



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:37 PM UTC

PDB ID : 1MSV / pdb_00001msv
Title : The S68A S-adenosylmethionine decarboxylase proenzyme processing mutant.
Authors : Tolbert, W.D.; Zhang, Y.; Bennett, E.M.; Cottet, S.E.; Ekstrom, J.L.; Pegg, A.E.; Ealick, S.E.
Deposited on : 2002-09-19
Resolution : 1.75 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

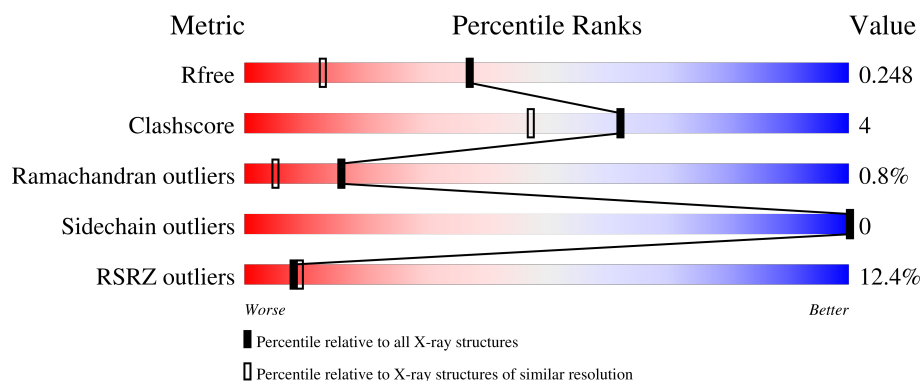
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.75 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	3183 (1.76-1.76)
Clashscore	190562	3299 (1.76-1.76)
Ramachandran outliers	187476	3274 (1.76-1.76)
Sidechain outliers	187428	3274 (1.76-1.76)
RSRZ outliers	180081	3183 (1.76-1.76)

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	354	14% 79% 11% 10%
1	B	354	8% 77% 12% 10%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5762 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 S-adenosylmethionine decarboxylase proenzyme.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	319	Total	C	N	O	S	0	14	0
			2683	1715	443	505	20			
1	B	320	Total	C	N	O	S	0	21	0
			2745	1753	453	518	21			

There are 42 discrepancies between the modelled and reference sequences:

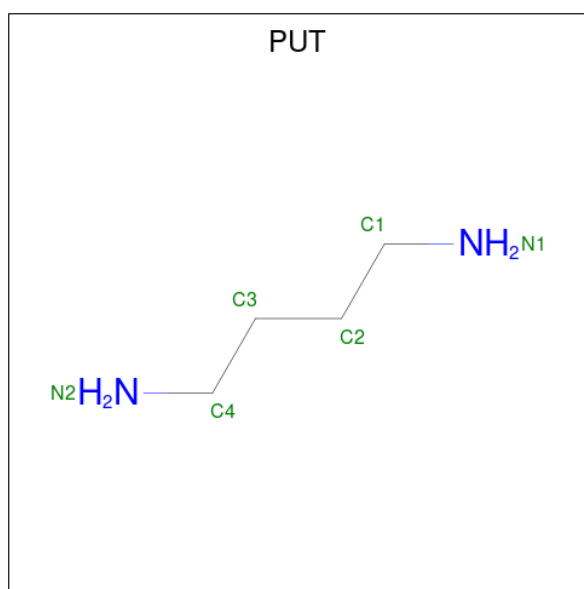
Chain	Residue	Modelled	Actual	Comment	Reference
A	68	ALA	SER	engineered mutation	UNP P17707
A	335	HIS	-	cloning artifact	UNP P17707
A	336	ILE	-	cloning artifact	UNP P17707
A	337	CYS	-	cloning artifact	UNP P17707
A	338	ARG	-	cloning artifact	UNP P17707
A	339	SER	-	cloning artifact	UNP P17707
A	330	GLN	-	cloning artifact	UNP P17707
A	341	MET	-	cloning artifact	UNP P17707
A	342	VAL	-	cloning artifact	UNP P17707
A	343	THR	-	cloning artifact	UNP P17707
A	344	SER	-	cloning artifact	UNP P17707
A	345	GLN	-	cloning artifact	UNP P17707
A	346	GLN	-	cloning artifact	UNP P17707
A	347	THR	-	cloning artifact	UNP P17707
A	348	SER	-	cloning artifact	UNP P17707
A	349	SER	-	cloning artifact	UNP P17707
A	350	VAL	-	cloning artifact	UNP P17707
A	351	VAL	-	cloning artifact	UNP P17707
A	352	ARG	-	cloning artifact	UNP P17707
A	353	GLN	-	cloning artifact	UNP P17707
A	354	THR	-	cloning artifact	UNP P17707
B	68	ALA	SER	engineered mutation	UNP P17707
B	335	HIS	-	cloning artifact	UNP P17707
B	336	ILE	-	cloning artifact	UNP P17707
B	337	CYS	-	cloning artifact	UNP P17707

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Chain	Residue	Modelled	Actual	Comment	Reference
B	338	ARG	-	cloning artifact	UNP P17707
B	339	SER	-	cloning artifact	UNP P17707
B	340	GLN	-	cloning artifact	UNP P17707
B	341	MET	-	cloning artifact	UNP P17707
B	342	VAL	-	cloning artifact	UNP P17707
B	343	THR	-	cloning artifact	UNP P17707
B	344	SER	-	cloning artifact	UNP P17707
B	345	GLN	-	cloning artifact	UNP P17707
B	346	GLN	-	cloning artifact	UNP P17707
B	347	THR	-	cloning artifact	UNP P17707
B	348	SER	-	cloning artifact	UNP P17707
B	349	SER	-	cloning artifact	UNP P17707
B	350	VAL	-	cloning artifact	UNP P17707
B	351	VAL	-	cloning artifact	UNP P17707
B	352	ARG	-	cloning artifact	UNP P17707
B	353	GLN	-	cloning artifact	UNP P17707
B	354	THR	-	cloning artifact	UNP P17707

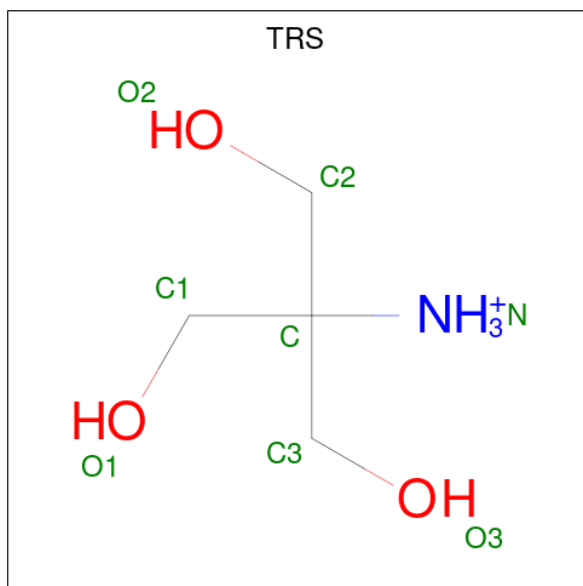
- Molecule 2 is 1,4-DIAMINOBUTANE (CCD ID: PUT) (formula: C₄H₁₂N₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	N	0	0
			6	4	2		
2	B	1	Total	C	N	0	0
			6	4	2		

- Molecule 3 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (CCD ID: TRS)

(formula: $C_4H_{12}NO_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			8	4	1	3		
3	B	1	Total	C	N	O	0	0
			8	4	1	3		

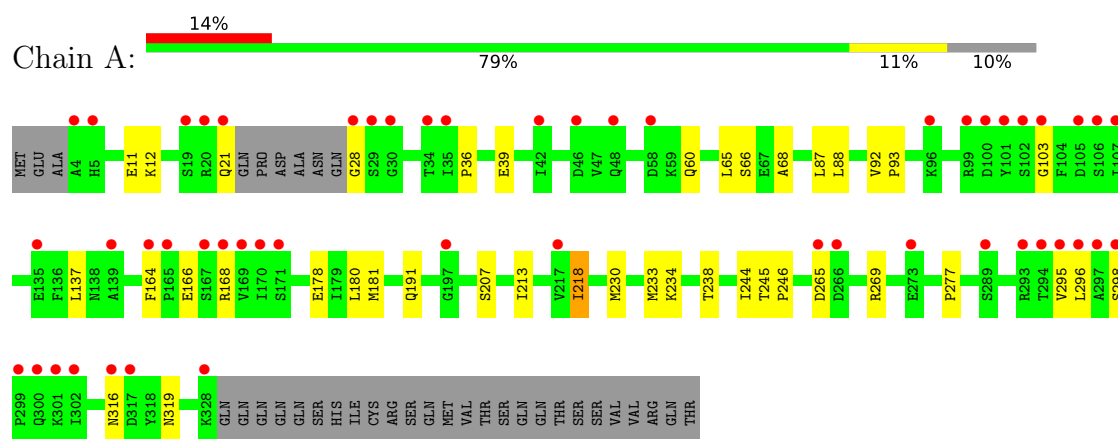
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	128	Total	O	0	0
			128	128		
4	B	178	Total	O	0	0
			178	178		

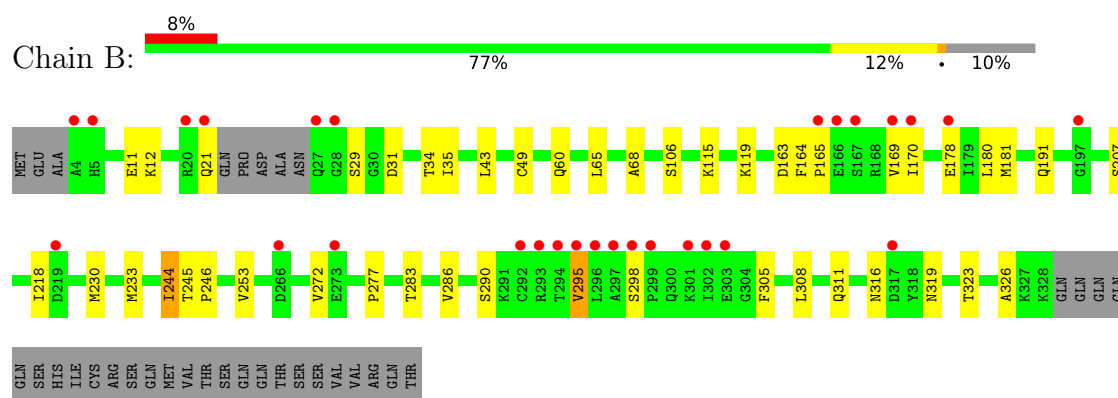
3 Residue-property plots [i](#)

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: S-adenosylmethionine decarboxylase proenzyme



- Molecule 1: S-adenosylmethionine decarboxylase proenzyme



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	73.94Å 56.38Å 99.19Å 90.00° 110.97° 90.00°	Depositor
Resolution (Å)	16.37 – 1.75 16.37 – 1.75	Depositor EDS
% Data completeness (in resolution range)	98.9 (16.37-1.75) 98.7 (16.37-1.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.76 (at 1.70Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.210 , 0.238 0.225 , 0.248	Depositor DCC
R_{free} test set	7623 reflections (9.20%)	wwPDB-VP
Wilson B-factor (Å ²)	22.0	Xtriage
Anisotropy	0.053	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.45 , 52.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.012 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5762	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.05% 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

5.1 Standard geometry

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

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	0.37	0/2746	0.86	10/3707 (0.3%)
1	B	0.38	0/2807	0.88	12/3789 (0.3%)
All	All	0.37	0/5553	0.87	22/7496 (0.3%)

There are no bond length outliers.

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	246	PRO	N-CA-C	8.57	124.58	113.86
1	B	246	PRO	N-CA-C	8.50	124.49	113.86
1	A	319	ASN	N-CA-C	-6.48	100.02	110.32
1	A	11	GLU	N-CA-C	6.36	120.27	110.42
1	B	12	LYS	N-CA-C	-6.34	98.91	109.24
1	B	11	GLU	N-CA-C	6.31	120.20	110.42
1	A	12	LYS	N-CA-C	-5.90	99.56	109.07
1	B	319	ASN	N-CA-C	-5.80	100.83	110.17
1	A	213	ILE	CA-C-N	5.66	125.66	119.89
1	A	213	ILE	C-N-CA	5.66	125.66	119.89
1	A	218	ILE	N-CA-C	5.51	116.53	108.53
1	A	245	THR	CA-C-N	5.50	125.11	119.56
1	A	245	THR	C-N-CA	5.50	125.11	119.56
1	A	244	ILE	N-CA-C	5.43	115.71	108.11
1	B	35	ILE	CA-C-N	5.37	125.31	119.78
1	B	35	ILE	C-N-CA	5.37	125.31	119.78
1	B	119	LYS	CA-C-N	5.31	125.37	119.32
1	B	119	LYS	C-N-CA	5.31	125.37	119.32
1	B	245	THR	CA-C-N	5.25	124.87	119.56
1	B	245	THR	C-N-CA	5.25	124.87	119.56
1	B	244	ILE	N-CA-C	5.25	115.46	108.11
1	B	115	LYS	N-CA-C	-5.04	103.75	110.55

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	2683	0	2619	19	0
1	B	2745	0	2683	25	0
2	A	6	0	12	0	0
2	B	6	0	12	0	0
3	A	8	0	11	0	0
3	B	8	0	11	0	0
4	A	128	0	0	1	0
4	B	178	0	0	2	0
All	All	5762	0	5348	44	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 (44) 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:164:PHE:H	1:A:316:ASN:HD21	1.31	0.78
1:B:164:PHE:H	1:B:316:ASN:HD21	1.30	0.76
1:B:164:PHE:H	1:B:316:ASN:ND2	1.93	0.67
1:A:164:PHE:H	1:A:316:ASN:ND2	1.98	0.61
1:A:295:VAL:HB	1:A:296:LEU:HD12	1.86	0.57
1:A:36:PRO:HG2	1:A:39:GLU:HG2	1.85	0.57
1:B:65:LEU:HB2	1:B:68:ALA:HB3	1.86	0.56
1:B:308:LEU:HD11	1:B:326:ALA:HB2	1.89	0.55
1:B:164:PHE:N	1:B:316:ASN:HD21	2.03	0.55
1:B:21:GLN:HB2	1:B:106:SER:HB2	1.89	0.55
1:B:272[A]:VAL:HG11	1:B:305:PHE:CE1	2.43	0.54
1:A:92:VAL:HB	1:A:93:PRO:HD3	1.90	0.53
1:B:191:GLN:HE22	1:B:207:SER:HA	1.75	0.51
1:B:181:MET:HB3	1:B:277:PRO:HB3	1.92	0.51
1:B:286[B]:VAL:HG13	1:B:290:SER:HB3	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:66:SER:HB3	4:A:875:HOH:O	2.11	0.50
1:A:181:MET:HB3	1:A:277:PRO:HB3	1.94	0.49
1:A:21:GLN:HE21	1:A:168:ARG:HH22	1.61	0.48
1:B:218:ILE:HG23	1:B:230:MET:HE2	1.96	0.47
1:A:234:LYS:HB2	1:A:238:THR:HB	1.96	0.47
1:B:49[B]:CYS:SG	1:B:65:LEU:HD13	2.55	0.47
1:B:43[B]:LEU:HD12	4:B:686:HOH:O	2.15	0.46
1:A:191:GLN:HE22	1:A:207:SER:HA	1.81	0.46
1:B:178:GLU:HG2	1:B:180:LEU:HD21	1.98	0.45
1:A:88:LEU:HD22	1:A:137:LEU:HD21	1.99	0.45
1:A:65:LEU:HB2	1:A:68:ALA:HB3	1.98	0.45
1:B:286[A]:VAL:HG21	1:B:295:VAL:HG21	1.98	0.44
1:B:163:ASP:HA	1:B:316:ASN:HD21	1.83	0.44
1:B:286[A]:VAL:CG2	1:B:295:VAL:HG21	2.48	0.43
1:B:178:GLU:HB3	1:B:283:THR:OG1	2.19	0.43
1:A:28:GLY:HA3	1:A:103:GLY:HA3	2.01	0.42
1:A:60:GLN:HG3	1:A:233:MET:SD	2.59	0.42
1:B:29:SER:CB	1:B:34[B]:THR:HG21	2.49	0.42
1:A:218:ILE:HG23	1:A:230:MET:HE2	2.02	0.42
1:B:31:ASP:O	1:B:34[A]:THR:HG22	2.19	0.42
1:A:178:GLU:HG2	1:A:180:LEU:HD21	2.01	0.42
1:B:244:ILE:HG23	1:B:253:VAL:HG22	2.02	0.42
1:B:323[B]:THR:HG21	4:B:906:HOH:O	2.19	0.41
1:A:296:LEU:HD12	1:A:296:LEU:N	2.36	0.41
1:B:169:VAL:HG22	1:B:170:ILE:N	2.36	0.40
1:A:87:LEU:HD23	1:A:87:LEU:C	2.47	0.40
1:A:265:ASP:O	1:A:269:ARG:HG3	2.21	0.40
1:B:60:GLN:HG3	1:B:233:MET:SD	2.62	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	329/354 (93%)	317 (96%)	10 (3%)	2 (1%)	21	8
1	B	337/354 (95%)	326 (97%)	8 (2%)	3 (1%)	14	4
All	All	666/708 (94%)	643 (96%)	18 (3%)	5 (1%)	16	5

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	165	PRO
1	A	166	GLU
1	B	298	SER
1	A	298	SER
1	B	295	VAL

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	303/322 (94%)	303 (100%)	0	100	100
1	B	311/322 (97%)	311 (100%)	0	100	100
All	All	614/644 (95%)	614 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	GLN
1	A	191	GLN
1	A	243	HIS
1	A	316	ASN
1	B	191	GLN
1	B	316	ASN

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 ⓘ

4 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	PUT	B	355	-	5,5,5	0.76	0	4,4,4	0.31	0
2	PUT	A	355	-	5,5,5	0.73	0	4,4,4	0.30	0
3	TRS	B	501	-	7,7,7	2.02	2 (28%)	9,9,9	2.17	3 (33%)
3	TRS	A	500	-	7,7,7	2.03	2 (28%)	9,9,9	2.19	3 (33%)

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	PUT	B	355	-	-	0/3/3/3	-
2	PUT	A	355	-	-	0/3/3/3	-
3	TRS	B	501	-	-	0/9/9/9	-
3	TRS	A	500	-	-	0/9/9/9	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	501	TRS	C2-C	-4.63	1.40	1.53
3	A	500	TRS	C2-C	-4.52	1.41	1.53
3	A	500	TRS	O1-C1	-2.75	1.33	1.42
3	B	501	TRS	O1-C1	-2.52	1.34	1.42

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	500	TRS	O3-C3-C	4.60	123.69	110.88
3	B	501	TRS	O3-C3-C	4.49	123.39	110.88
3	B	501	TRS	O2-C2-C	2.60	118.12	110.88
3	A	500	TRS	O2-C2-C	2.57	118.05	110.88
3	B	501	TRS	C2-C-C1	-2.27	104.61	110.66
3	A	500	TRS	C2-C-C1	-2.26	104.63	110.66

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	319/354 (90%)	0.81	51 (15%) 5 5	7, 25, 70, 98	14 (4%)
1	B	320/354 (90%)	0.37	28 (8%) 15 19	7, 20, 50, 84	21 (6%)
All	All	639/708 (90%)	0.59	79 (12%) 8 9	7, 23, 55, 98	35 (5%)

All (79) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	298	SER	7.2
1	A	294	THR	6.8
1	A	300[A]	GLN	6.6
1	A	296	LEU	6.2
1	A	299	PRO	6.1
1	B	294	THR	5.8
1	A	298	SER	5.4
1	A	297	ALA	4.9
1	A	165	PRO	4.7
1	A	295	VAL	4.7
1	B	295	VAL	4.4
1	B	165	PRO	4.2
1	A	4	ALA	4.1
1	A	169	VAL	4.1
1	B	293[A]	ARG	4.1
1	A	170	ILE	3.9
1	A	19	SER	3.9
1	A	167	SER	3.9
1	B	299	PRO	3.9
1	A	5[A]	HIS	3.6
1	A	317	ASP	3.6
1	B	167	SER	3.5
1	A	20	ARG	3.5
1	A	316	ASN	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	296	LEU	3.4
1	A	28	GLY	3.3
1	A	301	LYS	3.3
1	B	292	CYS	3.3
1	A	289	SER	3.2
1	A	21	GLN	3.2
1	B	266	ASP	3.2
1	B	297	ALA	3.2
1	A	164	PHE	3.1
1	B	21	GLN	3.1
1	A	293	ARG	3.1
1	A	106	SER	3.1
1	A	103	GLY	3.0
1	B	197	GLY	3.0
1	B	301	LYS	2.9
1	B	170	ILE	2.9
1	A	99[A]	ARG	2.8
1	B	273	GLU	2.8
1	B	5	HIS	2.8
1	B	20	ARG	2.8
1	B	166	GLU	2.8
1	A	101	TYR	2.8
1	A	266	ASP	2.8
1	B	27	GLN	2.7
1	A	135	GLU	2.6
1	A	273	GLU	2.6
1	B	169	VAL	2.6
1	A	42[A]	ILE	2.6
1	A	168	ARG	2.6
1	B	317	ASP	2.5
1	B	178	GLU	2.5
1	A	105	ASP	2.4
1	A	265	ASP	2.4
1	A	58	ASP	2.4
1	A	96	LYS	2.4
1	A	35	ILE	2.4
1	B	303	GLU	2.3
1	B	302	ILE	2.3
1	A	197	GLY	2.3
1	B	219	ASP	2.3
1	A	328	LYS	2.3
1	A	302	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	28	GLY	2.3
1	A	48	GLN	2.2
1	A	30	GLY	2.2
1	A	107	ILE	2.2
1	A	171	SER	2.2
1	B	4	ALA	2.1
1	A	29	SER	2.1
1	A	100	ASP	2.1
1	A	102	SER	2.1
1	A	34	THR	2.1
1	A	139	ALA	2.0
1	A	46	ASP	2.0
1	A	217[A]	VAL	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
3	TRS	A	500	8/8	0.91	0.10	24,27,32,35	0
3	TRS	B	501	8/8	0.92	0.11	23,25,26,32	0
2	PUT	A	355	6/6	0.94	0.07	20,22,25,26	0
2	PUT	B	355	6/6	0.95	0.07	15,16,18,18	0

6.5 Other polymers [i](#)

There are no such residues in this entry.