



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:57 AM UTC

PDB ID : 1DKI / pdb_00001dki
Title : CRYSTAL STRUCTURE OF THE ZYMOGEN FORM OF STREPTOCOCCAL PYROGENIC EXOTOXIN B ACTIVE SITE (C47S) MUTANT
Authors : Kagawa, T.F.; Cooney, J.C.; Baker, H.M.; McSweeney, S.; Liu, M.; Gubba, S.; Musser, J.M.; Baker, E.N.
Deposited on : 1999-12-07
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

<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

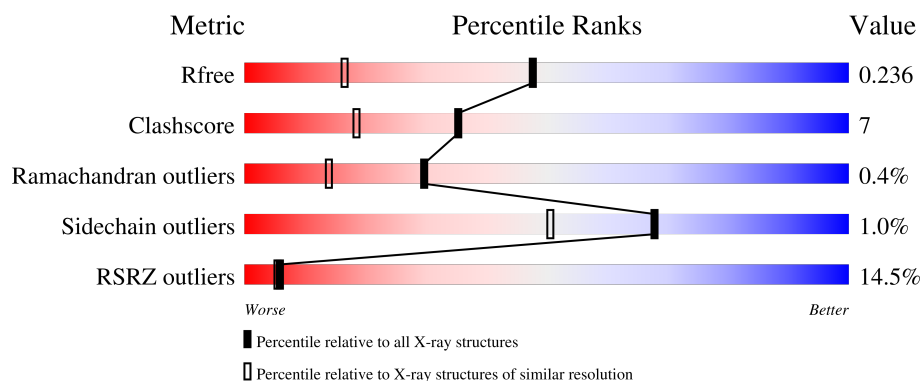
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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))
R_{free}	180053	4673 (1.60-1.60)
Clashscore	190562	4931 (1.60-1.60)
Ramachandran outliers	187476	4831 (1.60-1.60)
Sidechain outliers	187428	4830 (1.60-1.60)
RSRZ outliers	180081	4672 (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 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	371	15% 66% 15% • 18%
1	B	371	18% 65% 20% • 14%
1	C	371	9% 78% 13% 9%
1	D	371	9% 77% 12% 10%

2 Entry composition [i](#)

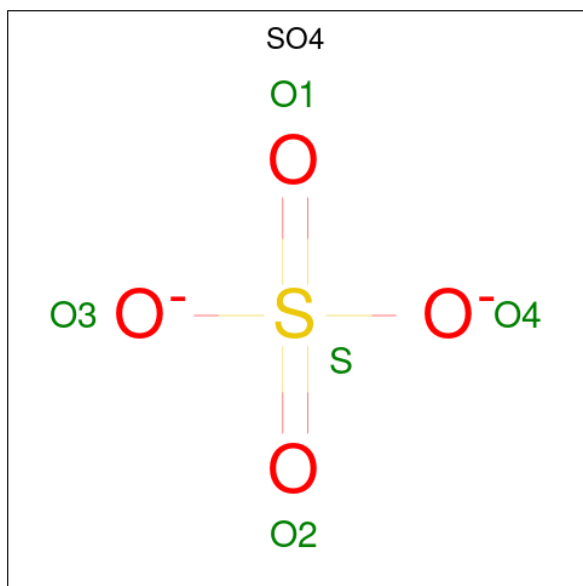
There are 3 unique types of molecules in this entry. The entry contains 10562 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PYROGENIC EXOTOXIN B ZYMOGEN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	303	Total	C	N	O	S	0	0	0
			2312	1463	400	443	6			
1	B	320	Total	C	N	O	S	0	0	2
			2426	1529	425	466	6			
1	C	339	Total	C	N	O	S	0	0	1
			2590	1638	446	500	6			
1	D	334	Total	C	N	O	S	0	0	0
			2575	1624	444	501	6			

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

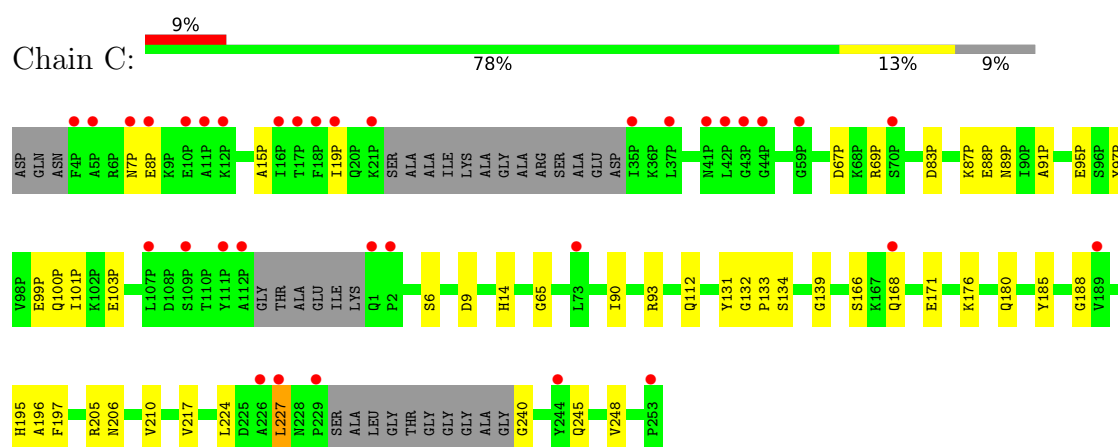
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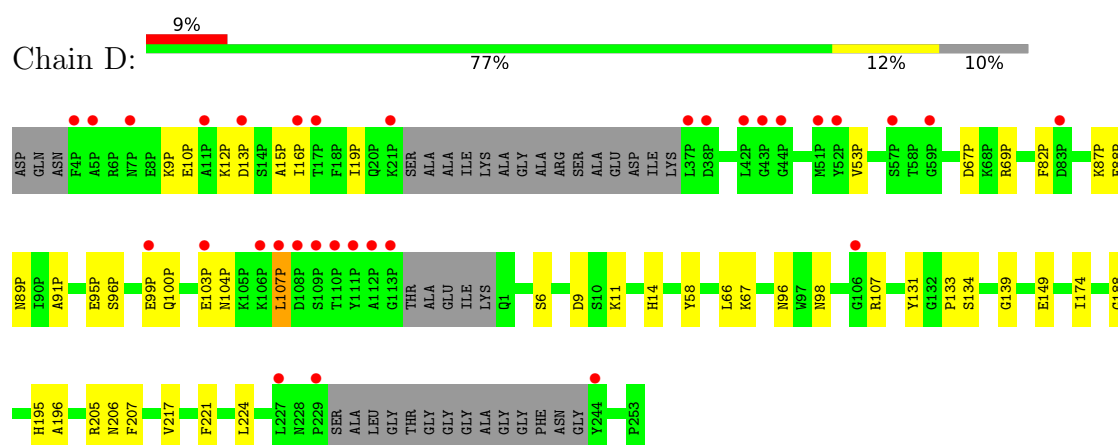
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	120	Total	O	0	0
			120	120		
3	B	98	Total	O	0	0
			98	98		
3	C	188	Total	O	0	0
			188	188		
3	D	233	Total	O	0	0
			233	233		



● Molecule 1: PYROGENIC EXOTOXIN B ZYMOGEN



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	46.66Å 116.56Å 144.73Å 90.00° 94.17° 90.00°	Depositor
Resolution (Å)	30.00 – 1.60 30.00 – 1.60	Depositor EDS
% Data completeness (in resolution range)	83.1 (30.00-1.60) 83.1 (30.00-1.60)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.48 (at 1.60Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.218 , 0.245 0.210 , 0.236	Depositor DCC
R_{free} test set	16726 reflections (9.93%)	wwPDB-VP
Wilson B-factor (Å ²)	29.6	Xtriage
Anisotropy	0.036	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 40.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	10562	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.47% 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: 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	0.53	0/2365	0.95	9/3195 (0.3%)
1	B	0.44	0/2480	0.94	12/3350 (0.4%)
1	C	0.54	0/2647	0.96	12/3576 (0.3%)
1	D	0.59	0/2632	0.96	8/3556 (0.2%)
All	All	0.53	0/10124	0.95	41/13677 (0.3%)

There are no bond length outliers.

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	166	SER	N-CA-C	-8.38	99.24	110.55
1	B	47(P)	SER	N-CA-C	7.88	119.78	111.03
1	B	133	PRO	N-CA-C	-7.55	103.45	113.65
1	B	166	SER	N-CA-C	-7.36	100.62	110.55
1	C	139	GLY	N-CA-C	7.14	119.92	111.35
1	D	139	GLY	N-CA-C	7.10	119.88	111.35
1	A	134	SER	N-CA-C	-7.02	98.41	109.50
1	D	134	SER	N-CA-C	-6.96	98.58	109.72
1	A	196	ALA	N-CA-C	6.96	120.83	110.52
1	C	133	PRO	N-CA-C	-6.96	103.86	113.53
1	C	196	ALA	N-CA-C	6.95	120.80	110.52
1	C	188	GLY	N-CA-C	-6.88	99.57	112.10
1	A	133	PRO	N-CA-C	-6.82	104.44	113.65
1	C	245	GLN	N-CA-C	6.75	119.54	110.35
1	A	166	SER	N-CA-C	-6.75	101.44	110.55
1	D	131	TYR	N-CA-C	6.58	120.02	110.28
1	D	133	PRO	N-CA-C	-6.56	104.79	113.65
1	D	149	GLU	N-CA-C	6.54	118.20	111.14
1	B	188	GLY	N-CA-C	-6.48	100.31	112.10
1	C	134	SER	N-CA-C	-6.46	99.38	109.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	131	TYR	N-CA-C	6.43	119.80	110.28
1	A	139	GLY	N-CA-C	6.42	119.06	111.35
1	B	134	SER	N-CA-C	-6.32	99.60	109.72
1	A	131	TYR	N-CA-C	6.18	119.43	110.28
1	B	38	PHE	N-CA-C	-6.13	104.44	112.41
1	D	188	GLY	N-CA-C	-6.08	101.04	112.10
1	B	131	TYR	N-CA-C	6.01	119.17	110.28
1	B	196	ALA	N-CA-C	5.92	119.59	110.42
1	C	93	ARG	N-CA-C	5.91	119.89	109.96
1	D	196	ALA	N-CA-C	5.90	119.57	110.42
1	B	245	GLN	N-CA-C	5.88	118.35	110.35
1	B	165	PHE	N-CA-C	5.86	118.61	109.52
1	D	88(P)	GLU	N-CA-C	5.85	118.44	111.71
1	A	188	GLY	N-CA-C	-5.77	101.60	112.10
1	A	66	LEU	N-CA-C	5.57	120.09	113.12
1	C	132	GLY	CA-C-N	5.45	125.09	119.05
1	C	132	GLY	C-N-CA	5.45	125.09	119.05
1	B	93	ARG	N-CA-C	5.37	118.30	109.76
1	C	88(P)	GLU	N-CA-C	5.26	117.76	111.71
1	B	223	ARG	N-CA-C	-5.08	102.96	110.48
1	A	184	VAL	N-CA-C	5.01	115.18	108.17

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	2312	0	2196	40	0
1	B	2426	0	2287	53	0
1	C	2590	0	2455	25	0
1	D	2575	0	2457	25	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
2	D	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	120	0	0	3	0
3	B	98	0	0	0	0
3	C	188	0	0	3	0
3	D	233	0	0	2	0
All	All	10562	0	9395	142	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 (142) 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:C:7(P):ASN:CA	1:C:8(P):GLU:N	1.79	1.44
1:B:6(P):ARG:HH12	1:B:72(P):GLU:HA	1.02	1.08
1:B:225:ASP:CA	1:B:226:ALA:N	2.21	1.04
1:B:99(P):GLU:CA	1:B:100(P):GLN:N	2.28	0.96
1:B:6(P):ARG:NH1	1:B:72(P):GLU:HA	1.84	0.90
1:B:154:ASN:HD22	1:B:156:SER:H	1.12	0.89
1:B:53(P):VAL:HG22	1:B:63(P):ILE:HD12	1.58	0.83
1:C:99(P):GLU:CA	1:C:100(P):GLN:N	2.45	0.79
1:A:39(P):LYS:HB2	1:A:39(P):LYS:NZ	1.97	0.78
1:A:66:LEU:O	1:A:67:LYS:HD2	1.84	0.77
1:B:102(P):LYS:HA	1:B:105(P):LYS:HD3	1.68	0.75
1:A:64(P):VAL:HG12	1:A:73(P):ILE:HD13	1.69	0.74
1:D:96:ASN:C	1:D:96:ASN:HD22	1.96	0.74
1:C:176:LYS:HE2	1:C:180:GLN:NE2	2.03	0.73
1:C:14:HIS:HD2	3:C:1149:HOH:O	1.74	0.71
1:B:12(P):LYS:HB2	1:B:12(P):LYS:NZ	2.06	0.70
1:A:114:MET:HE3	3:A:1095:HOH:O	1.92	0.70
1:A:101(P):ILE:HA	1:A:104(P):ASN:HD22	1.56	0.70
1:C:87(P):LYS:HE2	1:C:217:VAL:HG11	1.74	0.69
1:B:5(P):ALA:N	1:B:67(P):ASP:HA	2.09	0.68
1:B:66:LEU:O	1:B:67:LYS:HD2	1.94	0.66
1:A:106:GLY:HA2	3:A:1059:HOH:O	1.96	0.65
1:B:56(P):ILE:HB	1:B:59(P):GLY:O	1.96	0.64
1:A:39(P):LYS:HB2	1:A:39(P):LYS:HZ2	1.63	0.64
1:A:146:ALA:HA	1:A:150:ASN:HD22	1.64	0.62
1:C:176:LYS:HE2	1:C:180:GLN:HE22	1.64	0.61
1:B:205:ARG:O	1:B:206:ASN:HB2	1.99	0.61
1:B:241:PHE:CZ	1:B:243:GLY:HA2	2.35	0.61
1:A:73:LEU:HD12	3:A:1058:HOH:O	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:46(P):LEU:HD21	1:A:95(P):GLU:HG3	1.82	0.60
1:D:14:HIS:HD2	3:D:1052:HOH:O	1.85	0.59
1:A:203:ASP:OD2	1:A:207:PHE:HB2	2.03	0.58
1:B:154:ASN:ND2	1:B:156:SER:H	1.94	0.58
1:D:10(P):GLU:O	1:D:13(P):ASP:HB3	2.04	0.58
1:A:70(P):SER:HB3	1:A:104(P):ASN:ND2	2.19	0.58
1:B:61:TYR:CG	1:B:62:PRO:HA	2.39	0.58
1:D:87(P):LYS:HE2	1:D:217:VAL:HG11	1.86	0.57
1:B:1:GLN:HE22	1:B:202:ALA:CB	2.18	0.56
1:B:89(P):ASN:HA	1:B:195:HIS:HB3	1.87	0.56
1:B:67(P):ASP:OD2	1:B:69(P):ARG:HB2	2.05	0.56
1:B:61:TYR:CD1	1:B:62:PRO:HA	2.40	0.56
1:B:162:ARG:NH2	1:B:167:LYS:HG2	2.21	0.56
1:D:96(P):SER:O	1:D:100(P):GLN:HG3	2.05	0.56
1:A:207:PHE:HB3	1:A:221:PHE:HB3	1.89	0.55
1:B:154:ASN:HD22	1:B:156:SER:N	1.93	0.55
1:B:49(P):SER:HB2	1:B:52(P):TYR:OH	2.06	0.55
1:B:86:LEU:HD12	1:B:86:LEU:N	2.22	0.54
1:D:205:ARG:O	1:D:206:ASN:HB2	2.07	0.54
1:C:89(P):ASN:HA	1:C:195:HIS:HB3	1.89	0.54
1:A:101(P):ILE:HA	1:A:104(P):ASN:ND2	2.23	0.54
1:B:96(P):SER:O	1:B:100(P):GLN:HG3	2.08	0.54
1:B:1:GLN:HE22	1:B:202:ALA:HB1	1.72	0.53
1:D:89(P):ASN:HA	1:D:195:HIS:HB3	1.91	0.53
1:A:205:ARG:O	1:A:206:ASN:HB2	2.09	0.52
1:C:100(P):GLN:HB3	1:C:227:LEU:HD13	1.91	0.52
1:A:37(P):LEU:HB3	1:A:54(P):TYR:HB3	1.91	0.52
1:A:66:LEU:C	1:A:67:LYS:HD2	2.34	0.52
1:C:14:HIS:HE1	3:C:1105:HOH:O	1.92	0.52
1:C:168:GLN:H	1:C:168:GLN:CD	2.18	0.51
1:B:12(P):LYS:HB2	1:B:12(P):LYS:HZ3	1.74	0.51
1:D:207:PHE:HB3	1:D:221:PHE:HB3	1.92	0.51
1:B:207:PHE:HB3	1:B:221:PHE:HB3	1.92	0.51
1:B:94(P):MET:O	1:B:97(P):TYR:HB2	2.11	0.50
1:C:205:ARG:O	1:C:206:ASN:HB2	2.10	0.49
1:D:96:ASN:HD22	1:D:98:ASN:H	1.60	0.49
1:A:53(P):VAL:HG22	1:A:63(P):ILE:HG23	1.94	0.49
1:B:104(P):ASN:C	1:B:106(P):LYS:H	2.20	0.49
1:C:97(P):TYR:O	1:C:101(P):ILE:HG13	2.12	0.49
1:C:112:GLN:CD	3:C:1094:HOH:O	2.56	0.49
1:A:61:TYR:CG	1:A:62:PRO:HA	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:43(P):GLY:HA2	1:A:47(P):SER:HB3	1.95	0.48
1:B:41(P):ASN:C	1:B:41(P):ASN:HD22	2.21	0.48
1:B:47(P):SER:O	1:B:48(P):GLY:O	2.31	0.48
1:A:203:ASP:CG	1:A:207:PHE:HB2	2.39	0.48
1:C:6:SER:HB3	1:C:9:ASP:HB2	1.95	0.48
1:D:174:ILE:HD12	1:D:224:LEU:HD11	1.96	0.47
1:A:102(P):LYS:O	1:A:105(P):LYS:HB2	2.14	0.47
1:B:71:TYR:OH	1:B:86:LEU:HD22	2.14	0.47
1:D:14:HIS:HE1	3:D:1113:HOH:O	1.97	0.47
1:B:42(P):LEU:HD11	1:B:53(P):VAL:HG23	1.95	0.47
1:B:157:VAL:HG22	1:B:251:ILE:HA	1.97	0.47
1:D:91(P):ALA:O	1:D:95(P):GLU:HG3	2.14	0.46
1:A:186:TYR:CD1	1:A:199:ILE:HD12	2.51	0.46
1:C:15(P):ALA:O	1:C:19(P):ILE:HG13	2.15	0.46
1:D:6:SER:HB3	1:D:9:ASP:HB2	1.98	0.46
1:D:11:LYS:HD2	1:D:58:TYR:CD1	2.51	0.46
1:A:64(P):VAL:HG12	1:A:73(P):ILE:CD1	2.44	0.45
1:A:5:LYS:O	1:A:6:SER:C	2.59	0.45
1:D:96:ASN:HD21	1:D:98:ASN:HB2	1.81	0.45
1:B:40(P):VAL:HG22	1:B:55(P):ASN:ND2	2.32	0.45
1:D:15(P):ALA:O	1:D:19(P):ILE:HG13	2.17	0.45
1:D:96:ASN:C	1:D:96:ASN:ND2	2.71	0.45
1:D:96:ASN:ND2	1:D:98:ASN:H	2.15	0.45
1:A:214:TRP:CZ3	1:B:107:ARG:HD3	2.52	0.45
1:D:99(P):GLU:O	1:D:103(P):GLU:HB2	2.16	0.45
1:A:6:SER:HB3	1:A:9:ASP:HB2	1.98	0.45
1:B:64:LYS:HE3	1:B:64:LYS:HB3	1.79	0.45
1:C:168:GLN:CD	1:C:168:GLN:N	2.75	0.44
1:A:70(P):SER:HB3	1:A:104(P):ASN:HD22	1.82	0.44
1:A:39(P):LYS:HB2	1:A:39(P):LYS:HZ3	1.82	0.44
1:C:91(P):ALA:O	1:C:95(P):GLU:HG3	2.18	0.44
1:B:56(P):ILE:HB	1:B:59(P):GLY:C	2.42	0.44
1:C:100(P):GLN:HG2	1:C:227:LEU:HD22	2.00	0.44
1:B:104(P):ASN:C	1:B:106(P):LYS:N	2.75	0.43
1:B:176:LYS:HG2	1:B:180:GLN:NE2	2.33	0.43
1:A:87(P):LYS:HZ3	1:A:217:VAL:HG11	1.83	0.43
1:B:50(P):ASN:O	1:B:101(P):ILE:HD13	2.18	0.43
1:B:162:ARG:HH21	1:B:167:LYS:HG2	1.82	0.43
1:C:83(P):ASP:OD1	1:C:87(P):LYS:NZ	2.52	0.43
1:B:84(P):VAL:HG12	1:B:90(P):ILE:CG2	2.49	0.43
1:B:42(P):LEU:HB3	1:B:46(P):LEU:HB2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:12(P):LYS:HB2	1:B:12(P):LYS:HZ2	1.81	0.42
1:B:60(P):GLY:HA2	1:B:79(P):SER:O	2.18	0.42
1:C:103(P):GLU:OE1	1:C:240:GLY:HA2	2.19	0.42
1:D:67(P):ASP:OD2	1:D:69(P):ARG:HB2	2.19	0.42
1:B:154:ASN:ND2	1:B:155:GLN:N	2.68	0.42
1:C:65:GLY:HA3	1:C:90:ILE:O	2.19	0.42
1:A:46(P):LEU:CD2	1:A:95(P):GLU:HG3	2.48	0.42
1:A:84(P):VAL:HG12	1:A:90(P):ILE:CG2	2.50	0.42
1:B:41(P):ASN:HD22	1:B:43(P):GLY:H	1.68	0.42
1:B:66:LEU:C	1:B:67:LYS:HD2	2.44	0.42
1:A:69(P):ARG:HD2	1:A:104(P):ASN:O	2.19	0.42
1:D:9(P):LYS:O	1:D:13(P):ASP:HB2	2.20	0.41
1:B:97(P):TYR:O	1:B:101(P):ILE:HG13	2.20	0.41
1:D:53(P):VAL:HG11	1:D:82(P):PHE:CD2	2.55	0.41
1:D:66:LEU:O	1:D:67:LYS:HD3	2.20	0.41
1:D:12(P):LYS:O	1:D:16(P):ILE:HG13	2.21	0.41
1:A:61(P):PHE:CE2	1:A:77(P):SER:HB3	2.56	0.41
1:A:97(P):TYR:O	1:A:101(P):ILE:HG13	2.21	0.41
1:C:171:GLU:HG2	1:C:224:LEU:HD12	2.02	0.41
1:A:107:ARG:HH11	1:A:107:ARG:HD2	1.74	0.41
1:B:41(P):ASN:C	1:B:41(P):ASN:ND2	2.79	0.41
1:B:6:SER:HB3	1:B:9:ASP:HB2	2.03	0.41
1:C:197:PHE:CD1	1:C:210:VAL:HG13	2.56	0.41
1:D:104(P):ASN:O	1:D:107(P):LEU:HB2	2.21	0.41
1:A:95(P):GLU:O	1:A:99(P):GLU:HG3	2.21	0.41
1:C:67(P):ASP:OD1	1:C:69(P):ARG:HB2	2.20	0.41
1:C:185:TYR:HB3	1:C:248:VAL:HB	2.03	0.41
1:A:39(P):LYS:HB3	1:A:54(P):TYR:CE2	2.56	0.40
1:A:67(P):ASP:OD1	1:A:69(P):ARG:HB2	2.21	0.40
1:A:87(P):LYS:HZ3	1:A:217:VAL:CG1	2.34	0.40
1:B:154:ASN:HD22	1:B:155:GLN:N	2.19	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	295/371 (80%)	281 (95%)	13 (4%)	1 (0%)	36	20
1	B	308/371 (83%)	293 (95%)	11 (4%)	4 (1%)	9	2
1	C	327/371 (88%)	323 (99%)	4 (1%)	0	100	100
1	D	326/371 (88%)	319 (98%)	7 (2%)	0	100	100
All	All	1256/1484 (85%)	1216 (97%)	35 (3%)	5 (0%)	30	14

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	39(P)	LYS
1	B	43(P)	GLY
1	B	44(P)	GLY
1	B	48(P)	GLY
1	A	68(P)	LYS

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	241/300 (80%)	237 (98%)	4 (2%)	53	30
1	B	251/300 (84%)	248 (99%)	3 (1%)	63	43
1	C	270/300 (90%)	269 (100%)	1 (0%)	84	75
1	D	272/300 (91%)	270 (99%)	2 (1%)	76	63
All	All	1034/1200 (86%)	1024 (99%)	10 (1%)	68	50

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39(P)	LYS
1	A	39	VAL
1	A	112	GLN

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Mol	Chain	Res	Type
1	A	191	LYS
1	B	12(P)	LYS
1	B	1	GLN
1	B	154	ASN
1	C	227	LEU
1	D	107(P)	LEU
1	D	107	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	50(P)	ASN
1	A	104(P)	ASN
1	A	19	ASN
1	A	150	ASN
1	B	41(P)	ASN
1	B	50(P)	ASN
1	B	1	GLN
1	B	76	ASN
1	B	94	GLN
1	B	98	ASN
1	B	154	ASN
1	B	180	GLN
1	B	182	GLN
1	C	20(P)	GLN
1	C	14	HIS
1	C	19	ASN
1	C	159	GLN
1	C	161	ASN
1	C	180	GLN
1	D	50(P)	ASN
1	D	85(P)	ASN
1	D	14	HIS
1	D	63	ASN
1	D	94	GLN
1	D	96	ASN
1	D	159	GLN
1	D	161	ASN
1	D	173	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](#)

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	SO4	B	1002	-	4,4,4	0.43	0	6,6,6	0.19	0
2	SO4	D	1004	-	4,4,4	0.46	0	6,6,6	0.35	0
2	SO4	A	1001	-	4,4,4	0.45	0	6,6,6	0.28	0
2	SO4	C	1003	-	4,4,4	0.55	0	6,6,6	0.50	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

There are no such residues in this entry.

5.8 Polymer linkage issues

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	303/371 (81%)	1.00	54 (17%) 4 3	21, 36, 70, 73	0
1	B	320/371 (86%)	1.15	68 (21%) 2 2	21, 40, 68, 76	0
1	C	339/371 (91%)	0.68	34 (10%) 12 12	20, 34, 63, 72	0
1	D	334/371 (90%)	0.54	32 (9%) 13 13	18, 29, 62, 74	0
All	All	1296/1484 (87%)	0.84	188 (14%) 6 5	18, 35, 67, 76	0

All (188) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	64(P)	VAL	7.0
1	A	62(P)	VAL	6.2
1	C	112(P)	ALA	5.9
1	B	59(P)	GLY	5.7
1	C	7(P)	ASN	5.5
1	B	5(P)	ALA	5.1
1	A	244	TYR	5.1
1	A	53(P)	VAL	5.0
1	B	99(P)	GLU	4.9
1	D	113(P)	GLY	4.8
1	C	35(P)	ILE	4.7
1	D	7(P)	ASN	4.6
1	A	60(P)	GLY	4.5
1	D	103(P)	GLU	4.4
1	A	52(P)	TYR	4.4
1	D	4(P)	PHE	4.4
1	B	86(P)	GLY	4.3
1	A	204	GLY	4.3
1	B	204	GLY	4.2
1	A	98(P)	VAL	4.1
1	A	2	PRO	4.1

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Mol	Chain	Res	Type	RSRZ
1	A	55(P)	ASN	4.1
1	D	108(P)	ASP	4.0
1	C	5(P)	ALA	4.0
1	B	70(P)	SER	3.9
1	B	58(P)	THR	3.9
1	D	244	TYR	3.9
1	A	107(P)	LEU	3.8
1	A	54(P)	TYR	3.8
1	B	57(P)	SER	3.8
1	C	109(P)	SER	3.8
1	B	63(P)	ILE	3.8
1	D	112(P)	ALA	3.7
1	D	13(P)	ASP	3.7
1	A	3	VAL	3.7
1	B	229	PRO	3.7
1	B	2	PRO	3.6
1	C	4(P)	PHE	3.6
1	C	111(P)	TYR	3.6
1	C	226	ALA	3.5
1	A	73(P)	ILE	3.4
1	B	98(P)	VAL	3.3
1	D	16(P)	ILE	3.2
1	C	227	LEU	3.2
1	C	8(P)	GLU	3.2
1	D	106(P)	LYS	3.2
1	A	220	GLY	3.2
1	B	178	LEU	3.1
1	B	225	ASP	3.1
1	B	241	PHE	3.1
1	B	71(P)	PRO	3.1
1	B	56(P)	ILE	3.1
1	D	11(P)	ALA	3.1
1	A	38(P)	ASP	3.1
1	B	43(P)	GLY	3.1
1	A	46(P)	LEU	3.1
1	B	84(P)	VAL	3.0
1	A	107	ARG	3.0
1	A	63(P)	ILE	3.0
1	A	101(P)	ILE	3.0
1	D	107(P)	LEU	3.0
1	C	2	PRO	3.0
1	A	66(P)	GLY	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	4	VAL	3.0
1	A	106	GLY	3.0
1	B	3	VAL	3.0
1	A	104(P)	ASN	3.0
1	D	42(P)	LEU	3.0
1	A	65(P)	SER	2.9
1	A	105	SER	2.9
1	A	76(P)	TYR	2.9
1	A	61(P)	PHE	2.9
1	B	49(P)	SER	2.9
1	B	52(P)	TYR	2.9
1	B	54(P)	TYR	2.9
1	A	80(P)	GLY	2.8
1	B	60(P)	GLY	2.8
1	D	111(P)	TYR	2.8
1	A	37(P)	LEU	2.8
1	B	42(P)	LEU	2.8
1	C	16(P)	ILE	2.8
1	B	1	GLN	2.8
1	D	52(P)	TYR	2.8
1	A	227	LEU	2.8
1	A	42(P)	LEU	2.8
1	B	48(P)	GLY	2.8
1	D	229	PRO	2.7
1	C	42(P)	LEU	2.7
1	C	244	TYR	2.7
1	C	1	GLN	2.7
1	B	242	ASN	2.7
1	B	38(P)	ASP	2.7
1	B	172	ALA	2.7
1	A	68(P)	LYS	2.6
1	B	228	ASN	2.6
1	B	202	ALA	2.6
1	A	85(P)	ASN	2.6
1	B	192	VAL	2.6
1	B	107(P)	LEU	2.5
1	C	107(P)	LEU	2.5
1	B	40(P)	VAL	2.5
1	B	11(P)	ALA	2.5
1	B	7	LEU	2.5
1	D	110(P)	THR	2.5
1	B	61(P)	PHE	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	53(P)	VAL	2.5
1	C	19(P)	ILE	2.5
1	A	203	ASP	2.4
1	C	43(P)	GLY	2.4
1	D	17(P)	THR	2.4
1	B	68(P)	LYS	2.4
1	D	99(P)	GLU	2.4
1	C	37(P)	LEU	2.4
1	C	44(P)	GLY	2.4
1	C	229	PRO	2.4
1	C	168	GLN	2.4
1	B	176	LYS	2.4
1	A	41(P)	ASN	2.3
1	A	71(P)	PRO	2.3
1	C	253	PRO	2.3
1	D	51(P)	MET	2.3
1	B	97	TRP	2.3
1	C	189	VAL	2.3
1	D	106	GLY	2.3
1	A	178	LEU	2.3
1	A	82(P)	PHE	2.3
1	C	10(P)	GLU	2.3
1	B	106	GLY	2.3
1	B	167	LYS	2.3
1	B	78(P)	THR	2.3
1	D	37(P)	LEU	2.3
1	A	70(P)	SER	2.3
1	B	221	PHE	2.3
1	C	70(P)	SER	2.3
1	D	57(P)	SER	2.3
1	A	75(P)	GLY	2.3
1	A	103(P)	GLU	2.2
1	B	55(P)	ASN	2.2
1	D	109(P)	SER	2.2
1	B	9(P)	LYS	2.2
1	B	80(P)	GLY	2.2
1	D	44(P)	GLY	2.2
1	B	62(P)	VAL	2.2
1	C	17(P)	THR	2.2
1	B	79(P)	SER	2.2
1	B	66	LEU	2.2
1	B	44(P)	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	201	GLY	2.2
1	A	174	ILE	2.2
1	D	38(P)	ASP	2.2
1	B	74(P)	LEU	2.2
1	B	86	LEU	2.2
1	B	220	GLY	2.2
1	C	59(P)	GLY	2.2
1	A	97(P)	TYR	2.2
1	A	47(P)	SER	2.2
1	D	5(P)	ALA	2.1
1	B	69(P)	ARG	2.1
1	A	48(P)	GLY	2.1
1	B	190	GLY	2.1
1	B	7(P)	ASN	2.1
1	B	41(P)	ASN	2.1
1	B	208	TYR	2.1
1	A	40(P)	VAL	2.1
1	D	59(P)	GLY	2.1
1	C	12(P)	LYS	2.1
1	C	41(P)	ASN	2.1
1	D	227	LEU	2.1
1	C	18(P)	PHE	2.1
1	B	67	LYS	2.1
1	B	226	ALA	2.1
1	D	21(P)	LYS	2.1
1	D	43(P)	GLY	2.1
1	B	64(P)	VAL	2.1
1	A	49(P)	SER	2.1
1	B	227	LEU	2.1
1	C	73	LEU	2.1
1	A	212	TRP	2.1
1	A	207	PHE	2.1
1	B	67(P)	ASP	2.0
1	D	83(P)	ASP	2.0
1	A	205	ARG	2.0
1	A	44(P)	GLY	2.0
1	A	100(P)	GLN	2.0
1	B	73(P)	ILE	2.0
1	A	202	ALA	2.0
1	C	11(P)	ALA	2.0
1	C	21(P)	LYS	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	SO4	C	1003	5/5	0.96	0.08	28,29,31,31	0
2	SO4	B	1002	5/5	0.97	0.07	31,34,34,35	0
2	SO4	A	1001	5/5	0.97	0.08	33,35,36,37	0
2	SO4	D	1004	5/5	0.98	0.07	25,27,27,28	0

6.5 Other polymers [i](#)

There are no such residues in this entry.