.59	0.41
1:B:441:ARG:NH2	1:B:483:THR:CG2	2.78	0.41
1:A:361:TYR:C	1:A:361:TYR:CD1	2.94	0.41
1:B:493:LYS:HE2	1:B:495:ILE:CD1	2.48	0.41
1:B:361:TYR:O	1:B:362:LYS:C	2.59	0.41
1:B:410:PHE:HB3	1:B:467:ALA:HB1	2.02	0.41
1:B:356:TRP:CG	1:B:378:SER:HB3	2.55	0.41
1:B:388:TYR:N	1:B:422:LEU:HD11	2.35	0.41
1:A:354:LEU:HB2	1:A:381:PHE:CD2	2.56	0.41
1:A:418:TYR:O	1:A:422:LEU:HD13	2.20	0.41
1:B:441:ARG:CZ	1:B:483:THR:OG1	2.69	0.41
1:A:301:GLU:C	1:A:303:VAL:H	2.24	0.41
1:A:358:ILE:O	1:A:489:THR:HB	2.20	0.41
1:B:450:PRO:HB3	1:B:457:PHE:CD2	2.56	0.41
1:B:450:PRO:HB3	1:B:457:PHE:CE2	2.55	0.41
1:A:368:ALA:HB3	1:A:397:ASN:HB2	2.03	0.41
1:A:399:ASP:HA	1:A:403:LYS:CE	2.51	0.41
1:A:424:TRP:HB3	1:A:425:PRO:HA	2.02	0.40
1:B:425:PRO:CA	1:B:457:PHE:O	2.69	0.40
1:B:361:TYR:OH	1:B:479:LEU:HD23	2.20	0.40
1:B:305:LYS:HB2	1:B:305:LYS:HZ2	1.86	0.40
1:B:358:ILE:HD12	1:B:492:ILE:HD12	2.03	0.40
1:A:316:LEU:HD22	1:A:316:LEU:N	2.36	0.40
1:A:436:ASP:HA	1:A:484:TYR:CD1	2.55	0.40
1:B:419:ASP:OD1	1:B:424:TRP:HZ2	2.05	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	203/228 (89%)	162 (80%)	28 (14%)	13 (6%)	1 3
1	B	203/228 (89%)	165 (81%)	28 (14%)	10 (5%)	2 7
All	All	406/456 (89%)	327 (80%)	56 (14%)	23 (6%)	1 5

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	301	GLU
1	A	391	CYS
1	A	439	SER
1	A	440	SER
1	A	484	TYR
1	B	420	ALA
1	B	439	SER
1	B	440	SER
1	B	501	LEU
1	A	306	SER
1	A	420	ALA
1	A	463	GLU
1	A	502	PRO
1	B	499	SER
1	A	465	ASN
1	B	401	MET
1	B	502	PRO
1	A	398	GLY
1	A	446	ASP
1	B	402	GLY
1	A	460	PRO
1	B	348	ALA
1	B	460	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	A	179/201 (89%)	164 (92%)	15 (8%)	11 31
1	B	179/201 (89%)	163 (91%)	16 (9%)	9 28
All	All	358/402 (89%)	327 (91%)	31 (9%)	10 30

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	301	GLU
1	A	312	ARG
1	A	327	MET
1	A	329	SER
1	A	334	ARG
1	A	340	LEU
1	A	342	PHE
1	A	351	ASN
1	A	391	CYS
1	A	409	LEU
1	A	432	LEU
1	A	442	ARG
1	A	458	LYS
1	A	465	ASN
1	A	502	PRO
1	B	301	GLU
1	B	312	ARG
1	B	337	ASP
1	B	340	LEU
1	B	351	ASN
1	B	354	LEU
1	B	361	TYR
1	B	365	LYS
1	B	376	LEU
1	B	382	TYR
1	B	394	VAL
1	B	409	LEU

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Mol	Chain	Res	Type
1	B	432	LEU
1	B	470	CYS
1	B	487	ASP
1	B	502	PRO

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	343	GLN
1	A	351	ASN
1	A	428	GLN
1	A	465	ASN
1	B	309	GLN
1	B	343	GLN
1	B	351	ASN
1	B	406	HIS
1	B	428	GLN
1	B	465	ASN
1	B	476	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

There are no carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	A	205/228 (89%)	0.62	12 (5%)	22	14	26, 44, 89, 103	0
1	B	205/228 (89%)	0.62	12 (5%)	22	14	26, 44, 89, 103	0
All	All	410/456 (89%)	0.62	24 (5%)	22	14	26, 44, 90, 103	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	303	VAL	6.9
1	B	503	ASP	5.9
1	A	303	VAL	5.2
1	A	305	LYS	4.9
1	A	304	ASP	4.6
1	A	503	ASP	4.6
1	A	306	SER	4.4
1	B	306	SER	4.1
1	B	302	SER	4.1
1	B	300	LEU	4.0
1	A	300	LEU	4.0
1	B	307	ALA	3.6
1	B	310	VAL	3.5
1	B	504	PRO	3.3
1	A	307	ALA	3.1
1	B	304	ASP	2.9
1	A	504	PRO	2.7
1	A	302	SER	2.6
1	A	502	PRO	2.6
1	A	311	ALA	2.6
1	A	362	LYS	2.5
1	B	502	PRO	2.3
1	B	441	ARG	2.2
1	B	305	LYS	2.1

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

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.