



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:46 PM UTC

PDB ID : 3UE9 / pdb_00003ue9
Title : Crystal structure of Adenylosuccinate synthetase (AMPSase) (purA) from *Burkholderia thailandensis*
Authors : Seattle Structural Genomics Center for Infectious Disease (SSGCID)
Deposited on : 2011-10-28
Resolution : 1.95 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

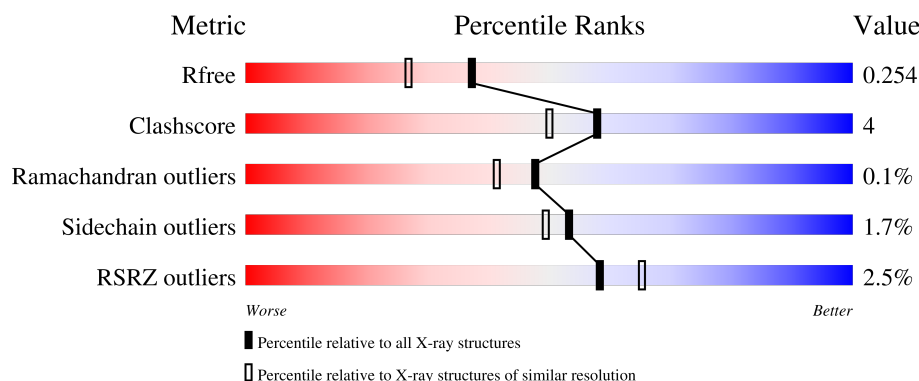
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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	452	3% 85% 9% 6%
1	B	452	3% 83% 9% 7%
1	C	452	0% 84% 10% 6%
1	D	452	2% 86% 8% 6%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 13889 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 Adenylosuccinate synthetase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	424	Total	C	N	O	S	0	6	0
			3202	2017	569	606	10			
1	B	421	Total	C	N	O	S	0	3	0
			3161	1989	569	592	11			
1	C	426	Total	C	N	O	S	0	3	0
			3203	2018	569	606	10			
1	D	425	Total	C	N	O	S	0	1	0
			3183	2003	566	604	10			

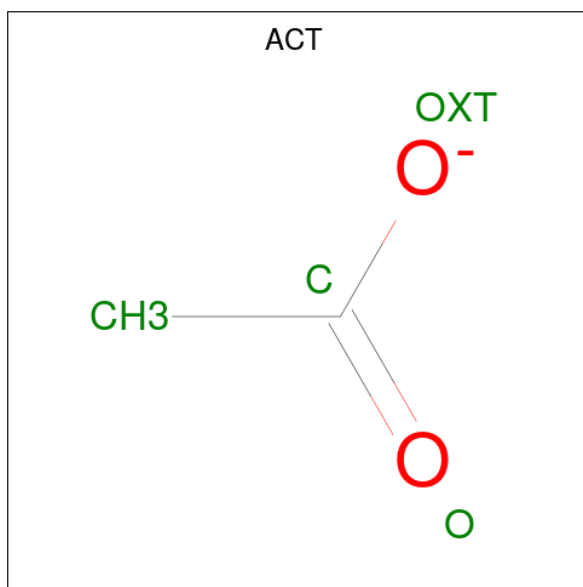
There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	expression tag	UNP Q2SWD3
A	-2	PRO	-	expression tag	UNP Q2SWD3
A	-1	GLY	-	expression tag	UNP Q2SWD3
A	0	SER	-	expression tag	UNP Q2SWD3
B	-3	GLY	-	expression tag	UNP Q2SWD3
B	-2	PRO	-	expression tag	UNP Q2SWD3
B	-1	GLY	-	expression tag	UNP Q2SWD3
B	0	SER	-	expression tag	UNP Q2SWD3
C	-3	GLY	-	expression tag	UNP Q2SWD3
C	-2	PRO	-	expression tag	UNP Q2SWD3
C	-1	GLY	-	expression tag	UNP Q2SWD3
C	0	SER	-	expression tag	UNP Q2SWD3
D	-3	GLY	-	expression tag	UNP Q2SWD3
D	-2	PRO	-	expression tag	UNP Q2SWD3
D	-1	GLY	-	expression tag	UNP Q2SWD3
D	0	SER	-	expression tag	UNP Q2SWD3

- Molecule 2 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Na 1 1	0	0
2	B	1	Total Na 1 1	0	0

- Molecule 3 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	C	1	Total C O 4 2 2	0	0
3	D	1	Total C O 4 2 2	0	0

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	C	1	Total Ca 1 1	0	0
4	D	1	Total Ca 1 1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	291	Total O 291 291	0	0

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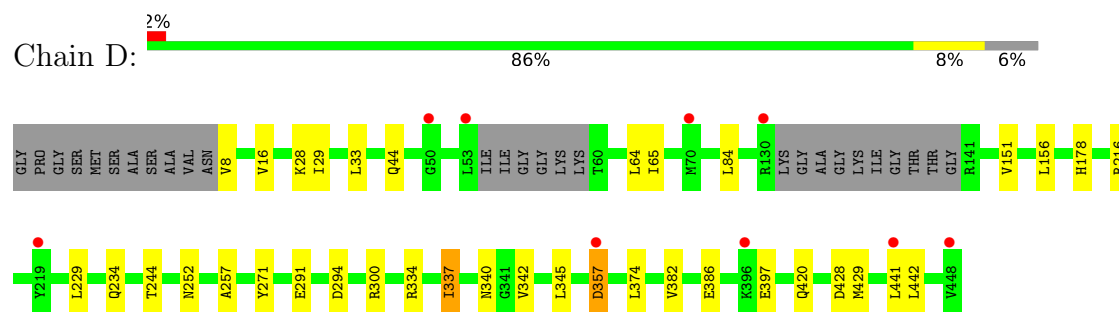
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	250	Total 250	O 250	0	0
5	C	334	Total 334	O 334	0	0
5	D	253	Total 253	O 253	0	0

- Molecule 1: Adenylosuccinate synthetase





- Molecule 1: Adenylosuccinate synthetase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	61.88Å 74.59Å 112.66Å 109.35° 90.01° 103.19°	Depositor
Resolution (Å)	50.00 – 1.95 50.00 – 1.95	Depositor EDS
% Data completeness (in resolution range)	96.2 (50.00-1.95) 96.2 (50.00-1.95)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.92 (at 1.95Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.201 , 0.247 0.207 , 0.254	Depositor DCC
R_{free} test set	6527 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	16.2	Xtriage
Anisotropy	0.110	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 42.6	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	0.93	EDS
Total number of atoms	13889	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 14.21% 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: ACT, NA, CA

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	1.07	1/3279 (0.0%)	0.90	1/4454 (0.0%)
1	B	1.00	2/3230 (0.1%)	0.87	0/4388
1	C	1.09	3/3272 (0.1%)	0.92	0/4450
1	D	1.01	2/3246 (0.1%)	0.86	2/4415 (0.0%)
All	All	1.04	8/13027 (0.1%)	0.89	3/17707 (0.0%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	427	ILE	N-CA	-5.83	1.39	1.46
1	D	257	ALA	C-O	-5.76	1.16	1.24
1	B	242	HIS	ND1-CE1	5.72	1.38	1.32
1	C	257	ALA	C-O	-5.60	1.16	1.24
1	A	242	HIS	ND1-CE1	5.45	1.38	1.32
1	C	43	PHE	C-O	5.29	1.30	1.24
1	C	18	GLY	C-O	5.25	1.29	1.24
1	D	178	HIS	CG-CD2	5.12	1.41	1.35

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	288	PHE	N-CA-CB	-5.79	103.49	111.77
1	D	65	ILE	CA-C-N	-5.09	114.48	119.92
1	D	65	ILE	C-N-CA	-5.09	114.48	119.92

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	3202	0	3114	29	0
1	B	3161	0	3063	32	0
1	C	3203	0	3126	32	0
1	D	3183	0	3089	27	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	C	4	0	3	0	0
3	D	4	0	3	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	291	0	0	3	0
5	B	250	0	0	3	0
5	C	334	0	0	0	0
5	D	253	0	0	1	0
All	All	13889	0	12398	112	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 4.

All (112) 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:151:VAL:HG11	1:D:151:VAL:CG1	1.77	1.14
1:B:151:VAL:CG1	1:D:151:VAL:HG11	1.77	1.12
1:D:29:ILE:HD12	1:D:429:MET:HE1	1.32	1.06
1:A:151[A]:VAL:HG11	1:C:151:VAL:HG11	1.03	1.03
1:A:151[A]:VAL:HG11	1:C:151:VAL:CG1	1.89	1.02
1:A:151[A]:VAL:CG1	1:C:151:VAL:HG11	1.90	1.00
1:D:29:ILE:HD12	1:D:429:MET:CE	1.93	0.96
1:B:29:ILE:HD12	1:B:429:MET:HE1	1.54	0.90
1:B:151:VAL:HG11	1:D:151:VAL:HG11	0.91	0.89
1:A:11:GLY:HA3	5:A:606:HOH:O	1.81	0.80
1:A:64:LEU:HD21	1:A:84:LEU:HD12	1.65	0.78
1:C:29:ILE:HD12	1:C:429:MET:CE	2.15	0.76
1:D:16:VAL:HG21	1:D:33:LEU:CD1	2.17	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:29:ILE:HD12	1:C:429:MET:HE1	1.69	0.73
1:C:16:VAL:HG21	1:C:33:LEU:CD1	2.18	0.73
1:D:16:VAL:HG21	1:D:33:LEU:HD13	1.71	0.71
1:C:16:VAL:HG21	1:C:33:LEU:HD13	1.72	0.70
1:B:16:VAL:HG21	1:B:33:LEU:CD1	2.23	0.69
1:C:283:VAL:HG22	1:C:323:ARG:HG2	1.76	0.66
1:A:16:VAL:HG23	1:A:229:LEU:HD11	1.78	0.65
1:B:220:GLU:OE2	5:B:574:HOH:O	2.14	0.65
1:D:64:LEU:HD21	1:D:84:LEU:HD12	1.77	0.64
1:C:16:VAL:HG23	1:C:229:LEU:HD11	1.81	0.63
1:B:29:ILE:HD12	1:B:429:MET:CE	2.25	0.62
1:D:16:VAL:HG23	1:D:229:LEU:HD11	1.82	0.62
1:B:283:VAL:HG12	5:B:1005:HOH:O	2.00	0.60
1:C:272:ILE:HD13	1:C:341:GLY:O	2.03	0.59
1:A:151[B]:VAL:HG23	5:A:473:HOH:O	2.03	0.58
1:A:395:TRP:C	1:A:396:LYS:HD3	2.28	0.58
1:D:16:VAL:CG2	1:D:33:LEU:HD13	2.33	0.58
1:A:337:ILE:HD12	1:A:342:VAL:HB	1.86	0.57
1:C:16:VAL:CG2	1:C:229:LEU:HD11	2.34	0.57
1:B:16:VAL:CG2	1:B:229:LEU:HD11	2.35	0.56
1:D:337:ILE:HD11	1:D:345:LEU:HD21	1.88	0.56
1:C:16:VAL:CG2	1:C:33:LEU:HD13	2.36	0.56
1:B:373:ASP:OD1	1:D:216:ARG:NH1	2.39	0.55
1:D:271:TYR:OH	1:D:428:ASP:OD2	2.23	0.55
1:D:16:VAL:CG2	1:D:229:LEU:HD11	2.38	0.54
1:B:16:VAL:HG23	1:B:229:LEU:HD11	1.90	0.54
1:A:38:GLN:HG2	1:A:226[A]:ARG:HD2	1.90	0.53
1:C:64:LEU:HD21	1:C:84:LEU:HD12	1.90	0.53
1:A:283:VAL:HG22	1:A:323:ARG:HG2	1.90	0.53
1:A:16:VAL:CG2	1:A:229:LEU:HD11	2.39	0.52
1:A:346:CYS:SG	1:A:429:MET:HE2	2.49	0.52
1:C:29:ILE:HD11	1:C:431:SER:HB2	1.91	0.51
1:B:31:ASP:HA	1:B:70:MET:HE2	1.93	0.51
1:B:49:ALA:HB3	1:B:63:ARG:O	2.11	0.51
1:A:95:GLU:OE2	5:A:852:HOH:O	2.19	0.51
1:D:28:LYS:HD3	1:D:28:LYS:C	2.37	0.50
1:D:337:ILE:HD12	1:D:342:VAL:HB	1.94	0.50
1:D:291:GLU:OE1	1:D:300[B]:ARG:NH1	2.45	0.50
1:B:28:LYS:NZ	1:B:438:GLU:O	2.45	0.49
1:C:334:ARG:HG2	1:C:374:LEU:HD21	1.93	0.49
1:A:28:LYS:NZ	1:A:438:GLU:O	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:346:CYS:SG	1:C:429:MET:HE2	2.54	0.48
1:B:64:LEU:HD21	1:B:84:LEU:HD12	1.96	0.48
1:D:29:ILE:CD1	1:D:429:MET:CE	2.79	0.48
1:D:16:VAL:CG2	1:D:33:LEU:CD1	2.89	0.48
1:B:16:VAL:HG21	1:B:33:LEU:HD13	1.93	0.47
1:C:105:GLU:O	1:C:105:GLU:HG3	2.14	0.47
1:C:275:ILE:HD13	1:C:346:CYS:HB3	1.95	0.47
1:B:43:PHE:CD2	1:B:258:ALA:HA	2.50	0.47
1:B:337:ILE:HD11	1:B:345:LEU:HD21	1.96	0.46
1:C:16:VAL:CG2	1:C:33:LEU:CD1	2.92	0.46
1:A:16:VAL:HG21	1:A:33:LEU:HD12	1.98	0.46
1:B:334:ARG:HG2	1:B:374:LEU:HD21	1.97	0.46
1:B:38:GLN:HG2	1:B:226[A]:ARG:HD2	1.97	0.46
1:A:420:GLN:HG2	1:A:427:ILE:HG13	1.98	0.46
1:A:29:ILE:HD11	1:A:431:SER:HB2	1.98	0.46
1:A:356:LEU:O	1:A:396:LYS:NZ	2.48	0.46
1:C:271:TYR:OH	1:C:428:ASP:OD2	2.33	0.45
1:C:337:ILE:HD13	1:C:337:ILE:N	2.32	0.45
1:C:18:GLY:O	1:C:236:THR:HG23	2.17	0.44
1:D:429:MET:HG2	1:D:441:LEU:HD12	1.98	0.44
1:A:69:ILE:HD11	1:A:103:VAL:HG13	1.99	0.44
1:D:357:ASP:HB2	5:D:511:HOH:O	2.16	0.44
1:D:244:THR:HG22	1:D:382:VAL:HG21	1.99	0.43
1:A:221[A]:GLU:OE1	1:A:226[A]:ARG:NE	2.50	0.43
1:B:31:ASP:CB	1:B:70:MET:HE2	2.48	0.43
1:B:399:THR:HA	1:B:402:ILE:HD12	2.01	0.43
1:D:291:GLU:OE2	1:D:294:ASP:HB2	2.19	0.43
1:A:216:ARG:HG2	1:A:220[B]:GLU:OE2	2.18	0.42
1:C:34:THR:HG22	1:C:71:ARG:HD3	2.00	0.42
1:C:39:GLY:HA2	1:C:75:ALA:O	2.20	0.42
1:B:443[B]:ARG:NH1	1:B:447:LYS:O	2.51	0.42
1:D:340:ASN:HB2	1:D:342:VAL:HG23	2.02	0.42
1:B:271:TYR:OH	1:B:428:ASP:OD2	2.36	0.42
1:A:334:ARG:HG2	1:A:374:LEU:CD2	2.50	0.41
1:C:44[A]:GLN:HG3	1:C:252:ASN:O	2.20	0.41
1:C:354:ASP:O	1:C:396:LYS:HA	2.20	0.41
1:A:197:LEU:HD11	1:A:201:LEU:HD11	2.01	0.41
1:B:16:VAL:HG21	1:B:229:LEU:HD11	2.02	0.41
1:B:16:VAL:CG2	1:B:33:LEU:CD1	2.95	0.41
1:D:334:ARG:HG2	1:D:374:LEU:HD21	2.02	0.41
1:A:153:ARG:HA	1:C:245:TYR:OH	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:44:GLN:HG3	1:D:252:ASN:O	2.20	0.41
1:C:44[B]:GLN:HE22	1:C:233:ALA:HB3	1.85	0.41
1:C:333:LEU:O	1:C:337:ILE:HD13	2.20	0.41
1:A:337:ILE:HD12	1:A:342:VAL:CB	2.51	0.41
1:B:34:THR:HG21	1:B:68:GLY:HA2	2.03	0.41
1:B:291:GLU:OE2	1:B:294:ASP:HB2	2.21	0.41
1:A:355:GLY:HA2	1:A:396:LYS:HD2	2.02	0.41
1:A:395:TRP:O	1:A:396:LYS:HD3	2.20	0.41
1:B:153:ARG:CZ	5:B:492:HOH:O	2.68	0.41
1:C:28:LYS:NZ	1:C:438:GLU:O	2.54	0.41
1:C:35:ASP:HA	1:C:71:ARG:CZ	2.51	0.41
1:C:281:THR:HA	1:C:324:ARG:O	2.22	0.40
1:A:111[B]:GLU:HB2	1:A:159:GLN:HB3	2.02	0.40
1:B:48:ASN:C	1:B:49:ALA:O	2.63	0.40
1:B:83:VAL:HG12	1:B:114:THR:HB	2.02	0.40
1:B:16:VAL:CG2	1:B:33:LEU:HD13	2.51	0.40
1:D:156:LEU:HD23	1:D:156:LEU:HA	1.97	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	422/452 (93%)	415 (98%)	7 (2%)	0	100	100
1	B	418/452 (92%)	407 (97%)	9 (2%)	2 (0%)	24	16
1	C	423/452 (94%)	417 (99%)	6 (1%)	0	100	100
1	D	420/452 (93%)	406 (97%)	14 (3%)	0	100	100
All	All	1683/1808 (93%)	1645 (98%)	36 (2%)	2 (0%)	48	41

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	49	ALA
1	B	50	GLY

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	322/352 (92%)	320 (99%)	2 (1%)	78	79
1	B	313/352 (89%)	310 (99%)	3 (1%)	68	67
1	C	324/352 (92%)	316 (98%)	8 (2%)	42	34
1	D	320/352 (91%)	312 (98%)	8 (2%)	42	34
All	All	1279/1408 (91%)	1258 (98%)	21 (2%)	53	52

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	LYS
1	A	442	LEU
1	B	70	MET
1	B	357	ASP
1	B	442	LEU
1	C	61	ILE
1	C	105	GLU
1	C	337	ILE
1	C	377	ARG
1	C	386	GLU
1	C	442	LEU
1	C	443	ARG
1	C	448	VAL
1	D	8	VAL
1	D	234	GLN
1	D	337	ILE
1	D	357	ASP
1	D	386	GLU
1	D	397	GLU
1	D	420	GLN

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Mol	Chain	Res	Type
1	D	442	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	227	ASN
1	A	338	GLN
1	B	303	GLN
1	B	338	GLN
1	D	338	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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 for Mogul analysis.

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	$\# Z > 2$
3	ACT	C	449	-	3,3,3	0.81	0	3,3,3	0.34	0
3	ACT	D	449	-	3,3,3	0.87	0	3,3,3	0.55	0

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.

No monomer is involved in short contacts.

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 ⓘ

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	424/452 (93%)	0.07	14 (3%)	49	55	6, 16, 30, 43	6 (1%)
1	B	421/452 (93%)	0.17	14 (3%)	49	55	9, 17, 34, 44	3 (0%)
1	C	426/452 (94%)	0.09	6 (1%)	73	79	5, 17, 30, 36	3 (0%)
1	D	425/452 (94%)	0.30	9 (2%)	63	70	9, 20, 35, 50	1 (0%)
All	All	1696/1808 (93%)	0.16	43 (2%)	58	65	5, 18, 33, 50	13 (0%)

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	49	ALA	4.5
1	D	53	LEU	4.1
1	B	318	VAL	3.8
1	B	188	GLY	3.3
1	B	319	THR	3.3
1	A	52	THR	3.2
1	A	10	PRO	3.2
1	A	57	GLY	3.1
1	B	52	THR	3.0
1	B	10	PRO	3.0
1	D	448	VAL	2.9
1	A	321	ARG	2.9
1	D	396	LYS	2.8
1	A	448	VAL	2.8
1	B	50	GLY	2.7
1	D	441	LEU	2.7
1	A	187	GLY	2.7
1	A	61	ILE	2.6
1	D	219	TYR	2.6
1	A	70	MET	2.6
1	B	61	ILE	2.6

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Mol	Chain	Res	Type	RSRZ
1	D	70	MET	2.5
1	D	130	ARG	2.5
1	C	140	GLY	2.5
1	A	284	GLY	2.4
1	C	357	ASP	2.4
1	A	51	HIS	2.4
1	B	317	SER	2.4
1	D	357	ASP	2.3
1	B	70	MET	2.3
1	C	219	TYR	2.3
1	A	396	LYS	2.3
1	A	67	SER	2.2
1	A	355	GLY	2.2
1	B	128	GLU	2.1
1	B	129	ALA	2.1
1	C	71	ARG	2.1
1	C	448	VAL	2.1
1	D	50	GLY	2.1
1	B	48	ASN	2.1
1	B	233	ALA	2.1
1	C	100	GLY	2.0
1	A	48	ASN	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(Å ²)	Q<0.9
3	ACT	C	449	4/4	0.77	0.15	31,33,33,34	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	ACT	D	449	4/4	0.80	0.14	28,29,29,32	0
2	NA	A	449	1/1	0.94	0.12	23,23,23,23	0
2	NA	B	449	1/1	0.98	0.05	21,21,21,21	0
4	CA	D	450	1/1	0.98	0.10	27,27,27,27	0
4	CA	C	450	1/1	0.99	0.09	23,23,23,23	0

6.5 Other polymers [i](#)

There are no such residues in this entry.