



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 12, 2024 – 07:21 PM EDT

PDB ID : 3SRN
Title : STRUCTURAL CHANGES THAT ACCOMPANY THE REDUCED CATALYTIC EFFICIENCY OF TWO SEMISYNTHETIC RIBONUCLEASE ANALOGS
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Deposited on : 1991-05-20
Resolution : 2.00 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	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

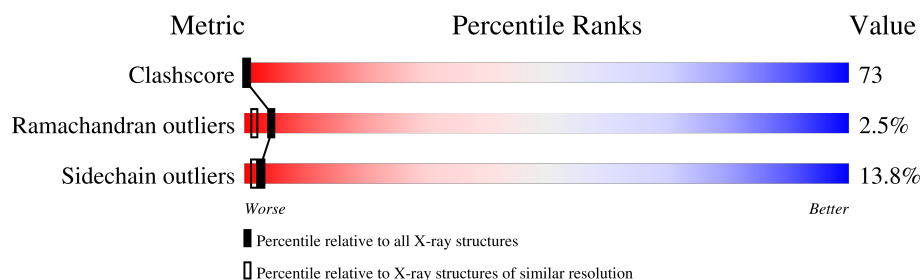
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 Å.


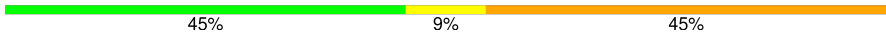
Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	180529	10737 (2.00-2.00)
Ramachandran outliers	177936	10628 (2.00-2.00)
Sidechain outliers	177891	10627 (2.00-2.00)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	113	 35% 44% 18% •
2	B	11	 45% 9% 45%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 1089 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 RIBONUCLEASE A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	113	863	516	158	177	12	0	0	0

- Molecule 2 is a protein called RIBONUCLEASE A.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	B	11	88	59	14	15	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	121	ASN	-	insertion	UNP P61823

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is water.

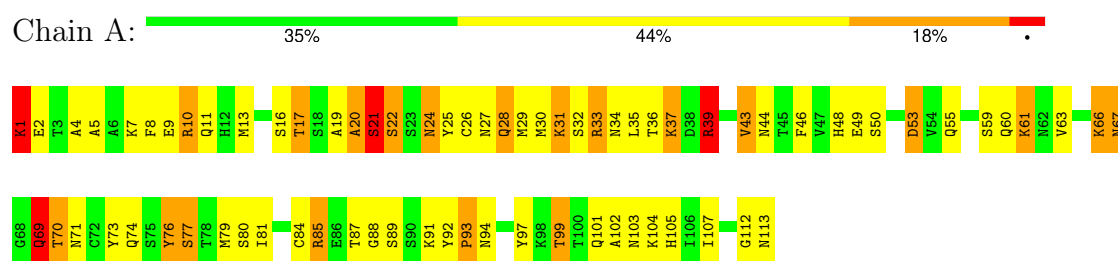
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	124	Total	O	0	0
			124	124		
4	B	9	Total	O	0	0
			9	9		

3 Residue-property plots [i](#)

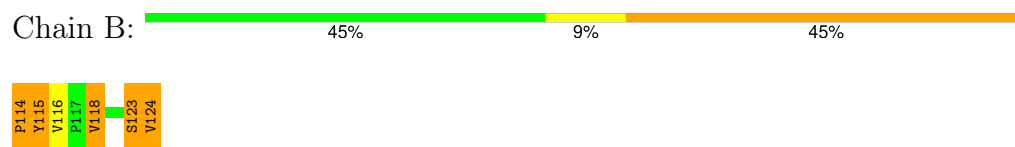
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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: RIBONUCLEASE A



• Molecule 2: RIBONUCLEASE A



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	64.70Å 64.70Å 64.90Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	5.00 – 2.00	Depositor
% Data completeness (in resolution range)	(Not available) (5.00-2.00)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROLSQ, X-PLOR	Depositor
R, R_{free}	0.186 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1089	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SO4

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	1.42	4/874 (0.5%)	2.64	29/1175 (2.5%)
2	B	1.30	0/92	2.31	6/125 (4.8%)
All	All	1.40	4/966 (0.4%)	2.61	35/1300 (2.7%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1	LYS	C-N	-16.62	0.95	1.34
1	A	39	ARG	CD-NE	-6.43	1.35	1.46
1	A	77	SER	CB-OG	-6.32	1.34	1.42
1	A	112	GLY	N-CA	5.68	1.54	1.46

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	39	ARG	CD-NE-CZ	38.68	177.75	123.60
1	A	1	LYS	O-C-N	-31.92	71.63	122.70
1	A	33	ARG	NE-CZ-NH1	27.17	133.88	120.30
1	A	1	LYS	CA-C-N	21.60	164.73	117.20
1	A	33	ARG	NE-CZ-NH2	-11.01	114.80	120.30
1	A	10	ARG	NE-CZ-NH2	10.52	125.56	120.30
1	A	61	LYS	CA-CB-CG	9.54	134.38	113.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	115	TYR	CA-CB-CG	9.48	131.42	113.40
2	B	123	SER	C-N-CA	9.30	144.96	121.70
1	A	85	ARG	NE-CZ-NH1	8.62	124.61	120.30
1	A	85	ARG	CG-CD-NE	8.03	128.65	111.80
1	A	33	ARG	NH1-CZ-NH2	-7.46	111.19	119.40
1	A	21	SER	C-N-CA	7.39	140.19	121.70
1	A	30	MET	CA-CB-CG	7.32	125.74	113.30
1	A	61	LYS	CB-CG-CD	7.22	130.36	111.60
1	A	39	ARG	NE-CZ-NH1	7.15	123.88	120.30
1	A	85	ARG	CD-NE-CZ	6.97	133.35	123.60
1	A	20	ALA	N-CA-CB	6.37	119.02	110.10
2	B	115	TYR	N-CA-C	-6.14	94.43	111.00
1	A	1	LYS	N-CA-C	-5.91	95.05	111.00
1	A	69	GLN	N-CA-CB	5.87	121.16	110.60
1	A	2	GLU	OE1-CD-OE2	5.85	130.31	123.30
1	A	2	GLU	CG-CD-OE2	-5.53	107.25	118.30
1	A	10	ARG	NE-CZ-NH1	-5.51	117.55	120.30
1	A	53	ASP	CB-CG-OD2	-5.48	113.37	118.30
1	A	43	VAL	O-C-N	5.44	131.41	122.70
1	A	39	ARG	CG-CD-NE	5.41	123.16	111.80
1	A	20	ALA	CA-C-N	-5.32	105.50	117.20
2	B	114	PRO	CA-C-O	5.31	132.95	120.20
1	A	31	LYS	C-N-CA	5.28	134.90	121.70
2	B	114	PRO	CA-C-N	-5.28	105.59	117.20
2	B	114	PRO	CA-N-CD	-5.27	104.12	111.50
1	A	112	GLY	N-CA-C	-5.18	100.15	113.10
1	A	76	TYR	CB-CG-CD1	-5.17	117.89	121.00
1	A	22	SER	N-CA-CB	-5.05	102.92	110.50

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1	LYS	Mainchain

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	863	0	823	125	0
2	B	88	0	81	22	0
3	B	5	0	0	0	0
4	A	124	0	0	65	0
4	B	9	0	0	9	0
All	All	1089	0	904	135	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 73.

All (135) 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:67:ASN:HD21	1:A:69:GLN:NE2	1.29	1.28
1:A:49:GLU:HB3	4:A:335:HOH:O	1.09	1.25
1:A:104:LYS:HB3	4:B:216:HOH:O	1.34	1.21
1:A:67:ASN:ND2	1:A:69:GLN:HE22	1.39	1.20
1:A:10:ARG:HD2	4:A:277:HOH:O	1.44	1.17
1:A:79:MET:HG3	4:A:172:HOH:O	1.45	1.16
1:A:101:GLN:HB2	4:A:229:HOH:O	1.46	1.14
1:A:16:SER:HB3	4:A:167:HOH:O	1.46	1.13
1:A:105:HIS:H	2:B:124:VAL:HG23	1.18	1.09
1:A:101:GLN:OE1	4:A:229:HOH:O	1.73	1.06
1:A:102:ALA:HB3	4:A:222:HOH:O	1.55	1.02
1:A:7:LYS:HE2	4:A:145:HOH:O	1.60	1.01
1:A:80:SER:O	4:A:288:HOH:O	1.81	0.98
2:B:116:VAL:HG22	4:B:346:HOH:O	1.67	0.95
1:A:80:SER:C	4:A:288:HOH:O	2.05	0.95
1:A:104:LYS:HD2	4:A:156:HOH:O	1.67	0.92
1:A:21:SER:OG	1:A:22:SER:N	2.01	0.88
1:A:55:GLN:HB3	4:A:196:HOH:O	1.73	0.87
1:A:28:GLN:OE1	4:A:130:HOH:O	1.92	0.86
1:A:77:SER:O	4:A:172:HOH:O	1.94	0.86
2:B:124:VAL:CG1	2:B:124:VAL:OXT	2.22	0.86
1:A:105:HIS:H	2:B:124:VAL:CG2	1.88	0.86
1:A:26:CYS:SG	1:A:99:THR:HB	2.18	0.84
1:A:105:HIS:N	2:B:124:VAL:HG23	1.95	0.81
1:A:67:ASN:N	4:A:175:HOH:O	2.12	0.80
1:A:66:LYS:HB2	1:A:66:LYS:NZ	1.97	0.79
2:B:124:VAL:OXT	2:B:124:VAL:HG12	1.82	0.79
4:A:264:HOH:O	2:B:115:TYR:CD2	2.37	0.78
1:A:17:THR:HG22	1:A:48:HIS:HD1	1.50	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:48:HIS:CD2	4:A:288:HOH:O	2.35	0.77
1:A:34:ASN:HA	1:A:37:LYS:HE3	1.65	0.77
1:A:66:LYS:C	4:A:175:HOH:O	2.22	0.77
1:A:1:LYS:O	4:A:184:HOH:O	2.02	0.76
1:A:69:GLN:HG2	4:A:173:HOH:O	1.87	0.75
1:A:70:THR:HB	4:A:233:HOH:O	1.86	0.75
1:A:20:ALA:HB3	4:A:229:HOH:O	1.87	0.74
1:A:85:ARG:NE	4:A:206:HOH:O	1.97	0.74
2:B:124:VAL:N	4:B:216:HOH:O	2.21	0.73
1:A:49:GLU:CB	4:A:335:HOH:O	1.89	0.72
1:A:20:ALA:CB	4:A:229:HOH:O	2.37	0.72
1:A:17:THR:HG22	1:A:48:HIS:ND1	2.05	0.72
1:A:48:HIS:HD2	4:A:288:HOH:O	1.70	0.71
1:A:36:THR:HA	1:A:39:ARG:O	1.89	0.71
1:A:5:ALA:O	4:A:151:HOH:O	2.08	0.71
1:A:67:ASN:ND2	1:A:69:GLN:NE2	2.13	0.71
1:A:113:ASN:O	4:A:347:HOH:O	2.08	0.71
4:A:264:HOH:O	2:B:115:TYR:CE2	2.45	0.70
1:A:70:THR:CB	4:A:233:HOH:O	2.39	0.69
1:A:34:ASN:HA	1:A:37:LYS:CE	2.23	0.68
1:A:69:GLN:NE2	4:A:173:HOH:O	2.26	0.68
1:A:63:VAL:HG12	4:A:306:HOH:O	1.93	0.67
1:A:39:ARG:HG3	4:A:276:HOH:O	1.93	0.67
1:A:4:ALA:HB1	2:B:118:VAL:HG22	1.77	0.66
1:A:71:ASN:HA	4:B:218:HOH:O	1.94	0.66
1:A:104:LYS:HD3	4:B:216:HOH:O	1.97	0.64
1:A:73:TYR:HH	2:B:115:TYR:HE2	1.45	0.64
1:A:32:SER:HB3	4:A:241:HOH:O	1.98	0.63
1:A:66:LYS:CA	4:A:175:HOH:O	2.46	0.63
1:A:27:ASN:HD21	1:A:97:TYR:H	1.45	0.63
1:A:113:ASN:HA	4:A:157:HOH:O	1.97	0.63
1:A:105:HIS:O	2:B:124:VAL:N	2.32	0.62
1:A:87:THR:O	1:A:89:SER:N	2.33	0.61
1:A:63:VAL:CG1	4:A:306:HOH:O	2.48	0.60
1:A:34:ASN:CA	1:A:37:LYS:HE3	2.30	0.60
1:A:103:ASN:ND2	4:A:213:HOH:O	2.37	0.58
1:A:20:ALA:CB	4:A:137:HOH:O	2.51	0.57
1:A:81:ILE:HD11	4:A:156:HOH:O	2.04	0.57
1:A:46:PHE:HD2	4:A:280:HOH:O	1.86	0.57
2:B:116:VAL:N	4:B:346:HOH:O	2.37	0.57
1:A:48:HIS:HB2	4:A:288:HOH:O	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:60:GLN:HG2	1:A:76:TYR:CE2	2.40	0.56
1:A:37:LYS:N	1:A:37:LYS:HE2	2.20	0.56
1:A:81:ILE:HD11	4:A:222:HOH:O	2.04	0.56
1:A:26:CYS:HB2	4:A:131:HOH:O	2.06	0.56
1:A:33:ARG:O	4:A:277:HOH:O	2.18	0.55
1:A:73:TYR:CE2	2:B:115:TYR:HE2	2.25	0.54
1:A:61:LYS:O	1:A:73:TYR:HA	2.08	0.54
1:A:26:CYS:H	1:A:99:THR:HG21	1.73	0.53
2:B:123:SER:C	4:B:216:HOH:O	2.47	0.53
1:A:34:ASN:HD22	1:A:37:LYS:HE3	1.74	0.52
1:A:66:LYS:HB2	1:A:66:LYS:HZ3	1.72	0.52
1:A:105:HIS:HB2	2:B:124:VAL:HA	1.91	0.52
2:B:124:VAL:OXT	2:B:124:VAL:HG13	2.09	0.50
1:A:17:THR:O	1:A:48:HIS:HB3	2.11	0.50
1:A:66:LYS:HB2	1:A:66:LYS:HZ2	1.74	0.50
1:A:21:SER:OG	1:A:22:SER:OG	2.24	0.50
1:A:26:CYS:HB2	1:A:99:THR:HG22	1.93	0.50
1:A:17:THR:HG23	1:A:19:ALA:O	2.12	0.50
1:A:66:LYS:NZ	1:A:66:LYS:CB	2.71	0.50
1:A:25:TYR:HE1	4:A:280:HOH:O	1.94	0.49
1:A:92:TYR:CD1	1:A:93:PRO:HA	2.47	0.49
1:A:66:LYS:HA	4:A:175:HOH:O	2.09	0.49
1:A:81:ILE:CD1	4:A:222:HOH:O	2.59	0.49
1:A:102:ALA:CB	4:A:222:HOH:O	2.34	0.49
1:A:48:HIS:CB	4:A:288:HOH:O	2.60	0.49
1:A:101:GLN:CG	4:A:229:HOH:O	2.60	0.49
1:A:25:TYR:CZ	1:A:29:MET:HG3	2.48	0.48
1:A:104:LYS:CD	4:A:156:HOH:O	2.42	0.48
1:A:21:SER:HG	1:A:22:SER:HG	1.56	0.48
1:A:104:LYS:CE	4:A:156:HOH:O	2.62	0.48
1:A:32:SER:OG	4:A:227:HOH:O	2.19	0.48
1:A:44:ASN:O	1:A:84:CYS:HB2	2.13	0.48
1:A:67:ASN:HD21	1:A:69:GLN:HE22	0.57	0.48
1:A:104:LYS:CD	4:B:216:HOH:O	2.59	0.48
1:A:67:ASN:HB3	4:A:159:HOH:O	2.15	0.47
2:B:115:TYR:OH	4:B:218:HOH:O	2.19	0.47
1:A:69:GLN:CG	4:A:173:HOH:O	2.54	0.47
1:A:4:ALA:CB	2:B:118:VAL:HG22	2.44	0.46
1:A:37:LYS:O	1:A:92:TYR:OH	2.20	0.46
1:A:11:GLN:HG2	1:A:35:LEU:HD21	1.97	0.46
1:A:49:GLU:CG	4:A:335:HOH:O	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:26:CYS:H	1:A:99:THR:CG2	2.27	0.46
1:A:31:LYS:NZ	4:A:128:HOH:O	2.49	0.46
1:A:34:ASN:HA	1:A:37:LYS:HZ1	1.81	0.46
1:A:59:SER:HA	4:A:199:HOH:O	2.14	0.46
1:A:39:ARG:CG	4:A:276:HOH:O	2.60	0.45
1:A:91:LYS:HG2	4:A:180:HOH:O	2.17	0.44
1:A:104:LYS:HB3	2:B:123:SER:HG	1.82	0.44
1:A:66:LYS:O	1:A:67:ASN:HB3	2.17	0.44
1:A:8:PHE:CD1	1:A:8:PHE:C	2.91	0.44
1:A:50:SER:O	1:A:53:ASP:N	2.51	0.43
1:A:87:THR:C	1:A:89:SER:N	2.71	0.43
1:A:94:ASN:HB3	4:A:226:HOH:O	2.18	0.43
1:A:24:ASN:O	1:A:28:GLN:HG3	2.18	0.42
1:A:66:LYS:O	4:A:159:HOH:O	2.21	0.42
1:A:43:VAL:HG22	1:A:85:ARG:HG3	2.00	0.42
1:A:31:LYS:HB2	1:A:31:LYS:HE2	1.65	0.42
1:A:73:TYR:CE2	2:B:115:TYR:CE2	3.06	0.42
1:A:9:GLU:HA	1:A:13:MET:HG2	2.01	0.41
1:A:105:HIS:O	2:B:124:VAL:HA	2.19	0.41
1:A:27:ASN:HD21	1:A:97:TYR:N	2.16	0.41
1:A:99:THR:HG22	4:A:131:HOH:O	2.20	0.41
1:A:61:LYS:O	1:A:74:GLN:N	2.43	0.40
1:A:34:ASN:HA	1:A:37:LYS:NZ	2.36	0.40
1:A:99:THR:OG1	4:A:137:HOH:O	2.04	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	111/113 (98%)	100 (90%)	8 (7%)	3 (3%)	4 1

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	9/11 (82%)	9 (100%)	0	0	100	100
All	All	120/124 (97%)	109 (91%)	8 (7%)	3 (2%)	4	1

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	21	SER
1	A	67	ASN
1	A	88	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	99/99 (100%)	87 (88%)	12 (12%)	4	2
2	B	10/10 (100%)	7 (70%)	3 (30%)	0	0
All	All	109/109 (100%)	94 (86%)	15 (14%)	3	1

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	THR
1	A	21	SER
1	A	24	ASN
1	A	28	GLN
1	A	37	LYS
1	A	39	ARG
1	A	66	LYS
1	A	69	GLN
1	A	70	THR
1	A	93	PRO
1	A	99	THR
1	A	107	ILE
2	B	114	PRO

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Mol	Chain	Res	Type
2	B	118	VAL
2	B	124	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	ASN
1	A	27	ASN
1	A	28	GLN
1	A	34	ASN
1	A	69	GLN
1	A	74	GLN
1	A	94	ASN
1	A	103	ASN

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](#)

1 ligand is 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	$\# Z > 2$
3	SO4	B	125	-	4,4,4	0.86	0	6,6,6	0.42	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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	1:LYS	C	2:GLU	N	0.95

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.