



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:08 PM GMT

PDB ID : 1S9H
Title : Crystal Structure of Adeno-associated virus Type 2 Rep40
Authors : James, J.A.; Escalante, C.R.; Yoon-Robarts, M.; Edwards, T.A.; Linden, R.M.; Aggarwal, A.K.
Deposited on : 2004-02-04
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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

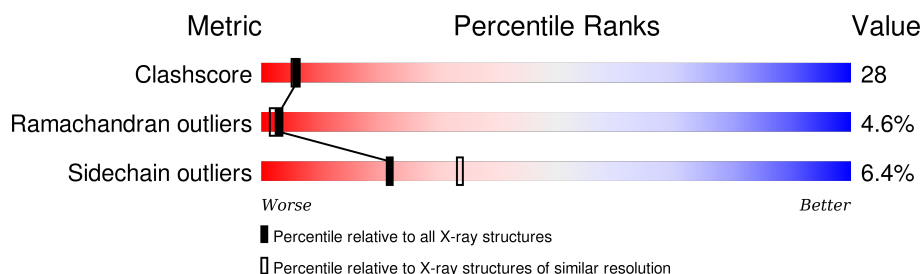
1 Overall quality at a glance

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))
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	268	 59% 34% 7%
1	B	268	 60% 34% 5%
1	C	268	 38% 30% 6% 26%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5963 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 Rep 40 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	268	Total	C	N	O	S	0	0	0
			2080	1329	350	391	10			
1	B	267	Total	C	N	O	S	0	0	0
			2064	1319	346	389	10			
1	C	199	Total	C	N	O	S	0	0	0
			1462	947	240	270	5			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	223	ALA	-	CLONING ARTIFACT	GB 2906019
A	224	GLY	-	CLONING ARTIFACT	GB 2906019
B	223	ALA	-	CLONING ARTIFACT	GB 2906019
B	224	GLY	-	CLONING ARTIFACT	GB 2906019
C	223	ALA	-	CLONING ARTIFACT	GB 2906019
C	224	GLY	-	CLONING ARTIFACT	GB 2906019

- Molecule 2 is water.

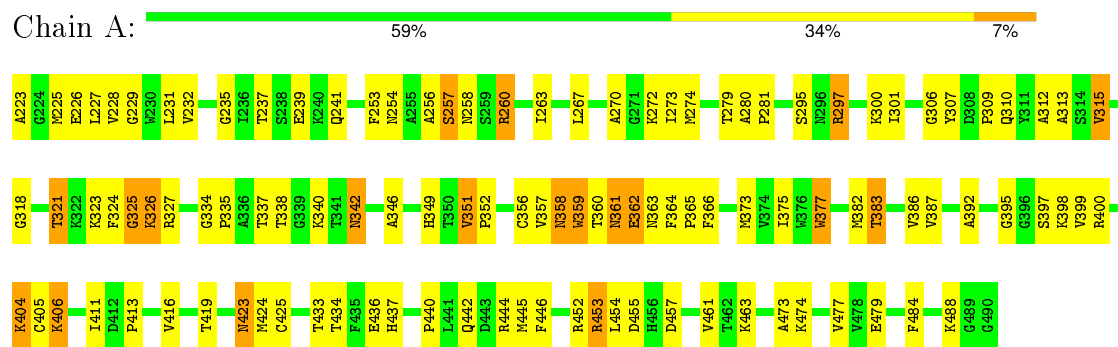
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	166	Total	O	0	0
			166	166		
2	B	117	Total	O	0	0
			117	117		
2	C	74	Total	O	0	0
			74	74		

3 Residue-property plots

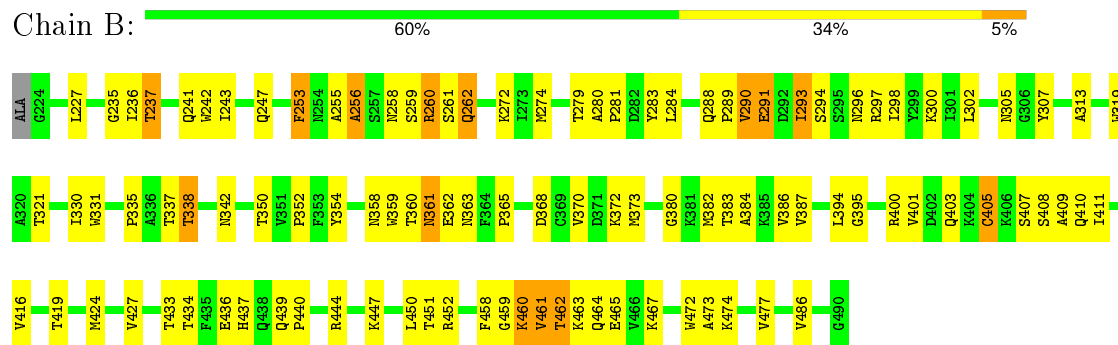
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according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

Note EDS was not executed.

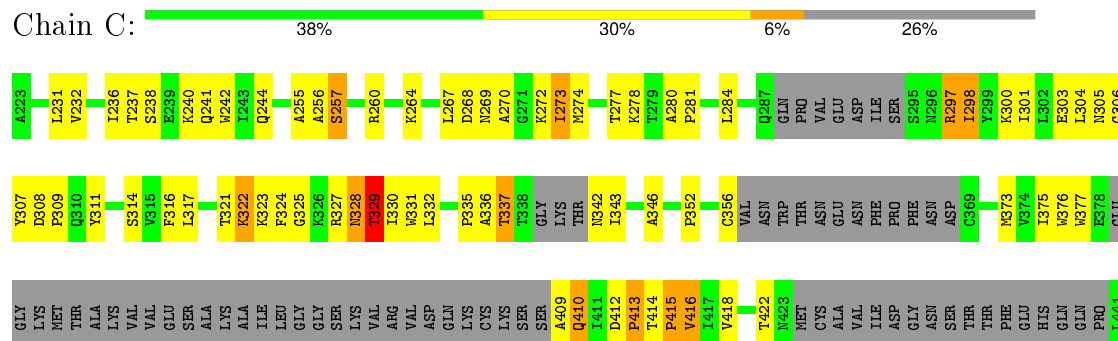
- Molecule 1: Rep 40 protein



- Molecule 1: Rep 40 protein



- Molecule 1: Rep 40 protein



Q442	D443	R444	R445
L450	Y461	Y462	K463
Q464	R465	Y466	K467
Y472	D475	H476	Y477
Y478	E483	F484	Y485
Y486	K487	K488	G489
G490			

4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	126.29 Å 126.29 Å 97.89 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.97 – 2.40	Depositor
% Data completeness (in resolution range)	85.5 (19.97-2.40)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.227 , 0.293	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5963	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

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	0.43	0/2130	0.68	0/2890
1	B	0.39	0/2113	0.61	0/2868
1	C	0.34	0/1494	0.60	1/2033 (0.0%)
All	All	0.40	0/5737	0.63	1/7791 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	335	PRO	N-CA-CB	5.45	109.83	103.30

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	2080	0	2003	112	0
1	B	2064	0	1969	102	0
1	C	1462	0	1314	99	0
2	A	166	0	0	12	0
2	B	117	0	0	2	0
2	C	74	0	0	3	0
All	All	5963	0	5286	310	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 28.

All (310) 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:A:434:THR:HG22	1:A:436:GLU:H	1.20	1.05
1:B:293:ILE:HD12	1:B:293:ILE:H	1.21	1.01
1:C:238:SER:OG	1:C:241:GLN:HG2	1.65	0.95
1:A:453:ARG:HE	1:A:453:ARG:H	1.17	0.88
1:C:307:TYR:O	1:C:309:PRO:HD3	1.72	0.88
1:B:342:ASN:HD21	1:B:460:LYS:H	1.23	0.86
1:A:357:VAL:HG12	1:A:377:TRP:HA	1.60	0.83
1:A:357:VAL:CG1	1:A:377:TRP:HA	2.08	0.82
1:C:322:LYS:NZ	1:C:322:LYS:HA	1.94	0.81
1:C:297:ARG:HG2	1:C:461:VAL:HA	1.62	0.81
1:B:459:GLY:O	1:B:460:LYS:HB2	1.83	0.79
1:A:359:TRP:CZ2	1:A:382:MET:HA	2.17	0.78
1:A:392:ALA:HB3	1:A:399:VAL:HG21	1.66	0.78
1:A:392:ALA:CB	1:A:399:VAL:HG21	2.14	0.78
1:A:337:THR:HG22	1:A:337:THR:O	1.82	0.78
1:B:305:ASN:HD21	1:B:451:THR:H	1.32	0.76
1:B:305:ASN:ND2	1:B:451:THR:H	1.84	0.76
1:B:280:ALA:HA	1:B:373:MET:HE3	1.68	0.76
1:B:462:THR:HG22	1:B:465:GLU:HG3	1.68	0.76
1:B:462:THR:CG2	1:B:465:GLU:H	1.99	0.75
1:C:463:LYS:N	1:C:463:LYS:HD3	2.02	0.75
1:A:231:LEU:N	2:A:496:HOH:O	2.20	0.74
1:A:334:GLY:O	1:A:340:LYS:HE2	1.88	0.74
1:A:228:VAL:O	2:A:496:HOH:O	2.06	0.74
1:C:269:ASN:O	1:C:273:ILE:HG23	1.88	0.74
1:C:273:ILE:O	1:C:277:THR:HB	1.90	0.72
1:A:279:THR:HG22	1:A:280:ALA:N	2.04	0.72
1:B:305:ASN:ND2	1:B:450:LEU:HA	2.04	0.72
1:C:321:THR:HG22	1:C:322:LYS:H	1.54	0.71
1:C:375:ILE:HB	1:C:416:VAL:HG22	1.71	0.71
1:C:305:ASN:HB3	1:C:488:LYS:NZ	2.05	0.71
1:A:453:ARG:H	1:A:453:ARG:NE	1.89	0.70
1:C:342:ASN:OD1	1:C:461:VAL:HG23	1.90	0.70
1:B:352:PRO:HD2	1:B:373:MET:HE2	1.73	0.69
1:B:294:SER:O	1:B:300:LYS:HE3	1.91	0.69
1:A:404:LYS:NZ	1:B:405:CYS:HB3	2.08	0.69
1:A:327:ARG:HD2	1:A:445:MET:O	1.93	0.68
1:C:463:LYS:H	1:C:463:LYS:HD3	1.58	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:392:ALA:HB1	1:A:399:VAL:CG2	2.23	0.68
1:B:338:THR:HG21	1:B:450:LEU:C	2.13	0.68
1:B:358:ASN:HB3	1:B:360:THR:HG23	1.76	0.68
1:B:279:THR:HG22	1:B:281:PRO:CD	2.24	0.68
1:C:280:ALA:HB3	1:C:281:PRO:HD3	1.76	0.67
1:B:403:GLN:HB2	1:B:407:SER:CB	2.24	0.67
1:A:357:VAL:HG11	1:A:377:TRP:CG	2.29	0.67
1:B:462:THR:HG23	1:B:465:GLU:H	1.59	0.67
1:B:359:TRP:CD1	1:B:380:GLY:HA3	2.29	0.67
1:A:325:GLY:O	1:A:326:LYS:HB2	1.93	0.67
1:B:361:ASN:C	1:B:361:ASN:HD22	1.97	0.66
1:A:232:VAL:HG22	2:A:491:HOH:O	1.95	0.66
1:A:312:ALA:O	1:A:315:VAL:HG13	1.96	0.66
1:B:290:VAL:O	1:B:291:GLU:HB3	1.94	0.66
1:B:261:SER:O	1:B:262:GLN:HB2	1.95	0.66
1:C:414:THR:O	1:C:416:VAL:HG23	1.97	0.65
1:B:321:THR:O	1:B:321:THR:HG22	1.96	0.65
1:C:322:LYS:HZ2	1:C:322:LYS:HA	1.61	0.64
1:B:293:ILE:CD1	1:B:293:ILE:H	1.96	0.64
1:B:382:MET:CE	1:B:387:VAL:HA	2.28	0.64
1:A:335:PRO:O	1:A:338:THR:HG22	1.96	0.64
1:A:382:MET:CE	1:A:387:VAL:HG23	2.26	0.64
1:C:308:ASP:HB2	1:C:490:GLY:HA2	1.80	0.64
1:A:352:PRO:HD2	1:A:373:MET:SD	2.37	0.64
1:B:384:ALA:HA	1:B:387:VAL:CG2	2.28	0.64
1:A:375:ILE:HB	1:A:416:VAL:HG22	1.79	0.64
1:B:462:THR:HG22	1:B:465:GLU:CG	2.27	0.63
1:C:332:LEU:HD22	1:C:450:LEU:HD11	1.81	0.63
1:A:297:ARG:HH11	1:A:297:ARG:CG	2.11	0.63
1:B:279:THR:HG22	1:B:281:PRO:HD2	1.82	0.62
1:B:439:GLN:HB3	1:B:440:PRO:HD3	1.82	0.62
1:A:357:VAL:HG11	1:A:377:TRP:CD1	2.34	0.62
1:A:404:LYS:HZ3	1:B:405:CYS:HB3	1.64	0.61
1:C:442:GLN:OE1	1:C:442:GLN:HA	2.00	0.61
1:B:342:ASN:ND2	1:B:460:LYS:H	1.92	0.61
1:B:236:ILE:HD13	1:B:242:TRP:HA	1.83	0.61
1:B:305:ASN:HD21	1:B:451:THR:N	1.99	0.60
1:A:254:ASN:HD21	1:A:263:ILE:HD11	1.66	0.60
1:B:474:LYS:O	1:B:477:VAL:HG23	2.01	0.60
1:B:437:HIS:O	1:B:440:PRO:HD2	2.02	0.60
1:C:332:LEU:HB3	1:C:450:LEU:HD13	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:279:THR:HG22	1:A:281:PRO:CD	2.32	0.60
1:B:447:LYS:O	1:B:486:VAL:HG23	2.02	0.60
1:C:232:VAL:HG21	1:C:273:ILE:CD1	2.32	0.60
1:A:392:ALA:CB	1:A:399:VAL:CG2	2.77	0.60
1:C:356:CYS:CB	1:C:376:TRP:HB3	2.31	0.60
1:A:337:THR:O	1:A:337:THR:CG2	2.50	0.60
1:B:313:ALA:HB1	1:B:473:ALA:HB2	1.84	0.59
1:B:280:ALA:HA	1:B:373:MET:CE	2.32	0.59
1:B:279:THR:HG22	1:B:280:ALA:N	2.18	0.59
1:A:229:GLY:C	2:A:496:HOH:O	2.40	0.59
1:C:321:THR:HG22	1:C:322:LYS:N	2.18	0.59
1:B:293:ILE:HG12	1:B:472:TRP:CD1	2.38	0.58
1:B:434:THR:HG22	1:B:436:GLU:H	1.68	0.58
1:C:322:LYS:HZ3	1:C:322:LYS:HA	1.66	0.58
1:A:307:TYR:O	1:A:309:PRO:HD3	2.03	0.58
1:A:411:ILE:HB	2:A:497:HOH:O	2.03	0.58
1:A:357:VAL:O	1:A:357:VAL:HG13	2.03	0.58
1:B:255:ALA:O	1:B:256:ALA:HB3	2.04	0.58
1:C:277:THR:HG22	1:C:278:LYS:HD2	1.85	0.58
1:B:227:LEU:HD22	1:B:253:PHE:CE2	2.38	0.58
1:A:279:THR:HG22	1:A:281:PRO:HD2	1.85	0.57
1:B:279:THR:HG22	1:B:281:PRO:HD3	1.85	0.57
1:B:298:ILE:O	1:B:302:LEU:HD13	2.04	0.57
1:C:475:ASP:C	1:C:476:HIS:ND1	2.58	0.57
1:C:304:LEU:HD12	1:C:305:ASN:HB2	1.87	0.57
1:B:370:VAL:O	1:B:372:LYS:HE2	2.04	0.57
1:C:484:PHE:O	1:C:485:TYR:HB2	2.05	0.57
1:C:305:ASN:HB3	1:C:488:LYS:HZ3	1.68	0.57
1:B:272:LYS:HA	1:B:272:LYS:HE2	1.87	0.56
1:A:325:GLY:O	1:A:326:LYS:CB	2.54	0.56
1:A:279:THR:HG22	1:A:280:ALA:H	1.68	0.56
1:A:232:VAL:N	2:A:496:HOH:O	2.01	0.56
1:A:254:ASN:ND2	1:A:263:ILE:HD11	2.20	0.56
1:A:223:ALA:HB3	1:A:226:GLU:CB	2.36	0.56
1:A:237:THR:H	1:A:241:GLN:NE2	2.04	0.56
1:C:317:LEU:HD23	1:C:317:LEU:C	2.26	0.56
1:A:321:THR:HG22	1:A:323:LYS:HB2	1.88	0.55
1:A:437:HIS:O	1:A:440:PRO:HD2	2.06	0.55
1:B:290:VAL:O	1:B:291:GLU:CB	2.55	0.55
1:B:462:THR:CG2	1:B:464:GLN:HB3	2.37	0.55
1:A:231:LEU:HD13	1:A:267:LEU:HD23	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:462:THR:CG2	1:B:465:GLU:HG3	2.35	0.55
1:A:349:HIS:HB2	1:A:463:LYS:HE3	1.89	0.55
1:A:382:MET:HE2	1:A:387:VAL:HG23	1.89	0.54
1:A:279:THR:CG2	1:A:280:ALA:N	2.70	0.54
1:C:304:LEU:HD12	1:C:305:ASN:N	2.23	0.54
1:C:329:THR:HA	1:C:416:VAL:O	2.07	0.54
1:C:346:ALA:O	1:C:466:VAL:HG21	2.08	0.54
1:B:237:THR:HB	1:B:241:GLN:OE1	2.07	0.54
1:C:304:LEU:HD12	1:C:305:ASN:CB	2.38	0.54
1:A:423:ASN:ND2	1:A:425:CYS:H	2.05	0.54
1:C:300:LYS:HG3	2:C:524:HOH:O	2.08	0.53
1:A:297:ARG:HH11	1:A:297:ARG:HG2	1.73	0.53
1:B:462:THR:HG23	1:B:464:GLN:N	2.24	0.53
1:A:280:ALA:HB3	1:A:281:PRO:HD3	1.89	0.53
1:B:335:PRO:HD2	1:B:450:LEU:O	2.09	0.53
1:B:280:ALA:CA	1:B:373:MET:HE3	2.37	0.53
1:A:256:ALA:O	1:A:257:SER:CB	2.56	0.53
1:A:398:LYS:HG3	2:B:514:HOH:O	2.09	0.53
1:C:375:ILE:O	1:C:416:VAL:HA	2.08	0.53
1:B:462:THR:HG22	1:B:465:GLU:H	1.71	0.53
1:C:336:ALA:O	1:C:337:THR:CB	2.57	0.53
1:C:463:LYS:H	1:C:463:LYS:CD	2.21	0.53
1:B:395:GLY:HA3	1:B:444:ARG:HH11	1.73	0.53
1:C:311:TYR:CE2	1:C:487:LYS:HB2	2.45	0.52
1:C:305:ASN:HB3	1:C:488:LYS:HZ2	1.75	0.52
1:B:335:PRO:HB2	1:B:338:THR:OG1	2.09	0.52
1:B:319:TRP:CD1	1:B:330:ILE:HG13	2.44	0.52
1:A:361:ASN:O	1:A:363:ASN:N	2.42	0.52
1:C:303:GLU:HA	1:C:303:GLU:OE1	2.10	0.52
1:A:225:MET:HG3	1:A:228:VAL:HG23	1.91	0.52
1:B:302:LEU:O	1:B:307:TYR:HB2	2.09	0.52
1:A:297:ARG:O	1:A:301:ILE:HG22	2.09	0.52
1:C:297:ARG:O	1:C:300:LYS:N	2.43	0.51
1:C:297:ARG:C	1:C:301:ILE:HG12	2.31	0.51
1:C:346:ALA:CB	1:C:461:VAL:HB	2.41	0.51
1:B:297:ARG:HG3	1:B:465:GLU:OE2	2.09	0.51
1:B:227:LEU:HD22	1:B:253:PHE:CZ	2.45	0.51
1:C:297:ARG:HD3	1:C:465:GLU:OE1	2.10	0.51
1:C:307:TYR:CD2	1:C:488:LYS:HA	2.46	0.51
1:C:232:VAL:HG21	1:C:273:ILE:HD11	1.91	0.51
1:C:415:PRO:O	1:C:416:VAL:HB	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:400:ARG:HG2	1:B:400:ARG:HH11	1.75	0.51
1:C:450:LEU:HD12	1:C:450:LEU:N	2.27	0.50
1:A:239:GLU:CD	1:A:260:ARG:HH22	2.14	0.50
1:A:434:THR:HG22	1:A:436:GLU:N	2.06	0.50
1:C:304:LEU:HD11	1:C:450:LEU:HA	1.94	0.50
1:B:293:ILE:N	1:B:293:ILE:HD12	2.06	0.50
1:C:461:VAL:O	1:C:462:THR:HB	2.11	0.50
1:C:377:TRP:CE3	1:C:418:VAL:HG22	2.47	0.50
1:A:279:THR:CG2	1:A:280:ALA:H	2.25	0.50
1:A:442:GLN:HG2	1:A:484:PHE:CZ	2.47	0.50
1:A:342:ASN:ND2	1:A:461:VAL:HG13	2.27	0.50
1:B:400:ARG:HA	1:B:409:ALA:O	2.12	0.49
1:C:311:TYR:CZ	1:C:487:LYS:HD3	2.47	0.49
1:B:368:ASP:O	1:B:372:LYS:HE3	2.11	0.49
1:C:316:PHE:CE2	1:C:330:ILE:HD12	2.47	0.49
1:A:346:ALA:CB	1:A:461:VAL:HG22	2.42	0.49
1:A:313:ALA:HB1	1:A:473:ALA:HB2	1.94	0.49
1:B:363:ASN:O	1:B:365:PRO:HD3	2.12	0.49
1:B:352:PRO:HD2	1:B:373:MET:CE	2.40	0.49
1:A:223:ALA:HB3	1:A:226:GLU:HB3	1.95	0.49
1:A:437:HIS:C	1:A:440:PRO:HD2	2.32	0.49
1:A:395:GLY:HA3	1:A:444:ARG:HD3	1.93	0.49
1:B:342:ASN:HD21	1:B:460:LYS:N	2.03	0.49
1:C:463:LYS:N	1:C:463:LYS:CD	2.74	0.49
1:A:273:ILE:HB	2:A:491:HOH:O	2.11	0.49
1:A:295:SER:HA	1:A:300:LYS:NZ	2.27	0.49
1:A:358:ASN:O	1:A:359:TRP:HB2	2.12	0.49
1:A:366:PHE:HE1	1:A:386:VAL:HG13	1.78	0.49
1:B:382:MET:HE3	1:B:387:VAL:HA	1.92	0.48
1:A:452:ARG:HA	1:A:453:ARG:HH21	1.76	0.48
1:C:297:ARG:N	1:C:297:ARG:HD2	2.28	0.48
1:A:453:ARG:HE	1:A:453:ARG:N	1.97	0.48
1:C:328:ASN:O	1:C:329:THR:HG23	2.12	0.48
1:C:255:ALA:O	1:C:256:ALA:HB3	2.14	0.48
1:C:444:ARG:O	1:C:445:MET:HB3	2.13	0.48
1:C:327:ARG:O	1:C:328:ASN:C	2.52	0.48
1:B:243:ILE:HG23	1:B:247:GLN:HG3	1.95	0.48
1:C:328:ASN:HB3	1:C:415:PRO:O	2.13	0.48
1:B:462:THR:HG23	1:B:464:GLN:HB3	1.94	0.48
1:A:398:LYS:HG2	1:A:413:PRO:HD3	1.95	0.48
1:B:331:TRP:CH2	1:B:424:MET:HB2	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:297:ARG:H	1:C:297:ARG:HD2	1.79	0.47
1:A:452:ARG:HG2	2:A:564:HOH:O	2.14	0.47
1:C:304:LEU:O	1:C:305:ASN:C	2.53	0.47
1:A:473:ALA:O	1:A:477:VAL:HG22	2.15	0.47
1:B:350:THR:HA	1:B:463:LYS:HE3	1.96	0.47
1:C:352:PRO:HD2	1:C:373:MET:SD	2.54	0.47
1:A:395:GLY:CA	1:A:444:ARG:HD3	2.45	0.47
1:A:267:LEU:O	1:A:270:ALA:O	2.33	0.47
1:C:484:PHE:O	1:C:485:TYR:CB	2.63	0.47
1:C:278:LYS:HB3	1:C:373:MET:CE	2.45	0.47
1:C:297:ARG:O	1:C:298:ILE:C	2.52	0.47
1:A:235:GLY:HA2	1:A:274:MET:SD	2.54	0.47
1:A:297:ARG:NH1	1:A:297:ARG:CG	2.75	0.47
1:A:301:ILE:HB	2:A:495:HOH:O	2.13	0.47
1:A:306:GLY:O	1:A:488:LYS:HB3	2.15	0.47
1:B:427:VAL:O	1:B:433:THR:HG23	2.14	0.47
1:B:284:LEU:O	1:B:467:LYS:HA	2.14	0.46
1:C:240:LYS:O	1:C:244:GLN:HG3	2.15	0.46
1:C:278:LYS:HB3	1:C:373:MET:HE1	1.97	0.46
1:A:423:ASN:HD22	1:A:424:MET:N	2.14	0.46
1:A:454:LEU:O	1:A:455:ASP:HB2	2.15	0.46
1:C:331:TRP:HA	1:C:418:VAL:O	2.15	0.46
1:A:295:SER:HA	1:A:300:LYS:HZ3	1.80	0.46
1:C:487:LYS:O	1:C:489:GLY:N	2.47	0.45
1:A:237:THR:H	1:A:241:GLN:HE22	1.65	0.45
1:C:236:ILE:HD13	1:C:242:TRP:HA	1.99	0.45
1:C:373:MET:O	1:C:415:PRO:HD2	2.16	0.45
1:A:239:GLU:OE2	1:A:260:ARG:NH2	2.49	0.45
1:C:413:PRO:HG2	2:C:548:HOH:O	2.16	0.45
1:C:322:LYS:O	1:C:324:PHE:N	2.50	0.45
1:B:361:ASN:ND2	1:B:361:ASN:C	2.68	0.45
1:B:401:VAL:O	1:B:408:SER:HA	2.16	0.45
1:C:377:TRP:HE3	1:C:418:VAL:HG22	1.81	0.45
1:C:256:ALA:O	1:C:257:SER:HB2	2.17	0.45
1:B:400:ARG:NH2	1:B:408:SER:OG	2.50	0.45
1:B:279:THR:CG2	1:B:281:PRO:HD2	2.46	0.44
1:A:227:LEU:HD13	1:A:253:PHE:CD2	2.52	0.44
1:A:405:CYS:O	1:A:406:LYS:C	2.56	0.44
1:B:337:THR:HG21	2:B:579:HOH:O	2.17	0.44
1:B:434:THR:CG2	1:B:436:GLU:HG2	2.47	0.44
1:A:321:THR:CG2	1:A:321:THR:O	2.66	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:279:THR:HG22	1:A:281:PRO:HD3	1.99	0.44
1:C:472:TRP:NE1	1:C:476:HIS:NE2	2.66	0.44
1:A:404:LYS:HA	1:A:404:LYS:HD3	1.76	0.44
1:B:437:HIS:C	1:B:440:PRO:HD2	2.38	0.44
1:C:238:SER:HB2	2:C:493:HOH:O	2.17	0.44
1:A:363:ASN:C	1:A:365:PRO:HD3	2.38	0.43
1:A:258:ASN:OD1	1:A:260:ARG:NH1	2.51	0.43
1:C:487:LYS:C	1:C:489:GLY:H	2.20	0.43
1:A:327:ARG:HG2	1:A:446:PHE:CE1	2.53	0.43
1:A:474:LYS:O	1:A:477:VAL:HG23	2.18	0.43
1:A:404:LYS:HZ2	1:B:405:CYS:HB3	1.82	0.43
1:C:331:TRP:CD1	1:C:332:LEU:N	2.87	0.43
1:B:296:ASN:O	1:B:300:LYS:HG3	2.18	0.43
1:A:256:ALA:O	1:A:257:SER:HB3	2.17	0.43
1:A:362:GLU:OE1	1:A:383:THR:HG21	2.19	0.43
1:B:460:LYS:O	1:B:461:VAL:O	2.36	0.43
1:B:258:ASN:C	1:B:260:ARG:H	2.22	0.43
1:C:343:ILE:HA	1:C:461:VAL:HG21	2.01	0.43
1:B:383:THR:CG2	1:B:386:VAL:HG22	2.48	0.43
1:C:346:ALA:HB2	1:C:461:VAL:HB	2.00	0.42
1:B:279:THR:CG2	1:B:280:ALA:N	2.82	0.42
1:A:442:GLN:HG2	1:A:484:PHE:CE1	2.54	0.42
1:B:394:LEU:HD23	1:B:416:VAL:HG11	2.01	0.42
1:C:231:LEU:HD13	1:C:267:LEU:HD23	2.01	0.42
1:C:314:SER:OG	1:C:477:VAL:HA	2.19	0.42
1:C:241:GLN:OE1	1:C:241:GLN:HA	2.20	0.42
1:C:486:VAL:O	1:C:488:LYS:N	2.53	0.42
1:A:274:MET:HG3	2:A:491:HOH:O	2.19	0.42
1:C:284:LEU:O	1:C:467:LYS:HA	2.20	0.42
1:C:260:ARG:NE	1:C:260:ARG:HA	2.34	0.42
1:B:283:TYR:HD1	1:B:373:MET:HE1	1.85	0.42
1:A:279:THR:CG2	1:A:281:PRO:HD2	2.48	0.42
1:A:361:ASN:C	1:A:363:ASN:N	2.72	0.42
1:C:409:ALA:O	1:C:410:GLN:CB	2.68	0.42
1:C:268:ASP:O	1:C:272:LYS:HB2	2.19	0.42
1:C:487:LYS:O	1:C:488:LYS:HB3	2.19	0.42
1:A:324:PHE:O	1:A:325:GLY:C	2.57	0.42
1:C:324:PHE:HB2	1:C:327:ARG:CB	2.50	0.41
1:C:462:THR:HG22	1:C:464:GLN:H	1.84	0.41
1:C:300:LYS:HA	1:C:303:GLU:HG2	2.01	0.41
1:B:294:SER:O	1:B:300:LYS:CE	2.64	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:255:ALA:C	1:C:257:SER:H	2.22	0.41
1:B:255:ALA:O	1:B:256:ALA:CB	2.67	0.41
1:B:386:VAL:O	1:B:386:VAL:HG23	2.19	0.41
1:A:399:VAL:HG12	1:A:400:ARG:N	2.36	0.41
1:B:462:THR:HG21	1:B:464:GLN:HB3	2.01	0.41
1:A:318:GLY:HA3	1:A:324:PHE:CE1	2.55	0.41
1:B:400:ARG:NH1	1:B:410:GLN:HG2	2.36	0.41
1:A:457:ASP:OD2	1:A:457:ASP:C	2.59	0.41
1:A:351:VAL:HG22	1:A:352:PRO:HD2	2.03	0.41
1:B:235:GLY:HA2	1:B:274:MET:SD	2.61	0.41
1:C:270:ALA:O	1:C:274:MET:HG3	2.21	0.41
1:A:397:SER:O	1:A:399:VAL:HG23	2.21	0.41
1:B:280:ALA:N	1:B:373:MET:HE3	2.36	0.41
1:A:274:MET:N	2:A:491:HOH:O	2.54	0.41
1:B:261:SER:O	1:B:262:GLN:CB	2.68	0.41
1:B:288:GLN:HA	1:B:289:PRO:HD2	1.89	0.41
1:B:313:ALA:HB1	1:B:473:ALA:CB	2.49	0.41
1:A:272:LYS:HA	1:A:272:LYS:HD3	1.95	0.41
1:B:462:THR:HG22	1:B:465:GLU:CB	2.51	0.41
1:B:370:VAL:HG22	1:B:411:ILE:HD13	2.03	0.41
1:C:264:LYS:HG2	1:C:268:ASP:OD2	2.21	0.40
1:A:351:VAL:HG22	1:A:373:MET:SD	2.62	0.40
1:B:400:ARG:HH11	1:B:410:GLN:HG2	1.87	0.40
1:C:412:ASP:O	1:C:414:THR:HG23	2.20	0.40
1:C:472:TRP:CD1	1:C:476:HIS:CE1	3.09	0.40
1:A:310:GLN:HB2	2:A:556:HOH:O	2.20	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	266/268 (99%)	243 (91%)	16 (6%)	7 (3%)	7	6
1	B	265/268 (99%)	238 (90%)	17 (6%)	10 (4%)	4	3
1	C	187/268 (70%)	145 (78%)	26 (14%)	16 (9%)	1	0
All	All	718/804 (89%)	626 (87%)	59 (8%)	33 (5%)	3	2

All (33) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	257	SER
1	A	325	GLY
1	A	326	LYS
1	A	406	LYS
1	C	298	ILE
1	C	323	LYS
1	C	328	ASN
1	C	329	THR
1	C	337	THR
1	C	477	VAL
1	A	358	ASN
1	A	362	GLU
1	A	404	LYS
1	B	256	ALA
1	B	262	GLN
1	B	291	GLU
1	B	452	ARG
1	B	461	VAL
1	C	306	GLY
1	C	416	VAL
1	C	422	THR
1	C	487	LYS
1	B	259	SER
1	B	405	CYS
1	B	460	LYS
1	C	297	ARG
1	C	413	PRO
1	B	260	ARG
1	C	325	GLY
1	C	410	GLN
1	B	290	VAL
1	C	257	SER
1	C	462	THR

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	217/229 (95%)	199 (92%)	18 (8%)	14	21
1	B	212/229 (93%)	202 (95%)	10 (5%)	32	50
1	C	132/229 (58%)	124 (94%)	8 (6%)	23	36
All	All	561/687 (82%)	525 (94%)	36 (6%)	22	34

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	260	ARG
1	A	297	ARG
1	A	315	VAL
1	A	321	THR
1	A	342	ASN
1	A	351	VAL
1	A	356	CYS
1	A	359	TRP
1	A	360	THR
1	A	361	ASN
1	A	364	PHE
1	A	377	TRP
1	A	383	THR
1	A	419	THR
1	A	423	ASN
1	A	433	THR
1	A	453	ARG
1	A	479	GLU
1	B	237	THR
1	B	253	PHE
1	B	293	ILE
1	B	338	THR
1	B	354	TYR
1	B	361	ASN
1	B	362	GLU
1	B	419	THR

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Mol	Chain	Res	Type
1	B	458	PHE
1	B	462	THR
1	C	237	THR
1	C	273	ILE
1	C	322	LYS
1	C	329	THR
1	C	415	PRO
1	C	463	LYS
1	C	478	VAL
1	C	483	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	241	GLN
1	A	254	ASN
1	A	269	ASN
1	A	287	GLN
1	A	342	ASN
1	A	349	HIS
1	A	361	ASN
1	A	423	ASN
1	A	476	HIS
1	B	305	ASN
1	B	361	ASN
1	B	363	ASN
1	B	438	GLN
1	B	464	GLN
1	B	482	HIS
1	C	305	ASN
1	C	482	HIS

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](#)

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](#)

EDS was not executed - this section will therefore be empty.

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

6.5 Other polymers [i](#)

EDS was not executed - this section will therefore be empty.