



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:36 PM GMT

PDB ID : 1UER
Title : Crystal structure of Porphyromonas gingivalis SOD
Authors : Yamakura, F.; Sugio, S.; Hiraoka, B.Y.; Yokota, T.; Ohmori, D.
Deposited on : 2003-05-20
Resolution : 1.60 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

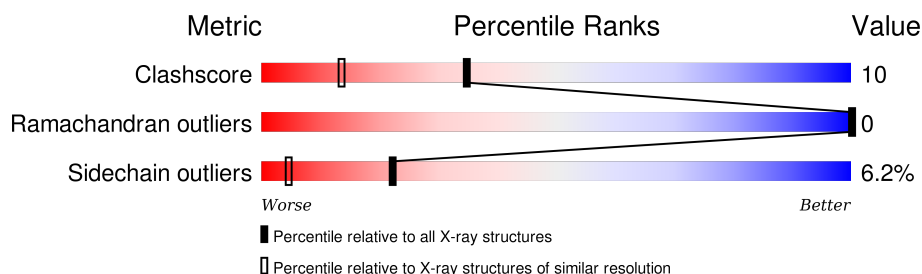
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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	102246	2732 (1.60-1.60)
Ramachandran outliers	100387	2654 (1.60-1.60)
Sidechain outliers	100360	2653 (1.60-1.60)

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. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	191	
1	B	191	
1	C	191	
1	D	191	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6612 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 superoxide dismutase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	191	Total	C	N	O	S	0	0	0
			1526	980	259	286	1			
1	B	191	Total	C	N	O	S	0	0	0
			1526	980	259	286	1			
1	C	191	Total	C	N	O	S	0	0	0
			1526	980	259	286	1			
1	D	191	Total	C	N	O	S	0	0	0
			1526	980	259	286	1			

- Molecule 2 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Fe	0	0
			1	1		
2	A	1	Total	Fe	0	0
			1	1		
2	D	1	Total	Fe	0	0
			1	1		
2	C	1	Total	Fe	0	0
			1	1		

- Molecule 3 is water.

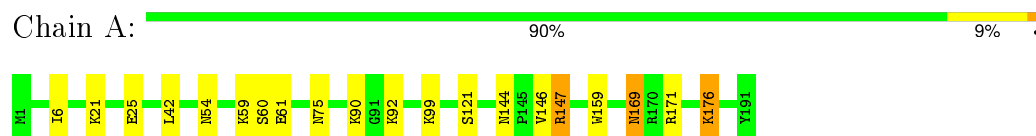
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	177	Total	O	0	0
			177	177		
3	B	156	Total	O	0	0
			156	156		
3	C	108	Total	O	0	0
			108	108		
3	D	63	Total	O	0	0
			63	63		

3 Residue-property plots [i](#)

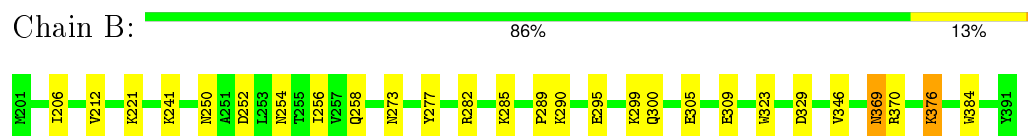
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

Note EDS was not executed.

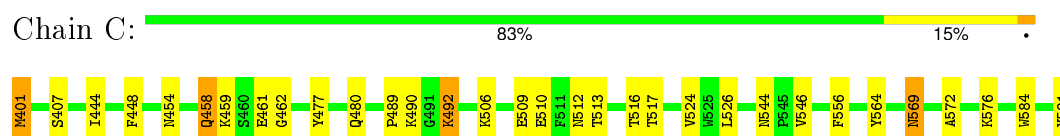
- Molecule 1: superoxide dismutase



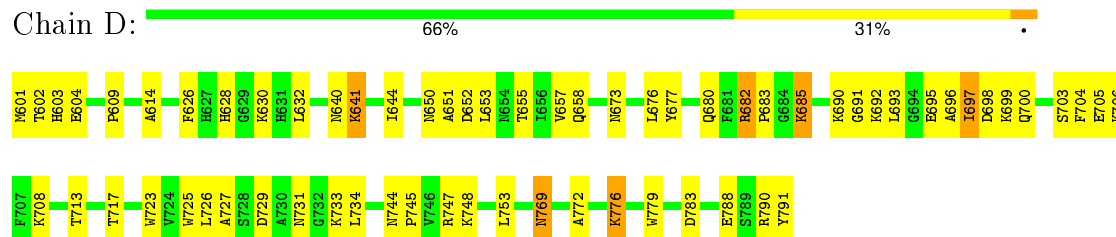
- Molecule 1: superoxide dismutase



- Molecule 1: superoxide dismutase



- Molecule 1: superoxide dismutase



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	75.32Å 95.93Å 101.39Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 1.60	Depositor
% Data completeness (in resolution range)	(Not available) (20.00-1.60)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 98.1	Depositor
R, R_{free}	0.231 , 0.256	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6612	wwPDB-VP
Average B, all atoms (Å ²)	33.0	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: FE

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	0.63	0/1567	0.76	1/2122 (0.0%)
1	B	0.62	0/1567	0.73	0/2122
1	C	0.58	0/1567	0.70	1/2122 (0.0%)
1	D	0.50	0/1567	0.66	0/2122
All	All	0.58	0/6268	0.71	2/8488 (0.0%)

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	B	0	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	171	ARG	NE-CZ-NH1	-5.74	117.43	120.30
1	C	526	LEU	N-CA-C	-5.04	97.39	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	277	TYR	Sidechain

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	1526	0	1479	18	0
1	B	1526	0	1476	23	0
1	C	1526	0	1476	26	0
1	D	1526	0	1476	53	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	177	0	0	9	0
3	B	156	0	0	9	0
3	C	108	0	0	4	0
3	D	63	0	0	4	0
All	All	6612	0	5907	119	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 10.

All (119) 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:D:682:ARG:HG2	1:D:682:ARG:HH11	1.22	1.01
1:B:376:LYS:HD2	3:B:534:HOH:O	1.72	0.89
1:C:492:LYS:HE2	1:C:492:LYS:O	1.73	0.88
1:A:6:ILE:HD13	1:A:75:ASN:HA	1.61	0.80
1:A:6:ILE:HG22	3:A:221:HOH:O	1.84	0.77
1:D:753:LEU:HD22	1:D:790:ARG:O	1.84	0.76
1:B:254:ASN:O	1:B:258:GLN:HG3	1.86	0.76
1:C:458:GLN:HE21	1:C:458:GLN:HA	1.50	0.74
1:D:682:ARG:CG	1:D:682:ARG:HH11	2.01	0.73
1:D:769:ASN:HD22	1:D:769:ASN:H	1.38	0.71
1:D:603:HIS:H	1:D:603:HIS:CD2	2.07	0.70
1:A:6:ILE:HB	3:A:270:HOH:O	1.92	0.70
1:D:626:PHE:CE1	1:D:630:LYS:HE2	2.28	0.69
1:C:461:GLU:HG2	1:C:462:GLY:H	1.59	0.68
1:D:682:ARG:HG2	1:D:682:ARG:NH1	2.02	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:691:GLY:O	1:D:695:GLU:HG3	1.95	0.66
1:C:461:GLU:HG2	1:C:462:GLY:N	2.10	0.65
1:A:169:ASN:HD22	1:A:169:ASN:H	1.46	0.64
1:D:693:LEU:O	1:D:697:ILE:HD12	1.97	0.64
1:C:492:LYS:HG3	1:C:591:TYR:HE2	1.64	0.63
1:D:609:PRO:HB2	1:D:682:ARG:HH21	1.64	0.62
1:B:241:LYS:HD3	1:B:241:LYS:N	2.13	0.62
1:C:458:GLN:HE21	1:C:458:GLN:CA	2.12	0.62
1:B:300:GLN:HB3	3:B:531:HOH:O	1.99	0.61
1:B:329:ASP:HB2	3:B:413:HOH:O	1.99	0.60
1:D:601:MET:N	1:D:650:ASN:HA	2.16	0.60
1:A:60:SER:O	1:A:61:GLU:HG3	2.02	0.60
1:B:369:ASN:H	1:B:369:ASN:HD22	1.51	0.59
1:D:693:LEU:HD23	1:D:788:GLU:HG3	1.84	0.59
1:D:677:TYR:O	1:D:680:GLN:HG2	2.03	0.58
1:D:685:LYS:NZ	1:D:685:LYS:HB2	2.19	0.57
1:D:788:GLU:C	1:D:790:ARG:H	2.08	0.57
1:D:753:LEU:CD2	1:D:790:ARG:O	2.52	0.57
1:D:685:LYS:HB2	1:D:685:LYS:HZ3	1.70	0.57
1:B:258:GLN:HG2	1:B:346:VAL:HG12	1.88	0.56
1:C:512:ASN:O	1:C:516:THR:HG23	2.04	0.56
1:D:685:LYS:HG2	1:D:783:ASP:HA	1.88	0.56
1:D:696:ALA:O	1:D:700:GLN:HG3	2.05	0.56
1:D:713:THR:O	1:D:717:THR:HG23	2.07	0.55
1:D:704:PHE:CZ	1:D:708:LYS:HD2	2.41	0.55
1:C:572:ALA:O	1:C:576:LYS:HG3	2.06	0.55
1:D:693:LEU:HD22	1:D:788:GLU:HA	1.89	0.54
1:C:569:ASN:HD22	1:C:569:ASN:H	1.54	0.54
1:D:703:SER:OG	1:D:706:LYS:HG3	2.06	0.54
1:B:299:LYS:CD	3:B:414:HOH:O	2.55	0.54
1:D:673:ASN:HB3	1:D:723:TRP:CZ2	2.43	0.54
1:C:524:VAL:CG1	1:C:556:PHE:HB2	2.38	0.54
1:D:708:LYS:HG2	1:D:779:TRP:CE2	2.42	0.54
1:B:285:LYS:HG2	3:B:546:HOH:O	2.08	0.53
1:C:513:THR:O	1:C:517:THR:HG23	2.08	0.53
1:B:285:LYS:NZ	3:B:546:HOH:O	2.42	0.53
1:C:458:GLN:NE2	1:C:546:VAL:O	2.43	0.52
1:C:458:GLN:HA	1:C:458:GLN:NE2	2.22	0.52
1:A:21:LYS:O	1:A:25:GLU:HG3	2.10	0.52
1:B:299:LYS:HD3	3:B:414:HOH:O	2.10	0.51
1:C:448:PHE:HE1	1:C:459:LYS:HG3	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:144:ASN:OD1	1:A:146:VAL:HG22	2.10	0.51
1:C:569:ASN:ND2	1:D:630:LYS:HD2	2.27	0.50
1:A:6:ILE:CD1	1:A:75:ASN:HA	2.40	0.49
1:D:601:MET:N	3:D:375:HOH:O	2.45	0.49
1:D:685:LYS:NZ	1:D:685:LYS:CB	2.76	0.49
1:C:444:ILE:HD12	3:C:697:HOH:O	2.11	0.49
1:D:640:ASN:CB	3:D:40:HOH:O	2.61	0.49
1:D:700:GLN:HE22	1:D:734:LEU:HG	1.79	0.48
1:B:289:PRO:HB3	1:B:384:TRP:CE3	2.48	0.48
1:B:370:ARG:NH2	3:B:533:HOH:O	2.33	0.48
1:D:734:LEU:HD21	1:D:791:TYR:CE1	2.49	0.47
1:D:696:ALA:HB1	1:D:700:GLN:HE21	1.81	0.46
1:D:769:ASN:HD22	1:D:769:ASN:N	2.08	0.46
1:C:492:LYS:HD3	3:C:632:HOH:O	2.16	0.46
1:B:241:LYS:N	1:B:241:LYS:CD	2.79	0.46
1:D:727:ALA:O	1:D:734:LEU:HA	2.16	0.46
1:B:295:GLU:O	1:B:299:LYS:HG2	2.16	0.46
1:D:772:ALA:O	1:D:776:LYS:HD3	2.17	0.45
1:D:640:ASN:HB3	3:D:40:HOH:O	2.16	0.45
1:A:147:ARG:CZ	3:A:207:HOH:O	2.64	0.45
1:D:641:LYS:N	3:D:40:HOH:O	2.49	0.45
1:D:614:ALA:O	1:D:683:PRO:HB3	2.17	0.45
1:D:652:ASP:OD1	1:D:655:THR:CB	2.65	0.45
1:B:299:LYS:HD3	1:B:299:LYS:N	2.30	0.45
1:A:6:ILE:HG23	3:A:195:HOH:O	2.17	0.45
1:D:731:ASN:H	1:D:731:ASN:HD22	1.65	0.44
1:B:273:ASN:HB3	1:B:323:TRP:CZ2	2.52	0.44
1:D:729:ASP:OD1	1:D:729:ASP:C	2.56	0.44
1:C:544:ASN:OD1	1:C:546:VAL:HG13	2.18	0.44
1:A:176:LYS:HE3	3:A:355:HOH:O	2.16	0.44
1:D:609:PRO:HB2	1:D:682:ARG:NH2	2.31	0.44
1:C:506:LYS:O	1:C:510:GLU:HG3	2.17	0.44
1:A:121:SER:HB3	1:A:159:TRP:CD2	2.53	0.43
1:D:685:LYS:CG	1:D:783:ASP:HA	2.49	0.43
1:B:212:VAL:HG23	1:B:221:LYS:HG3	2.00	0.43
1:D:632:LEU:HD12	1:D:632:LEU:O	2.19	0.43
1:C:477:TYR:O	1:C:480:GLN:HG2	2.17	0.43
1:D:628:HIS:O	1:D:632:LEU:HB3	2.19	0.43
1:C:489:PRO:HB3	1:C:584:TRP:CE3	2.53	0.43
1:A:147:ARG:NE	3:A:207:HOH:O	2.50	0.43
1:A:147:ARG:NH2	3:A:207:HOH:O	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:744:ASN:CG	1:D:745:PRO:HD2	2.39	0.43
1:C:458:GLN:CD	1:C:546:VAL:O	2.57	0.43
1:C:524:VAL:HG13	1:C:556:PHE:HB2	2.00	0.43
1:D:652:ASP:OD1	1:D:655:THR:HB	2.19	0.42
1:D:682:ARG:CG	1:D:682:ARG:NH1	2.67	0.42
1:C:401:MET:HB3	3:C:603:HOH:O	2.20	0.42
1:C:444:ILE:CD1	3:C:697:HOH:O	2.66	0.42
1:D:676:LEU:HB2	1:D:725:TRP:HH2	1.85	0.42
1:B:206:ILE:HD12	1:B:206:ILE:O	2.19	0.42
1:B:282:ARG:HB2	1:B:285:LYS:HD2	2.02	0.42
1:D:726:LEU:HG	1:D:753:LEU:HD12	2.02	0.41
1:D:653:LEU:O	1:D:657:VAL:HG23	2.20	0.41
1:B:282:ARG:CB	1:B:285:LYS:HE3	2.50	0.41
1:A:6:ILE:HG21	1:A:75:ASN:OD1	2.20	0.41
1:A:54:ASN:HB2	3:A:327:HOH:O	2.19	0.41
1:D:602:THR:HG22	1:D:602:THR:O	2.20	0.41
1:A:6:ILE:CG2	1:A:75:ASN:OD1	2.69	0.41
1:C:458:GLN:CA	1:C:458:GLN:NE2	2.83	0.41
1:A:6:ILE:CG2	3:A:270:HOH:O	2.68	0.41
1:B:252:ASP:O	1:B:256:ILE:HG13	2.21	0.41
1:B:370:ARG:NE	3:B:533:HOH:O	2.34	0.41
1:D:601:MET:HG3	1:D:651:ALA:O	2.22	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	189/191 (99%)	183 (97%)	6 (3%)	0	100	100
1	B	189/191 (99%)	183 (97%)	6 (3%)	0	100	100
1	C	189/191 (99%)	184 (97%)	5 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	189/191 (99%)	178 (94%)	11 (6%)	0	100	100
All	All	756/764 (99%)	728 (96%)	28 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	160/160 (100%)	152 (95%)	8 (5%)	30	7
1	B	160/160 (100%)	154 (96%)	6 (4%)	40	13
1	C	160/160 (100%)	151 (94%)	9 (6%)	26	6
1	D	160/160 (100%)	143 (89%)	17 (11%)	8	1
All	All	640/640 (100%)	600 (94%)	40 (6%)	22	4

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	42	LEU
1	A	59	LYS
1	A	90	LYS
1	A	92	LYS
1	A	99	LYS
1	A	147	ARG
1	A	169	ASN
1	A	176	LYS
1	B	250	ASN
1	B	290	LYS
1	B	305	GLU
1	B	309	GLU
1	B	369	ASN
1	B	376	LYS
1	C	401	MET
1	C	407	SER

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Mol	Chain	Res	Type
1	C	454	ASN
1	C	458	GLN
1	C	490	LYS
1	C	492	LYS
1	C	509	GLU
1	C	564	TYR
1	C	569	ASN
1	D	604	GLU
1	D	641	LYS
1	D	644	ILE
1	D	658	GLN
1	D	682	ARG
1	D	685	LYS
1	D	690	LYS
1	D	692	LYS
1	D	697	ILE
1	D	698	ASP
1	D	699	LYS
1	D	705	GLU
1	D	733	LYS
1	D	747	ARG
1	D	748	LYS
1	D	769	ASN
1	D	776	LYS

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

Mol	Chain	Res	Type
1	A	169	ASN
1	B	369	ASN
1	C	458	GLN
1	C	569	ASN
1	D	603	HIS
1	D	675	ASN
1	D	731	ASN
1	D	769	ASN

5.3.3 RNA ⓘ

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - 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.

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

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates

EDS was not executed - this section will therefore be empty.

6.4 Ligands

EDS was not executed - this section will therefore be empty.

6.5 Other polymers

EDS was not executed - this section will therefore be empty.