



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 04:58 PM UTC

PDB ID : 2C0P / pdb_00002c0p
Title : Aged form of mouse acetylcholinesterase inhibited by tabun
Authors : Ekstrom, F.; Akfur, C.; Tunemalm, A.-K.; Lundberg, S.
Deposited on : 2005-09-07
Resolution : 2.50 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

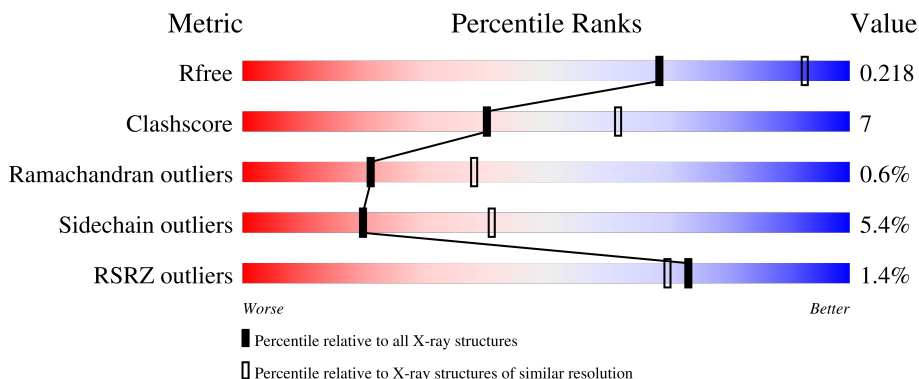
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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 $\leq 5\%$. 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.

Mol	Chain	Length	Quality of chain
1	A	548	% 79% 18% ..
1	B	548	% 81% 15% ..

2 Entry composition [i](#)

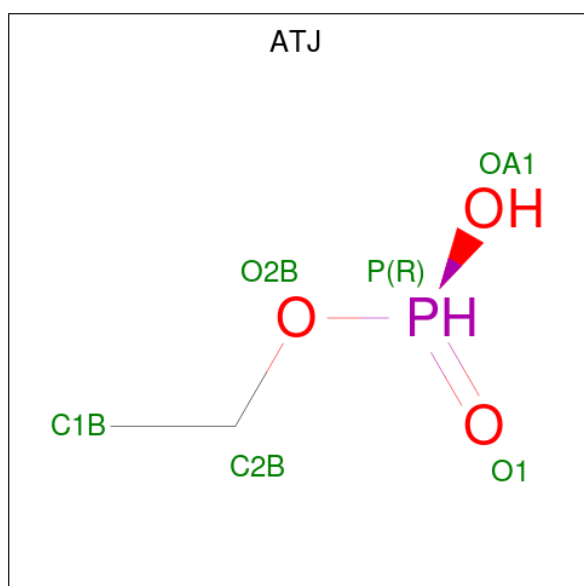
There are 3 unique types of molecules in this entry. The entry contains 8771 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 ACETYLCHOLINESTERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	535	Total	C	N	O	S	0	0	0
			4177	2679	725	759	14			
1	B	534	Total	C	N	O	S	0	0	1
			4159	2670	719	756	14			

- Molecule 2 is ETHYL HYDROGEN PHOSPHONATE (CCD ID: ATJ) (formula: $C_2H_7O_3P$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	P	0	0
			6	2	3	1		
2	B	1	Total	C	O	P	0	0
			6	2	3	1		

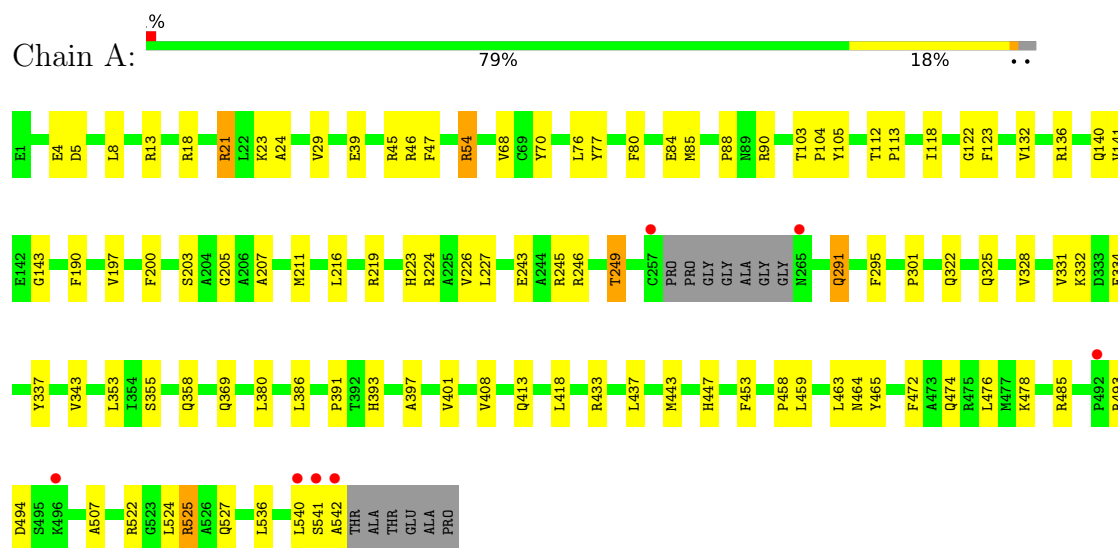
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	258	Total 258	O 258	0	0
3	B	165	Total 165	O 165	0	0

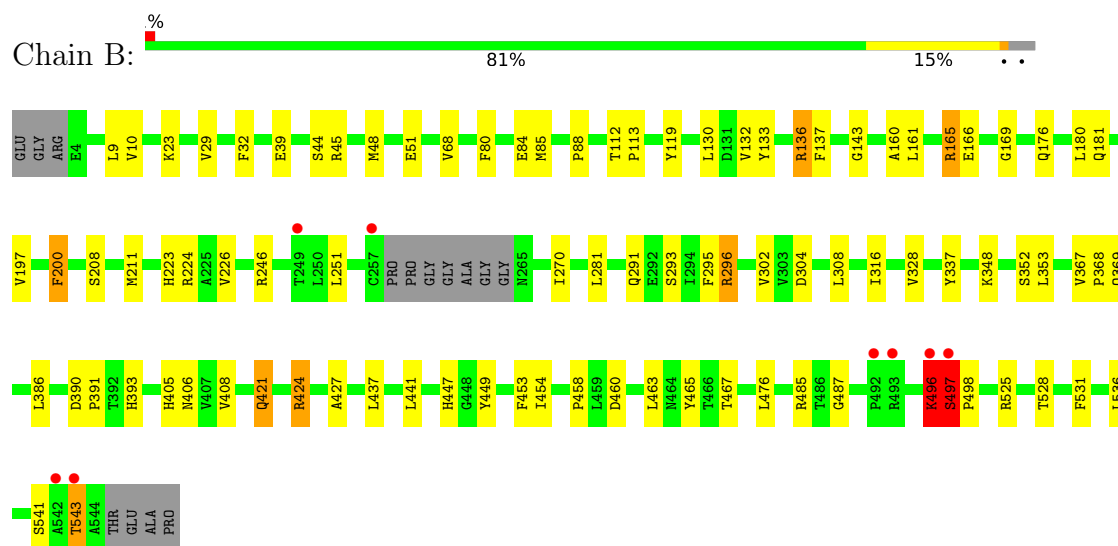
3 Residue-property plots [i](#)

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: ACETYLCHOLINESTERASE



• Molecule 1: ACETYLCHOLINESTERASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	79.02Å 110.88Å 226.38Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.82 – 2.50 28.82 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.2 (28.82-2.50) 99.1 (28.82-2.50)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.49 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.182 , 0.220 0.181 , 0.218	Depositor DCC
R_{free} test set	1381 reflections (2.00%)	wwPDB-VP
Wilson B-factor (Å ²)	49.6	Xtriage
Anisotropy	0.117	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 33.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8771	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.73% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ 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](#)

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

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	0.68	0/4300	0.91	1/5875 (0.0%)
1	B	0.69	1/4282 (0.0%)	0.91	3/5853 (0.1%)
All	All	0.68	1/8582 (0.0%)	0.91	4/11728 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	543	THR	C-N	-6.80	1.23	1.33

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	39	GLU	CA-C-N	-5.95	114.25	120.38
1	B	39	GLU	C-N-CA	-5.95	114.25	120.38
1	A	343	VAL	N-CA-C	5.51	113.01	107.55
1	B	454	ILE	CB-CA-C	-5.41	104.84	112.14

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	4177	0	4064	57	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	4159	0	4046	50	0
2	A	6	0	5	1	0
2	B	6	0	5	1	0
3	A	258	0	0	6	0
3	B	165	0	0	4	0
All	All	8771	0	8120	108	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 7.

All (108) 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:424:ARG:HH11	1:B:424:ARG:HG3	1.14	1.10
1:B:197:VAL:H	1:B:223:HIS:HD2	1.10	0.96
1:A:54:ARG:HH11	1:A:54:ARG:HG2	1.28	0.94
1:B:424:ARG:HG3	1:B:424:ARG:NH1	1.81	0.85
1:B:197:VAL:H	1:B:223:HIS:CD2	1.97	0.83
1:A:433:ARG:HD3	3:A:2225:HOH:O	1.78	0.81
1:A:4:GLU:OE2	1:A:18:ARG:HD3	1.80	0.80
1:A:369:GLN:HB2	3:A:2179:HOH:O	1.81	0.80
1:A:197:VAL:H	1:A:223:HIS:HD2	1.32	0.77
1:B:424:ARG:HH11	1:B:424:ARG:CG	1.97	0.75
1:B:496:LYS:HD3	1:B:496:LYS:H	1.52	0.72
1:B:112:THR:HG21	1:B:143:GLY:O	1.89	0.71
1:A:337:TYR:HA	1:A:443:MET:HE3	1.73	0.71
1:B:460:ASP:HB3	1:B:463:LEU:HD12	1.74	0.70
1:A:54:ARG:HH11	1:A:54:ARG:CG	2.03	0.69
1:B:458:PRO:HA	1:B:465:TYR:CD2	2.28	0.68
1:B:211:MET:HG2	1:B:308:LEU:HD21	1.77	0.66
1:A:245:ARG:O	1:A:249:THR:HG22	1.96	0.65
1:B:296:ARG:HG2	3:B:2094:HOH:O	1.96	0.64
1:A:54:ARG:HG2	1:A:54:ARG:NH1	2.07	0.63
1:B:496:LYS:HD3	1:B:496:LYS:N	2.10	0.63
1:A:113:PRO:HG2	1:A:485:ARG:HG2	1.83	0.61
1:A:245:ARG:O	1:A:249:THR:CG2	2.49	0.60
1:A:76:LEU:HD23	1:A:77:TYR:CE2	2.37	0.60
1:B:316:ILE:O	1:B:421:GLN:NE2	2.33	0.60
1:A:337:TYR:HA	1:A:443:MET:CE	2.32	0.59
1:B:447:HIS:HE1	2:B:1544:ATJ:P	2.26	0.58
1:A:224:ARG:HG2	1:A:325:GLN:HB2	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:112:THR:HG21	1:A:143:GLY:O	2.03	0.57
1:B:296:ARG:NH2	1:B:406:ASN:OD1	2.37	0.57
1:A:227:LEU:HB2	1:A:328:VAL:HG12	1.87	0.57
1:A:291:GLN:HE22	1:A:369:GLN:NE2	2.03	0.57
1:B:80:PHE:O	1:B:84:GLU:HG2	2.05	0.56
1:B:166:GLU:HG2	1:B:270:ILE:HD13	1.88	0.55
1:A:104:PRO:HG3	1:A:143:GLY:HA2	1.90	0.53
2:A:1543:ATJ:H2B2	3:A:2258:HOH:O	2.07	0.53
1:B:224:ARG:HD3	1:B:487:GLY:HA2	1.91	0.53
1:B:85:MET:CE	1:B:132:VAL:HG11	2.39	0.53
1:A:207:ALA:O	1:A:211:MET:HG2	2.09	0.52
1:B:328:VAL:O	1:B:427:ALA:HA	2.10	0.52
1:B:224:ARG:HD3	1:B:487:GLY:CA	2.40	0.52
1:B:160:ALA:HB2	1:B:169:GLY:HA3	1.91	0.51
1:B:29:VAL:HG21	1:B:136:ARG:HB2	1.93	0.51
1:A:24:ALA:HB3	1:A:140:GLN:HG3	1.92	0.50
1:B:224:ARG:HH11	1:B:224:ARG:HG3	1.75	0.50
1:B:165:ARG:HH11	1:B:165:ARG:HB3	1.77	0.49
1:B:293:SER:HB3	1:B:368:PRO:HB3	1.94	0.49
1:B:437:LEU:HD11	1:B:449:TYR:CD2	2.47	0.49
1:B:200:PHE:CB	1:B:226:VAL:HB	2.42	0.49
1:A:103:THR:HG21	1:A:190:PHE:HB3	1.95	0.49
1:A:507:ALA:O	1:A:522:ARG:NH1	2.46	0.49
1:B:44:SER:O	1:B:48:MET:HE3	2.14	0.48
1:A:5:ASP:HB3	1:A:8:LEU:HD12	1.96	0.48
1:B:224:ARG:HG3	1:B:224:ARG:NH1	2.28	0.48
1:B:390:ASP:OD2	1:B:393:HIS:ND1	2.47	0.48
1:B:68:VAL:HG11	1:B:88:PRO:HB3	1.96	0.48
1:A:458:PRO:HA	1:A:465:TYR:CD2	2.49	0.47
1:B:160:ALA:HB2	1:B:169:GLY:CA	2.44	0.47
1:A:203:SER:HB2	1:A:447:HIS:NE2	2.28	0.47
1:A:433:ARG:CZ	1:A:437:LEU:HD23	2.45	0.47
1:A:219:ARG:HA	1:A:219:ARG:HD2	1.65	0.47
1:A:29:VAL:HG21	1:A:136:ARG:HB2	1.97	0.47
1:A:80:PHE:O	1:A:84:GLU:HG2	2.14	0.47
1:B:165:ARG:HB3	1:B:165:ARG:NH1	2.30	0.46
1:B:113:PRO:HG2	1:B:485:ARG:HG2	1.98	0.46
1:B:353:LEU:HB3	1:B:391:PRO:HB2	1.98	0.45
1:B:393:HIS:HB3	3:B:2116:HOH:O	2.17	0.45
1:A:200:PHE:CB	1:A:226:VAL:HB	2.46	0.45
1:A:291:GLN:HE22	1:A:369:GLN:HE22	1.63	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:68:VAL:HG23	1:A:90:ARG:HB2	1.99	0.45
1:A:291:GLN:NE2	1:A:369:GLN:NE2	2.64	0.45
1:A:39:GLU:OE2	1:A:54:ARG:HB2	2.17	0.45
1:B:136:ARG:HG2	1:B:137:PHE:N	2.32	0.45
1:B:176:GLN:OE1	1:B:208:SER:HB3	2.18	0.44
1:A:211:MET:HE3	1:A:301:PRO:HG2	1.99	0.44
1:B:369:GLN:HE22	1:B:405:HIS:CE1	2.36	0.44
1:A:141:VAL:HG21	1:A:459:LEU:HD23	2.00	0.43
1:A:397:ALA:O	1:A:401:VAL:HG23	2.17	0.43
1:A:453:PHE:HB3	1:A:476:LEU:HD12	2.00	0.43
1:B:453:PHE:HB3	1:B:476:LEU:HD12	1.99	0.43
1:A:355:SER:OG	1:A:358:GLN:HG3	2.18	0.43
1:A:393:HIS:HB2	3:A:2209:HOH:O	2.17	0.43
1:A:459:LEU:HD11	1:A:474:GLN:HG3	2.00	0.43
1:B:48:MET:SD	3:B:2066:HOH:O	2.61	0.43
1:A:46:ARG:HD3	1:A:47:PHE:CZ	2.54	0.43
1:A:21:ARG:HG3	1:A:105:TYR:CE1	2.54	0.43
1:A:105:TYR:CD2	1:A:105:TYR:C	2.97	0.43
1:A:527:GLN:HG3	3:A:2252:HOH:O	2.19	0.42
1:B:528:THR:O	1:B:531:PHE:HB3	2.20	0.42
1:A:472:PHE:CZ	1:A:476:LEU:HD11	2.54	0.42
1:B:45:ARG:HD2	1:B:51:GLU:OE1	2.20	0.42
1:A:122:GLY:O	1:A:123:PHE:HB2	2.20	0.42
1:A:243:GLU:OE1	1:A:246:ARG:NH1	2.53	0.42
1:A:541:SER:O	1:A:542:ALA:HB2	2.20	0.42
1:B:10:VAL:HG23	1:B:32:PHE:CE2	2.55	0.42
1:A:85:MET:CE	1:A:132:VAL:HG11	2.49	0.42
1:A:118:ILE:O	1:A:205:GLY:HA3	2.20	0.41
1:B:302:VAL:HG13	1:B:304:ASP:HB3	2.02	0.41
1:B:408:VAL:HG11	1:B:525:ARG:HG3	2.01	0.41
1:A:331:VAL:HG22	1:A:334:GLU:CD	2.44	0.41
1:A:88:PRO:HD3	3:A:2060:HOH:O	2.20	0.41
1:B:130:LEU:HD12	1:B:133:TYR:CE2	2.56	0.41
1:A:408:VAL:HG11	1:A:525:ARG:HG3	2.03	0.41
1:B:497:SER:H	1:B:498:PRO:HA	1.84	0.41
1:A:463:LEU:O	1:A:464:ASN:HB2	2.21	0.40
1:A:353:LEU:HB3	1:A:391:PRO:HB2	2.03	0.40
1:A:54:ARG:CG	1:A:54:ARG:NH1	2.72	0.40
1:B:119:TYR:HB2	3:B:2057:HOH:O	2.20	0.40

There are no symmetry-related clashes.

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	531/548 (97%)	506 (95%)	24 (4%)	1 (0%)	43	63
1	B	530/548 (97%)	500 (94%)	25 (5%)	5 (1%)	14	27
All	All	1061/1096 (97%)	1006 (95%)	49 (5%)	6 (1%)	21	38

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	494	ASP
1	B	337	TYR
1	B	543	THR
1	B	496	LYS
1	B	497	SER
1	B	541	SER

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	439/446 (98%)	417 (95%)	22 (5%)	22	44
1	B	438/446 (98%)	413 (94%)	25 (6%)	18	39
All	All	877/892 (98%)	830 (95%)	47 (5%)	20	41

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	ARG
1	A	21	ARG
1	A	23	LYS
1	A	45	ARG
1	A	54	ARG
1	A	70	TYR
1	A	216	LEU
1	A	249	THR
1	A	291	GLN
1	A	295	PHE
1	A	322	GLN
1	A	332	LYS
1	A	380	LEU
1	A	386	LEU
1	A	413	GLN
1	A	418	LEU
1	A	478	LYS
1	A	493	ARG
1	A	524	LEU
1	A	525	ARG
1	A	536	LEU
1	A	540	LEU
1	B	9	LEU
1	B	23	LYS
1	B	136	ARG
1	B	161	LEU
1	B	165	ARG
1	B	180	LEU
1	B	181	GLN
1	B	200	PHE
1	B	246	ARG
1	B	251	LEU
1	B	281	LEU
1	B	291	GLN
1	B	295	PHE
1	B	296	ARG
1	B	348	LYS
1	B	352	SER
1	B	367	VAL
1	B	386	LEU
1	B	421	GLN
1	B	424	ARG
1	B	441	LEU

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Mol	Chain	Res	Type
1	B	467	THR
1	B	496	LYS
1	B	497	SER
1	B	536	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	265	ASN
1	A	284	HIS
1	A	287	HIS
1	A	291	GLN
1	A	421	GLN
1	A	474	GLN
1	B	87	ASN
1	B	181	GLN
1	B	223	HIS
1	B	291	GLN
1	B	405	HIS
1	B	447	HIS
1	B	474	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	ATJ	A	1543	1	2,5,5	0.56	0	1,5,5	0.35	0
2	ATJ	B	1544	1	2,5,5	0.87	0	1,5,5	0.73	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ATJ	A	1543	1	-	0/1/3/3	-
2	ATJ	B	1544	1	-	0/1/3/3	-

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.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1543	ATJ	1	0
2	B	1544	ATJ	1	0

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	535/548 (97%)	-0.51	7 (1%) 75 71	29, 45, 70, 114	0
1	B	534/548 (97%)	-0.32	8 (1%) 72 68	31, 51, 77, 117	0
All	All	1069/1096 (97%)	-0.41	15 (1%) 73 70	29, 47, 75, 117	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	542	ALA	4.7
1	A	542	ALA	3.8
1	B	496	LYS	3.8
1	A	496	LYS	3.5
1	A	540	LEU	3.5
1	B	543	THR	3.1
1	B	257	CYS	3.0
1	A	257	CYS	2.9
1	B	497	SER	2.7
1	B	492	PRO	2.3
1	A	265	ASN	2.3
1	A	541	SER	2.2
1	B	493	ARG	2.2
1	A	492	PRO	2.1
1	B	249	THR	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

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. 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	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	ATJ	B	1544	6/6	0.90	0.17	42,45,61,63	0
2	ATJ	A	1543	6/6	0.96	0.13	38,40,56,59	0

6.5 Other polymers [i](#)

There are no such residues in this entry.