



Full wwPDB X-ray Structure Validation Report

Feb 26, 2014 – 10:08 PM GMT

PDB ID : 1ANV
Title : ADENOVIRUS 5 DBP/URANYL FLUORIDE SOAK
Authors : Kanellopoulos, P.N.; Tsernoglou, D.; Van Der Vliet, P.C.; Tucker, P.A.
Deposited on : 1996-03-14
Resolution : 2.70 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

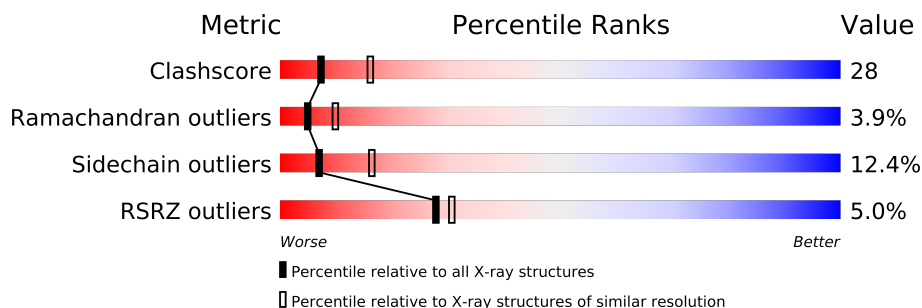
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.70 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	79885	1939 (2.70-2.70)
Ramachandran outliers	78287	1905 (2.70-2.70)
Sidechain outliers	78261	1905 (2.70-2.70)
RSRZ outliers	66119	1559 (2.70-2.70)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	356	

2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 2500 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

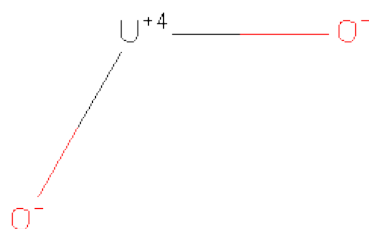
- Molecule 1 is a protein called ADENOVIRUS SINGLE-STRANDED DNA-BINDING PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	314	Total	C	N	O	S	0	0	0
			2463	1562	432	446	23			

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Zn	0	0
			2	2		

- Molecule 3 is URANYL (VI) ION (three-letter code: IUM) (formula: O₂U).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	U	0	0
			1	1		
3	A	1	Total	U	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	U	0	0
			1	1		
3	A	1	Total	U	0	0
			1	1		

- Molecule 4 is water.

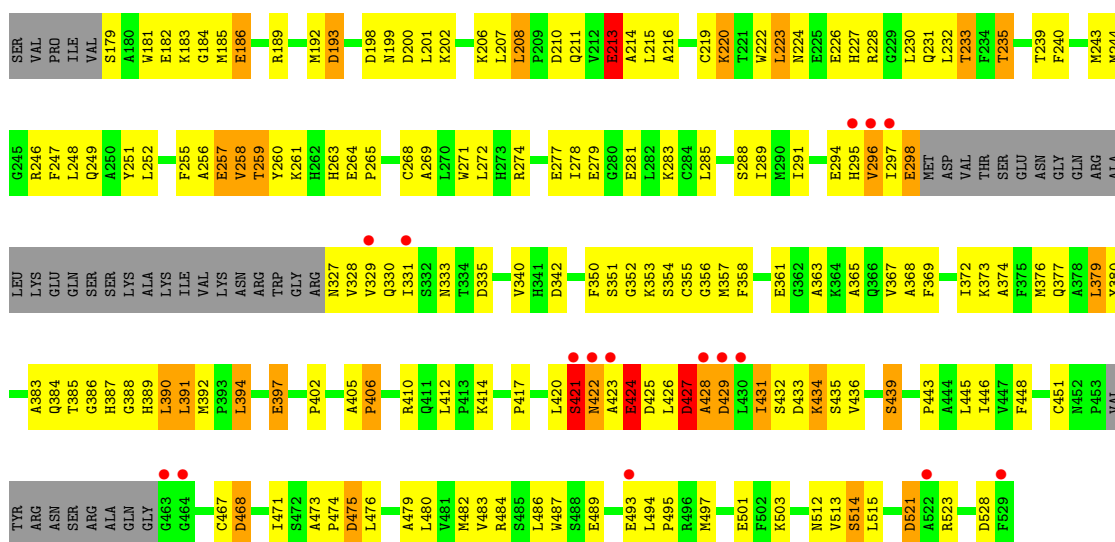
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	31	Total	O	0	0
			31	31		

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.

• Molecule 1: ADENOVIRUS SINGLE-STRANDED DNA-BINDING PROTEIN

Chain A: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	79.70Å 75.60Å 60.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.70 14.93 – 2.70	Depositor EDS
% Data completeness (in resolution range)	98.7 (15.00-2.70) 93.8 (14.93-2.70)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.79 (at 2.69Å)	Xtriage
Refinement program	TNT	Depositor
R, R_{free}	0.204 , 0.310 0.194 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	28.6	Xtriage
Anisotropy	0.144	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 71.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 10312 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	2500	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.73% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, IUM

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	0.77	13/2525 (0.5%)	1.09	23/3417 (0.7%)

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	493	GLU	CD-OE1	5.91	1.32	1.25
1	A	226	GLU	CD-OE2	5.79	1.32	1.25
1	A	361	GLU	CD-OE2	5.67	1.31	1.25
1	A	277	GLU	CD-OE1	5.61	1.31	1.25
1	A	294	GLU	CD-OE2	5.61	1.31	1.25
1	A	186	GLU	CD-OE1	5.54	1.31	1.25
1	A	489	GLU	CD-OE1	5.46	1.31	1.25
1	A	213	GLU	CD-OE2	5.44	1.31	1.25
1	A	264	GLU	CD-OE2	5.39	1.31	1.25
1	A	279	GLU	CD-OE2	5.35	1.31	1.25
1	A	182	GLU	CD-OE1	5.34	1.31	1.25
1	A	424	GLU	CD-OE2	5.30	1.31	1.25
1	A	397	GLU	CD-OE1	5.20	1.31	1.25

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	193	ASP	CB-CG-OD2	-8.15	110.96	118.30
1	A	475	ASP	CB-CG-OD1	-7.82	111.27	118.30
1	A	475	ASP	CB-CG-OD2	7.36	124.92	118.30
1	A	193	ASP	CB-CG-OD1	7.20	124.78	118.30
1	A	521	ASP	CB-CG-OD1	-6.42	112.52	118.30
1	A	198	ASP	CB-CG-OD1	-6.28	112.65	118.30
1	A	198	ASP	CB-CG-OD2	6.26	123.94	118.30
1	A	468	ASP	CB-CG-OD1	6.25	123.92	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	429	ASP	CB-CG-OD2	6.17	123.85	118.30
1	A	433	ASP	CB-CG-OD2	-6.08	112.83	118.30
1	A	468	ASP	CB-CG-OD2	-5.97	112.92	118.30
1	A	246	ARG	NE-CZ-NH1	5.90	123.25	120.30
1	A	200	ASP	CB-CG-OD2	5.90	123.61	118.30
1	A	200	ASP	CB-CG-OD1	-5.86	113.02	118.30
1	A	335	ASP	CB-CG-OD2	-5.85	113.03	118.30
1	A	429	ASP	CB-CG-OD1	-5.85	113.03	118.30
1	A	521	ASP	CB-CG-OD2	5.59	123.33	118.30
1	A	528	ASP	CB-CG-OD2	-5.46	113.39	118.30
1	A	210	ASP	CB-CG-OD1	-5.46	113.39	118.30
1	A	433	ASP	CB-CG-OD1	5.43	123.19	118.30
1	A	427	ASP	CB-CG-OD2	5.18	122.96	118.30
1	A	425	ASP	CB-CG-OD2	-5.11	113.70	118.30
1	A	528	ASP	CB-CG-OD1	5.04	122.83	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2463	0	2400	137	1
2	A	2	0	0	0	0
3	A	4	0	0	0	1
4	A	31	0	0	5	0
All	All	2500	0	2400	137	1

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 28.

All (137) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:269:ALA:HB2	1:A:394:LEU:HD21	1.38	1.05

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:235:THR:HB	1:A:471:ILE:HA	1.39	1.01
1:A:263:HIS:CE1	1:A:402:PRO:HG2	2.07	0.89
1:A:379:LEU:HB3	1:A:380:TYR:CD1	2.16	0.81
1:A:269:ALA:CB	1:A:394:LEU:HD21	2.12	0.78
1:A:222:TRP:CD1	1:A:486:LEU:HD13	2.23	0.74
1:A:181:TRP:HZ2	1:A:206:LYS:HD2	1.53	0.73
1:A:272:LEU:HD23	1:A:274:ARG:CZ	2.19	0.72
1:A:374:ALA:HA	1:A:377:GLN:HE21	1.54	0.71
1:A:295:HIS:HB2	1:A:333:ASN:HB2	1.74	0.70
1:A:230:LEU:HD12	1:A:231:GLN:H	1.57	0.70
1:A:199:ASN:HD22	1:A:202:LYS:HE3	1.56	0.69
1:A:373:LYS:O	1:A:377:GLN:HG3	1.92	0.69
1:A:514:SER:O	1:A:515:LEU:HD23	1.93	0.69
1:A:424:GLU:OE2	1:A:443:PRO:HB3	1.94	0.68
1:A:263:HIS:HE1	1:A:402:PRO:HG2	1.57	0.67
1:A:479:ALA:O	1:A:483:VAL:HG23	1.95	0.66
1:A:342:ASP:OD1	1:A:354:SER:HB2	1.96	0.65
1:A:291:ILE:N	1:A:291:ILE:HD13	2.13	0.64
1:A:473:ALA:HB3	1:A:474:PRO:HD3	1.79	0.63
1:A:421:SER:O	1:A:423:ALA:N	2.30	0.63
1:A:431:ILE:HG21	1:A:436:VAL:HG11	1.81	0.62
1:A:255:PHE:HB2	1:A:495:PRO:HG3	1.81	0.62
1:A:184:GLY:HA3	1:A:207:LEU:HD11	1.81	0.62
1:A:363:ALA:O	1:A:367:VAL:HG23	1.99	0.62
1:A:446:ILE:N	4:A:23:HOH:O	2.32	0.61
1:A:391:LEU:HD13	1:A:476:LEU:HG	1.83	0.61
1:A:278:ILE:HB	1:A:281:GLU:HB3	1.81	0.61
1:A:355:CYS:SG	1:A:357:MET:HB3	2.42	0.60
1:A:420:LEU:O	1:A:421:SER:C	2.39	0.60
1:A:247:PHE:HB2	1:A:483:VAL:HG21	1.84	0.60
1:A:353:LYS:NZ	4:A:2:HOH:O	2.32	0.60
1:A:181:TRP:CZ2	1:A:206:LYS:HD2	2.35	0.59
1:A:327:ASN:O	1:A:328:VAL:HG13	2.03	0.58
1:A:421:SER:O	1:A:422:ASN:C	2.41	0.58
1:A:421:SER:O	1:A:424:GLU:N	2.38	0.57
1:A:206:LYS:HB2	1:A:208:LEU:HD22	1.87	0.57
1:A:397:GLU:HB3	1:A:467:CYS:SG	2.44	0.57
1:A:424:GLU:HG2	1:A:424:GLU:O	2.04	0.57
1:A:487:TRP:CD1	1:A:494:LEU:HD12	2.39	0.57
1:A:434:LYS:HZ2	1:A:434:LYS:HB2	1.70	0.56
1:A:397:GLU:HB3	1:A:467:CYS:CB	2.36	0.56
1:A:372:ILE:O	1:A:376:MET:HG3	2.06	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:331:ILE:N	1:A:331:ILE:HD12	2.21	0.55
1:A:394:LEU:HA	1:A:412:LEU:HD22	1.88	0.55
1:A:379:LEU:HD22	1:A:380:TYR:HE1	1.72	0.55
1:A:379:LEU:HB3	1:A:380:TYR:CE1	2.41	0.55
1:A:223:LEU:O	1:A:227:HIS:N	2.40	0.55
1:A:199:ASN:HD22	1:A:202:LYS:CE	2.20	0.55
1:A:445:LEU:HA	4:A:23:HOH:O	2.06	0.55
1:A:230:LEU:HD12	1:A:231:GLN:N	2.22	0.54
1:A:365:ALA:O	1:A:369:PHE:HD2	1.90	0.54
1:A:514:SER:C	1:A:515:LEU:HD23	2.28	0.54
1:A:386:GLY:O	1:A:388:GLY:N	2.35	0.54
1:A:392:MET:CE	1:A:394:LEU:HD11	2.38	0.53
1:A:183:LYS:HD3	1:A:256:ALA:HA	1.89	0.53
1:A:185:MET:O	1:A:189:ARG:HB2	2.09	0.53
1:A:476:LEU:O	1:A:476:LEU:HD12	2.10	0.52
1:A:235:THR:HG21	1:A:475:ASP:CB	2.39	0.52
1:A:230:LEU:HD21	1:A:482:MET:CE	2.39	0.52
1:A:410:ARG:HG2	1:A:412:LEU:O	2.10	0.51
1:A:272:LEU:O	1:A:274:ARG:HD2	2.11	0.50
1:A:232:LEU:HB3	1:A:235:THR:HG23	1.94	0.50
1:A:379:LEU:HD22	1:A:380:TYR:CE1	2.46	0.50
1:A:368:ALA:HB1	1:A:446:ILE:HG13	1.94	0.50
1:A:268:CYS:HB3	1:A:392:MET:O	2.12	0.49
1:A:208:LEU:HB2	1:A:211:GLN:OE1	2.11	0.49
1:A:420:LEU:O	1:A:422:ASN:OD1	2.30	0.49
1:A:421:SER:H	1:A:443:PRO:HB2	1.78	0.49
1:A:369:PHE:O	1:A:372:ILE:HG22	2.12	0.49
1:A:383:ALA:O	1:A:385:THR:N	2.45	0.49
1:A:260:TYR:HB2	1:A:263:HIS:HB2	1.95	0.48
1:A:480:LEU:O	1:A:484:ARG:HG3	2.12	0.48
1:A:421:SER:OG	1:A:422:ASN:N	2.45	0.48
1:A:417:PRO:HA	1:A:445:LEU:O	2.13	0.48
1:A:256:ALA:O	1:A:257:GLU:HB2	2.14	0.48
1:A:298:GLU:OE2	1:A:330:GLN:OE1	2.30	0.48
1:A:271:TRP:CE2	1:A:285:LEU:HD23	2.49	0.48
1:A:263:HIS:ND1	1:A:402:PRO:HG2	2.27	0.47
1:A:431:ILE:HG21	1:A:436:VAL:CG1	2.43	0.47
1:A:296:VAL:C	1:A:297:ILE:HG12	2.35	0.47
1:A:272:LEU:HG	1:A:272:LEU:O	2.15	0.47
1:A:431:ILE:HG22	1:A:432:SER:N	2.29	0.47
1:A:352:GLY:CA	1:A:512:ASN:ND2	2.78	0.47
1:A:350:PHE:N	1:A:350:PHE:CD1	2.83	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:434:LYS:NZ	1:A:434:LYS:HB2	2.30	0.46
1:A:219:CYS:O	1:A:223:LEU:HB2	2.15	0.46
1:A:239:THR:O	1:A:243:MET:HG2	2.16	0.46
1:A:480:LEU:O	1:A:480:LEU:HD12	2.16	0.46
1:A:222:TRP:O	1:A:223:LEU:C	2.54	0.45
1:A:340:VAL:HG12	4:A:1:HOH:O	2.16	0.45
1:A:206:LYS:O	1:A:211:GLN:OE1	2.34	0.45
1:A:281:GLU:OE2	1:A:283:LYS:HE3	2.16	0.45
1:A:405:ALA:O	1:A:406:PRO:C	2.55	0.45
1:A:295:HIS:CE1	1:A:358:PHE:CZ	3.05	0.45
1:A:352:GLY:N	1:A:512:ASN:ND2	2.65	0.45
1:A:230:LEU:HD21	1:A:482:MET:HE1	1.98	0.44
1:A:352:GLY:HA2	1:A:512:ASN:HD22	1.82	0.44
1:A:427:ASP:O	1:A:428:ALA:HB2	2.17	0.44
1:A:296:VAL:HG13	1:A:330:GLN:HG3	1.99	0.44
1:A:251:TYR:CE1	1:A:497:MET:HB2	2.53	0.44
1:A:189:ARG:HA	1:A:192:MET:HE2	2.00	0.44
1:A:261:LYS:HA	4:A:20:HOH:O	2.17	0.44
1:A:269:ALA:HB2	1:A:394:LEU:CD2	2.28	0.44
1:A:392:MET:HE3	1:A:394:LEU:CD1	2.47	0.44
1:A:208:LEU:HD12	1:A:208:LEU:HA	1.73	0.43
1:A:410:ARG:HG2	1:A:412:LEU:H	1.82	0.43
1:A:258:VAL:HG23	1:A:259:THR:N	2.33	0.43
1:A:414:LYS:O	1:A:448:PHE:HA	2.18	0.43
1:A:521:ASP:C	1:A:523:ARG:H	2.21	0.43
1:A:379:LEU:HB3	1:A:380:TYR:HD1	1.78	0.43
1:A:220:LYS:HE3	1:A:220:LYS:HB2	1.56	0.43
1:A:513:VAL:HG22	1:A:515:LEU:HD21	2.00	0.43
1:A:420:LEU:CD2	1:A:439:SER:HB3	2.49	0.42
1:A:329:VAL:O	1:A:331:ILE:HD12	2.19	0.42
1:A:352:GLY:HA2	1:A:512:ASN:ND2	2.35	0.42
1:A:215:LEU:HD12	1:A:249:GLN:HB2	2.02	0.42
1:A:230:LEU:HD21	1:A:482:MET:HE3	2.02	0.41
1:A:420:LEU:HD23	1:A:424:GLU:OE1	2.18	0.41
1:A:240:PHE:O	1:A:244:MET:HG3	2.20	0.41
1:A:223:LEU:HD12	1:A:227:HIS:HB2	2.02	0.41
1:A:383:ALA:C	1:A:385:THR:H	2.23	0.41
1:A:350:PHE:CD2	1:A:356:GLY:HA2	2.55	0.41
1:A:295:HIS:CE1	1:A:358:PHE:HZ	2.38	0.41
1:A:422:ASN:OD1	1:A:423:ALA:N	2.44	0.41
1:A:271:TRP:O	1:A:389:HIS:HA	2.21	0.41
1:A:426:LEU:O	1:A:427:ASP:HB3	2.20	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:392:MET:HE2	1:A:394:LEU:HD11	2.03	0.41
1:A:412:LEU:O	1:A:451:CYS:HA	2.21	0.41
1:A:271:TRP:HE3	1:A:390:LEU:HD23	1.85	0.41
1:A:410:ARG:O	1:A:451:CYS:HB3	2.22	0.40
1:A:230:LEU:CD2	1:A:482:MET:HE3	2.51	0.40
1:A:213:GLU:CD	1:A:214:ALA:H	2.25	0.40
1:A:213:GLU:O	1:A:216:ALA:N	2.55	0.40
1:A:392:MET:CE	1:A:394:LEU:CD1	3.00	0.40
1:A:222:TRP:CG	1:A:486:LEU:HD13	2.57	0.40
1:A:230:LEU:CD2	1:A:482:MET:CE	2.99	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:193:ASP:OD1	3:A:532:IUM:U[2_665]	2.12	0.08

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

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	308/356 (86%)	262 (85%)	34 (11%)	12 (4%)	5	10

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	422	ASN
1	A	428	ALA
1	A	431	ILE
1	A	233	THR
1	A	384	GLN
1	A	427	ASP
1	A	429	ASP
1	A	224	ASN
1	A	387	HIS

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Mol	Chain	Res	Type
1	A	421	SER
1	A	296	VAL
1	A	265	PRO

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	267/303 (88%)	234 (88%)	33 (12%)	7 16

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	179	SER
1	A	186	GLU
1	A	201	LEU
1	A	208	LEU
1	A	213	GLU
1	A	220	LYS
1	A	223	LEU
1	A	228	ARG
1	A	233	THR
1	A	235	THR
1	A	248	LEU
1	A	252	LEU
1	A	257	GLU
1	A	258	VAL
1	A	259	THR
1	A	288	SER
1	A	289	ILE
1	A	298	GLU
1	A	351	SER
1	A	379	LEU
1	A	390	LEU
1	A	391	LEU
1	A	394	LEU
1	A	406	PRO
1	A	421	SER

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Mol	Chain	Res	Type
1	A	424	GLU
1	A	434	LYS
1	A	435	SER
1	A	439	SER
1	A	468	ASP
1	A	501	GLU
1	A	503	LYS
1	A	514	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	199	ASN
1	A	295	HIS
1	A	327	ASN
1	A	333	ASN
1	A	377	GLN
1	A	382	ASN
1	A	512	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

Of 6 ligands modelled in this entry, 2 are monoatomic and 4 are modelled with single atom - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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, 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	314/356 (88%)	-0.25	16 (5%)	27 30	6, 26, 70, 99	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	297	ILE	7.0
1	A	428	ALA	6.5
1	A	421	SER	5.3
1	A	430	LEU	3.9
1	A	296	VAL	3.6
1	A	422	ASN	3.2
1	A	429	ASP	3.0
1	A	329	VAL	3.0
1	A	493	GLU	2.8
1	A	463	GLY	2.7
1	A	423	ALA	2.3
1	A	522	ALA	2.2
1	A	331	ILE	2.2
1	A	529	PHE	2.2
1	A	464	GLY	2.0
1	A	295	HIS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	ZN	A	530	1/1	0.06	-1.80	21,21,21,21	0
2	ZN	A	531	1/1	0.02	-2.06	25,25,25,25	0
3	IUM	A	532	1/3	0.04	-2.45	42,42,42,42	0
3	IUM	A	533	1/3	0.03	-2.65	56,56,56,56	0
3	IUM	A	535	1/3	0.03	-3.51	57,57,57,57	0
3	IUM	A	534	1/3	0.04	-4.99	61,61,61,61	0

6.5 Other polymers

There are no such residues in this entry.