



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:35 PM UTC

PDB ID : 3KSK / pdb_00003ksk
Title : Crystal Structure of single chain PvuII
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Deposited on : 2009-11-23
Resolution : 2.35 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

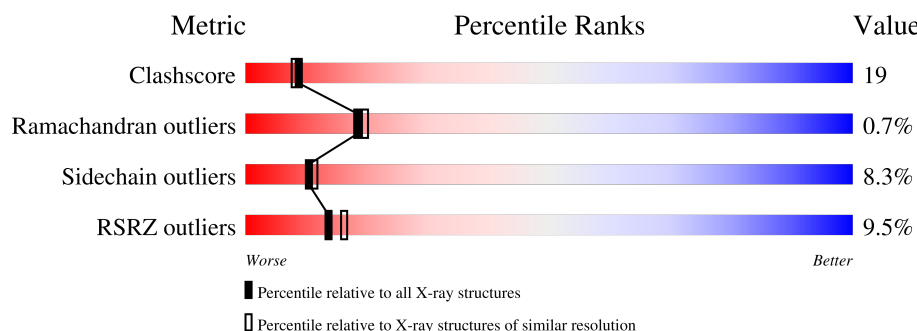
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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	324	
1	B	324	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	MPD	A	328	-	-	X	-
4	TRS	B	328	-	X	-	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5633 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 Type-2 restriction enzyme PvuII.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	313	Total	C	N	O	S	0	1	0
			2572	1670	435	463	4			
1	B	311	Total	C	N	O	S	0	0	0
			2543	1651	428	460	4			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	158	GLY	-	linker	UNP P23657
A	159	SER	-	linker	UNP P23657
A	160	GLY	-	linker	UNP P23657
A	161	GLY	-	linker	UNP P23657
A	318	GLY	-	expression tag	UNP P23657
A	319	SER	-	expression tag	UNP P23657
A	320	HIS	-	expression tag	UNP P23657
A	321	HIS	-	expression tag	UNP P23657
A	322	HIS	-	expression tag	UNP P23657
A	323	HIS	-	expression tag	UNP P23657
A	324	HIS	-	expression tag	UNP P23657
A	325	HIS	-	expression tag	UNP P23657
B	158	GLY	-	linker	UNP P23657
B	159	SER	-	linker	UNP P23657
B	160	GLY	-	linker	UNP P23657
B	161	GLY	-	linker	UNP P23657
B	318	GLY	-	expression tag	UNP P23657
B	319	SER	-	expression tag	UNP P23657
B	320	HIS	-	expression tag	UNP P23657
B	321	HIS	-	expression tag	UNP P23657
B	322	HIS	-	expression tag	UNP P23657
B	323	HIS	-	expression tag	UNP P23657
B	324	HIS	-	expression tag	UNP P23657
B	325	HIS	-	expression tag	UNP P23657

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



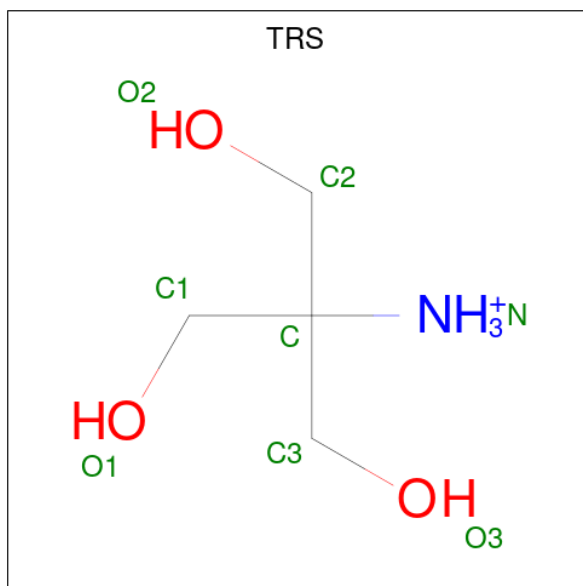
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is (4S)-2-METHYL-2,4-PENTANEDIOL (CCD ID: MPD) (formula: C₆H₁₄O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			8	6	2		

- Molecule 4 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (CCD ID: TRS) (formula: $C_4H_{12}NO_3$).



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

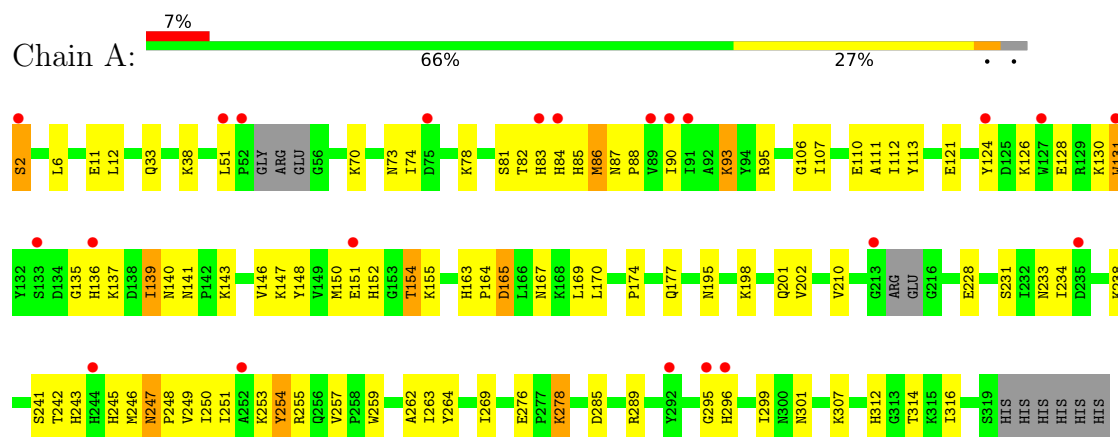
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	274	Total 274	O 274	0	0
5	B	195	Total 195	O 195	0	0

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: Type-2 restriction enzyme PvuII



4 Data and refinement statistics

Property	Value	Source
Space group	P 42	Depositor
Cell constants a, b, c, α , β , γ	101.92Å 101.92Å 100.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	102.06 – 2.35 101.92 – 2.35	Depositor EDS
% Data completeness (in resolution range)	98.6 (102.06-2.35) 98.6 (101.92-2.35)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.95 (at 2.34Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.215 , 0.283 0.215 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	46.0	Xtriage
Anisotropy	0.340	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 47.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.011 for -h,-l,-k 0.000 for -h,l,k 0.000 for l,-k,h 0.003 for -l,-k,-h 0.023 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5633	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.84% 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: MPD, TRS, 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	0.66	0/2650	1.26	7/3595 (0.2%)
1	B	0.68	1/2616 (0.0%)	1.34	15/3553 (0.4%)
All	All	0.67	1/5266 (0.0%)	1.30	22/7148 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	165	ASP	C-N	-5.16	1.27	1.34

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	24	LEU	CA-CB-CG	9.12	148.21	116.30
1	B	58	ASP	N-CA-C	8.42	121.39	111.71
1	A	257	VAL	N-CA-C	7.18	117.06	108.45
1	B	165	ASP	CA-CB-CG	6.86	119.46	112.60
1	B	133	SER	N-CA-C	6.64	118.21	110.97
1	A	165	ASP	CA-CB-CG	6.63	119.23	112.60
1	A	131	TRP	N-CA-C	-6.31	104.32	111.14
1	B	123	TYR	N-CA-C	6.07	119.16	111.69
1	A	254	TYR	N-CA-C	6.05	118.38	111.11
1	B	16	ILE	N-CA-C	-5.83	105.05	110.53
1	B	118	LYS	N-CA-C	5.54	118.84	111.75
1	B	48	LEU	N-CA-C	5.35	118.40	110.48
1	A	143	LYS	N-CA-C	5.28	118.83	109.96
1	B	201	GLN	N-CA-C	-5.26	105.44	111.07
1	A	228	GLU	N-CA-C	-5.25	101.13	109.96
1	B	278	LYS	N-CA-C	5.22	117.71	111.71
1	B	234	ILE	N-CA-C	5.21	118.35	111.17
1	B	314	THR	CB-CA-C	5.20	118.33	109.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	202	VAL	N-CA-C	5.15	115.77	110.36
1	B	146	VAL	CB-CA-C	-5.04	105.26	112.22
1	B	93	LYS	N-CA-C	-5.02	104.77	111.24
1	A	93	LYS	N-CA-C	-5.01	105.73	111.14

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	2572	0	2529	100	0
1	B	2543	0	2484	92	0
2	A	10	0	0	1	0
2	B	15	0	0	0	0
3	A	8	0	14	15	0
4	B	16	0	24	1	0
5	A	274	0	0	14	0
5	B	195	0	0	14	0
All	All	5633	0	5051	191	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 19.

All (191) 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:155:LYS:HE3	3:A:328:MPD:C5	1.69	1.20
1:A:155:LYS:HE3	3:A:328:MPD:H53	1.29	1.09
1:A:155:LYS:CE	3:A:328:MPD:H53	1.87	1.05
1:B:3:HIS:CD2	5:B:379:HOH:O	2.13	1.01
1:B:29:ASN:ND2	5:B:519:HOH:O	1.93	1.01
1:B:29:ASN:ND2	1:B:29:ASN:H	1.54	0.99
1:A:238:LYS:HE3	1:B:49:THR:HG21	1.46	0.98
1:B:49:THR:HG22	5:B:404:HOH:O	1.64	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:81:SER:HB2	1:A:141:ASN:HD21	1.33	0.92
1:A:155:LYS:HE3	3:A:328:MPD:H51	1.47	0.92
1:A:246:MET:HE2	1:A:251:ILE:HG13	1.53	0.91
1:A:111:ALA:HA	3:A:328:MPD:C5	2.01	0.90
1:B:241:SER:HB2	1:B:301:ASN:HD21	1.36	0.90
1:B:274:ARG:HD3	1:B:276:GLU:OE2	1.73	0.89
1:B:29:ASN:HD22	1:B:29:ASN:N	1.68	0.88
1:A:247:ASN:HD22	1:A:249:VAL:H	1.21	0.88
1:B:70:LYS:HD3	1:B:81:SER:O	1.74	0.86
1:B:91:ILE:HG22	1:B:95:ARG:HD2	1.60	0.84
1:B:29:ASN:H	1:B:29:ASN:HD22	0.87	0.82
1:A:198:LYS:HZ2	1:A:233:ASN:HD21	1.28	0.81
1:B:90:ILE:HG13	5:B:496:HOH:O	1.80	0.80
1:B:180:GLN:HE21	1:B:191:ILE:HG12	1.48	0.78
1:A:111:ALA:HA	3:A:328:MPD:C4	2.13	0.78
1:B:147:LYS:N	1:B:147:LYS:HD3	1.98	0.78
1:B:147:LYS:HD3	1:B:147:LYS:H	1.48	0.76
1:A:198:LYS:NZ	1:A:233:ASN:HD21	1.84	0.76
1:A:246:MET:CE	1:A:251:ILE:HG13	2.16	0.75
1:B:242:THR:H	1:B:301:ASN:ND2	1.85	0.75
1:B:236:LEU:HD12	5:B:459:HOH:O	1.86	0.75
1:A:111:ALA:HA	3:A:328:MPD:H51	1.67	0.75
1:A:33:GLN:HG2	5:A:476:HOH:O	1.87	0.74
1:B:115:LEU:HD22	1:B:153:GLY:HA2	1.69	0.74
1:B:3:HIS:HD2	5:B:379:HOH:O	1.56	0.73
1:A:2:SER:N	5:A:350:HOH:O	2.21	0.73
1:A:233:ASN:HB3	5:A:329:HOH:O	1.87	0.73
1:B:242:THR:H	1:B:301:ASN:HD21	1.36	0.73
1:B:29:ASN:ND2	1:B:29:ASN:N	2.32	0.72
1:B:115:LEU:HD22	1:B:153:GLY:CA	2.19	0.71
1:A:246:MET:HE2	1:A:251:ILE:CG1	2.21	0.70
1:A:111:ALA:HA	3:A:328:MPD:H4	1.73	0.70
1:A:163:HIS:HD2	1:A:165:ASP:H	1.40	0.69
1:A:74:ILE:O	1:A:78:LYS:NZ	2.25	0.69
1:A:276:GLU:CB	1:A:278:LYS:HE2	2.23	0.69
1:A:247:ASN:ND2	1:A:249:VAL:H	1.90	0.69
1:A:174:PRO:HA	1:A:177:GLN:HE21	1.58	0.67
1:B:314:THR:CG2	5:B:440:HOH:O	2.41	0.67
1:B:86:MET:HG2	5:B:518:HOH:O	1.93	0.67
1:B:57:ASN:HB3	1:B:66:GLU:HG2	1.75	0.66
1:B:131:TRP:HD1	1:B:132:TYR:CD1	2.13	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:91:ILE:O	1:B:95:ARG:HG3	1.94	0.66
1:B:315:LYS:HE2	1:B:318:GLY:HA3	1.77	0.66
1:A:33:GLN:CG	5:A:476:HOH:O	2.42	0.65
1:B:180:GLN:NE2	1:B:191:ILE:HG12	2.11	0.65
1:B:241:SER:HB2	1:B:301:ASN:ND2	2.10	0.65
1:A:81:SER:HB2	1:A:141:ASN:ND2	2.09	0.65
1:A:285:ASP:OD2	1:A:289:ARG:NH1	2.30	0.64
1:B:180:GLN:HE22	1:B:191:ILE:H	1.44	0.64
1:A:312:HIS:HE1	5:A:349:HOH:O	1.80	0.63
1:A:195:ASN:OD1	1:A:198:LYS:HE3	1.98	0.63
1:A:155:LYS:HE2	3:A:328:MPD:H53	1.79	0.63
1:A:124:TYR:O	1:A:128:GLU:HG2	1.98	0.62
1:A:111:ALA:CB	3:A:328:MPD:H51	2.30	0.62
1:A:86:MET:CE	1:A:90:ILE:CG2	2.78	0.62
1:B:126:LYS:NZ	1:B:126:LYS:HB3	2.12	0.62
1:B:57:ASN:CB	1:B:66:GLU:HG2	2.29	0.62
1:A:111:ALA:CA	3:A:328:MPD:H51	2.31	0.61
1:B:314:THR:HG23	5:B:440:HOH:O	2.00	0.61
1:A:95:ARG:HG2	1:A:124:TYR:CD2	2.36	0.60
1:A:131:TRP:CD1	1:A:137:LYS:O	2.54	0.60
1:B:147:LYS:H	1:B:147:LYS:CD	2.15	0.60
1:A:86:MET:HE3	1:A:90:ILE:HG21	1.83	0.60
1:A:163:HIS:CD2	1:A:165:ASP:H	2.20	0.60
1:A:242:THR:H	1:A:301:ASN:ND2	2.00	0.60
1:A:154:THR:HG23	5:A:602:HOH:O	2.00	0.59
1:A:155:LYS:CE	3:A:328:MPD:C5	2.54	0.58
1:A:82:THR:H	1:A:141:ASN:ND2	2.01	0.58
1:B:91:ILE:O	1:B:95:ARG:CG	2.51	0.58
1:A:238:LYS:CE	1:B:49:THR:HG21	2.30	0.57
1:B:48:LEU:HD21	1:B:100:ILE:CD1	2.34	0.57
1:A:233:ASN:ND2	5:A:329:HOH:O	2.36	0.57
1:A:276:GLU:HB2	1:A:278:LYS:HE2	1.86	0.57
1:A:82:THR:HG22	1:A:83:HIS:N	2.19	0.56
1:A:85:HIS:HD2	5:A:356:HOH:O	1.87	0.56
1:A:147:LYS:HE2	1:A:147:LYS:H	1.70	0.56
1:A:247:ASN:HD22	1:A:249:VAL:N	1.96	0.56
1:B:139:ILE:O	5:B:518:HOH:O	2.17	0.56
1:A:139:ILE:H	1:A:139:ILE:HD12	1.71	0.56
1:A:247:ASN:HB2	1:A:248:PRO:HD2	1.87	0.56
1:A:245:HIS:O	1:A:250:ILE:CD1	2.54	0.56
1:A:110:GLU:O	3:A:328:MPD:H52	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:86:MET:CE	1:A:90:ILE:HG22	2.36	0.55
1:A:276:GLU:HB3	1:A:278:LYS:HE2	1.87	0.55
1:B:299:ILE:HG22	1:B:302:PRO:HD3	1.89	0.55
1:B:84:HIS:HA	1:B:141:ASN:H	1.73	0.54
1:A:202:VAL:HG22	1:A:269:ILE:HD13	1.88	0.54
1:B:139:ILE:H	1:B:139:ILE:HD12	1.72	0.54
1:A:198:LYS:NZ	1:A:233:ASN:ND2	2.55	0.54
1:A:70:LYS:HD2	1:A:81:SER:O	2.08	0.54
1:A:95:ARG:HH22	1:A:128:GLU:CD	2.16	0.54
1:A:246:MET:HE3	1:A:247:ASN:H	1.73	0.54
1:A:87:ASN:HB2	1:A:88:PRO:CD	2.38	0.53
1:B:73:ASN:C	1:B:73:ASN:HD22	2.14	0.53
1:B:141:ASN:HB3	5:B:360:HOH:O	2.07	0.53
1:B:167:ASN:O	1:B:171:GLU:HG3	2.08	0.53
1:B:247:ASN:OD1	1:B:249:VAL:HG13	2.09	0.53
1:B:209:THR:HG22	1:B:220:VAL:HG13	1.91	0.53
1:B:78:LYS:O	1:B:146:VAL:HG23	2.09	0.53
1:B:87:ASN:HB2	1:B:88:PRO:CD	2.39	0.53
1:B:20:GLN:HE21	1:B:31:ILE:HG12	1.75	0.52
1:B:211:LEU:HD13	1:B:220:VAL:HG12	1.90	0.52
1:A:38:LYS:HE2	1:A:107:ILE:HA	1.92	0.51
1:B:180:GLN:NE2	1:B:191:ILE:H	2.08	0.51
1:A:251:ILE:HG22	1:A:255:ARG:HH22	1.76	0.51
1:A:90:ILE:O	1:A:93:LYS:HB3	2.11	0.51
1:A:278:LYS:H	1:A:278:LYS:NZ	2.09	0.50
1:B:80:PHE:HE1	1:B:146:VAL:HG13	1.77	0.50
1:B:80:PHE:O	1:B:81:SER:C	2.55	0.49
1:B:312:HIS:HE1	5:B:332:HOH:O	1.94	0.49
1:B:82:THR:OG1	1:B:142:PRO:HD2	2.12	0.49
1:B:129:ARG:NE	1:B:129:ARG:HA	2.28	0.49
1:A:246:MET:HB3	1:A:299:ILE:HB	1.93	0.49
1:B:131:TRP:HH2	5:B:373:HOH:O	1.95	0.49
1:B:306:VAL:O	1:B:310:MET:HG3	2.12	0.49
1:A:164:PRO:O	1:A:167:ASN:ND2	2.46	0.49
1:B:80:PHE:CD2	1:B:101:PHE:HD2	2.31	0.49
1:A:137:LYS:O	1:A:137:LYS:HD3	2.13	0.48
1:B:25:LYS:HE3	1:B:165:ASP:OD2	2.13	0.48
1:B:91:ILE:O	1:B:95:ARG:CD	2.61	0.48
1:A:295:GLY:O	1:A:296:HIS:HB2	2.12	0.48
1:A:86:MET:CE	1:A:90:ILE:HG21	2.41	0.48
1:A:233:ASN:OD1	5:A:482:HOH:O	2.20	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:122:PHE:CZ	1:B:126:LYS:HD2	2.49	0.48
1:A:139:ILE:HD12	1:A:139:ILE:N	2.28	0.48
2:A:326:SO4:O2	5:A:556:HOH:O	2.20	0.48
1:A:201:GLN:NE2	1:A:269:ILE:HG22	2.30	0.47
1:B:170:LEU:HD23	4:B:329:TRS:H31	1.96	0.47
1:B:147:LYS:N	1:B:147:LYS:CD	2.73	0.47
1:A:73:ASN:HD21	1:A:106:GLY:H	1.63	0.47
1:A:198:LYS:HZ2	1:A:233:ASN:ND2	2.05	0.47
1:A:251:ILE:HG22	1:A:255:ARG:NH2	2.30	0.47
1:A:234:ILE:HB	1:A:263:ILE:HG23	1.98	0.46
1:B:91:ILE:HA	1:B:94:TYR:HB2	1.96	0.46
1:B:91:ILE:HG21	1:B:124:TYR:HB3	1.97	0.46
1:A:146:VAL:O	1:A:150:MET:HG3	2.15	0.46
1:A:241:SER:HB3	1:A:301:ASN:HD21	1.81	0.46
1:B:244:HIS:O	1:B:300:ASN:HA	2.16	0.46
1:A:86:MET:HE1	1:A:90:ILE:HG22	1.96	0.46
1:B:83:HIS:HB3	1:B:86:MET:HB3	1.98	0.46
1:B:91:ILE:HG22	1:B:95:ARG:HH11	1.81	0.45
1:A:247:ASN:HD22	1:A:247:ASN:C	2.24	0.45
1:A:247:ASN:HB2	1:A:248:PRO:CD	2.46	0.45
1:B:125:ASP:O	1:B:129:ARG:HG2	2.17	0.45
1:B:126:LYS:HB3	1:B:126:LYS:HZ3	1.80	0.45
1:B:48:LEU:HD21	1:B:100:ILE:HD12	1.98	0.45
1:B:3:HIS:CE1	1:B:6:LEU:HB2	2.51	0.45
1:B:90:ILE:HG22	1:B:91:ILE:N	2.32	0.45
1:A:113:TYR:CE1	1:A:155:LYS:HB2	2.52	0.44
3:A:328:MPD:H4	3:A:328:MPD:H12	1.66	0.44
1:B:88:PRO:HG2	5:B:480:HOH:O	2.17	0.44
1:A:231:SER:HA	1:A:262:ALA:O	2.18	0.43
1:B:41:GLN:HE21	1:B:109:ILE:HG22	1.83	0.43
1:B:121:GLU:OE2	1:B:121:GLU:HA	2.18	0.43
1:B:274:ARG:CD	1:B:276:GLU:OE2	2.56	0.43
1:A:316:ILE:HD13	1:A:316:ILE:HG21	1.51	0.43
1:B:91:ILE:CG2	1:B:124:TYR:HB3	2.48	0.43
1:A:254:TYR:CD1	1:A:259:TRP:HZ2	2.36	0.43
1:A:169:LEU:C	1:A:169:LEU:HD23	2.43	0.42
1:A:95:ARG:NH2	1:A:128:GLU:CD	2.77	0.42
1:B:129:ARG:HA	1:B:129:ARG:HE	1.84	0.42
1:A:82:THR:H	1:A:141:ASN:HD21	1.67	0.42
1:B:38:LYS:HB3	1:B:104:TYR:CE2	2.55	0.42
1:B:40:LEU:HD22	1:B:44:LEU:HD22	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:76:LEU:HD12	1:B:77:THR:HG22	2.02	0.42
1:B:95:ARG:HG2	1:B:124:TYR:CD1	2.53	0.42
1:B:86:MET:CE	1:B:142:PRO:HG2	2.49	0.42
1:A:152:HIS:HE1	5:A:369:HOH:O	2.03	0.42
1:A:264:TYR:HD2	5:A:482:HOH:O	2.02	0.41
5:A:366:HOH:O	1:B:52:PRO:HG2	2.20	0.41
1:A:242:THR:OG1	1:A:243:HIS:N	2.52	0.41
1:B:23:ALA:HB1	1:B:28:ILE:HB	2.01	0.41
1:A:148:TYR:HA	1:A:151:GLU:HG2	2.02	0.41
1:B:84:HIS:O	1:B:140:ASN:HA	2.20	0.41
1:B:105:ARG:HD2	1:B:110:GLU:OE2	2.21	0.41
1:A:112:ILE:HG13	3:A:328:MPD:H12	2.01	0.41
1:A:238:LYS:HE3	1:B:49:THR:CG2	2.35	0.41
1:A:95:ARG:NH1	1:A:121:GLU:OE2	2.53	0.41
1:A:242:THR:H	1:A:301:ASN:HD21	1.67	0.41
1:A:314:THR:HG23	5:A:578:HOH:O	2.21	0.41
1:B:211:LEU:HD13	1:B:220:VAL:CG1	2.51	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	308/324 (95%)	298 (97%)	8 (3%)	2 (1%)	21	24
1	B	305/324 (94%)	284 (93%)	19 (6%)	2 (1%)	18	20
All	All	613/648 (95%)	582 (95%)	27 (4%)	4 (1%)	18	20

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	135	GLY

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Mol	Chain	Res	Type
1	B	81	SER
1	A	136	HIS
1	B	294	ASP

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	275/284 (97%)	259 (94%)	16 (6%)	18	22
1	B	270/284 (95%)	241 (89%)	29 (11%)	6	6
All	All	545/568 (96%)	500 (92%)	45 (8%)	10	11

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	SER
1	A	6	LEU
1	A	11	GLU
1	A	12	LEU
1	A	51	LEU
1	A	86	MET
1	A	126	LYS
1	A	130	LYS
1	A	139	ILE
1	A	154	THR
1	A	170	LEU
1	A	210	VAL
1	A	247	ASN
1	A	253	LYS
1	A	278	LYS
1	A	307	LYS
1	B	17	GLN
1	B	24	LEU
1	B	29	ASN
1	B	40	LEU

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Mol	Chain	Res	Type
1	B	42	VAL
1	B	44	LEU
1	B	48	LEU
1	B	49	THR
1	B	50	VAL
1	B	70	LYS
1	B	73	ASN
1	B	90	ILE
1	B	115	LEU
1	B	120	LEU
1	B	126	LYS
1	B	139	ILE
1	B	146	VAL
1	B	147	LYS
1	B	149	VAL
1	B	176	ILE
1	B	182	LEU
1	B	199	LEU
1	B	209	THR
1	B	211	LEU
1	B	237	THR
1	B	249	VAL
1	B	269	ILE
1	B	278	LYS
1	B	314	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	57	ASN
1	A	73	ASN
1	A	85	HIS
1	A	163	HIS
1	A	167	ASN
1	A	177	GLN
1	A	180	GLN
1	A	201	GLN
1	A	217	ASN
1	A	222	ASN
1	A	233	ASN
1	A	247	ASN
1	A	256	GLN

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Mol	Chain	Res	Type
1	A	296	HIS
1	A	300	ASN
1	A	301	ASN
1	A	312	HIS
1	B	7	ASN
1	B	20	GLN
1	B	29	ASN
1	B	41	GLN
1	B	73	ASN
1	B	140	ASN
1	B	152	HIS
1	B	180	GLN
1	B	217	ASN
1	B	225	GLN
1	B	244	HIS
1	B	301	ASN
1	B	312	HIS

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](#)

8 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	327	-	4,4,4	0.36	0	6,6,6	0.64	0
2	SO4	A	326	-	4,4,4	0.23	0	6,6,6	0.46	0
2	SO4	B	1	-	4,4,4	0.31	0	6,6,6	0.31	0
2	SO4	B	326	-	4,4,4	0.30	0	6,6,6	0.19	0
3	MPD	A	328	-	7,7,7	0.40	0	9,10,10	1.24	1 (11%)
2	SO4	A	327	-	4,4,4	0.23	0	6,6,6	0.58	0
4	TRS	B	328	-	7,7,7	0.47	0	9,9,9	1.24	1 (11%)
4	TRS	B	329	-	7,7,7	0.36	0	9,9,9	1.17	0

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
3	MPD	A	328	-	-	5/5/5/5	-
4	TRS	B	329	-	-	3/9/9/9	-
4	TRS	B	328	-	-	9/9/9/9	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	328	MPD	O2-C2-CM	2.48	115.73	107.99
4	B	328	TRS	C3-C-C1	2.12	116.31	110.66

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	328	MPD	C1-C2-C3-C4
3	A	328	MPD	O2-C2-C3-C4
4	B	328	TRS	C2-C-C1-O1
4	B	328	TRS	N-C-C1-O1
4	B	329	TRS	C2-C-C1-O1
4	B	329	TRS	C3-C-C1-O1
4	B	328	TRS	C3-C-C1-O1
4	B	328	TRS	N-C-C2-O2
4	B	328	TRS	C2-C-C3-O3
4	B	328	TRS	N-C-C3-O3

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Mol	Chain	Res	Type	Atoms
4	B	329	TRS	N-C-C1-O1
4	B	328	TRS	C1-C-C2-O2
4	B	328	TRS	C1-C-C3-O3
3	A	328	MPD	C2-C3-C4-C5
4	B	328	TRS	C3-C-C2-O2
3	A	328	MPD	C2-C3-C4-O4
3	A	328	MPD	CM-C2-C3-C4

There are no ring outliers.

3 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	326	SO4	1	0
3	A	328	MPD	15	0
4	B	329	TRS	1	0

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	313/324 (96%)	0.42	22 (7%)	22 25	28, 42, 67, 82	1 (0%)
1	B	311/324 (95%)	0.73	37 (11%)	9 10	29, 45, 78, 92	0
All	All	624/648 (96%)	0.57	59 (9%)	14 16	28, 43, 74, 92	1 (0%)

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	131	TRP	4.8
1	B	134	ASP	4.8
1	B	133	SER	4.4
1	B	88	PRO	4.3
1	B	127	TRP	4.1
1	B	86	MET	4.1
1	A	213	GLY	4.0
1	B	90	ILE	3.5
1	B	139	ILE	3.5
1	B	100	ILE	3.4
1	A	131	TRP	3.4
1	B	141	ASN	3.4
1	A	84[A]	HIS	3.4
1	B	89	VAL	3.2
1	A	90	ILE	3.0
1	B	93	LYS	3.0
1	B	292	TYR	3.0
1	B	85	HIS	3.0
1	B	83	HIS	2.9
1	B	137	LYS	2.9
1	B	62	ASN	2.8
1	B	82	THR	2.8
1	A	124	TYR	2.8
1	B	124	TYR	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	97	VAL	2.7
1	B	143	LYS	2.7
1	B	299	ILE	2.6
1	A	133	SER	2.6
1	B	136	HIS	2.6
1	A	292	TYR	2.5
1	A	91	ILE	2.5
1	B	132	TYR	2.5
1	A	127	TRP	2.4
1	B	142	PRO	2.4
1	B	216	GLY	2.4
1	A	52	PRO	2.4
1	B	84	HIS	2.3
1	A	235	ASP	2.3
1	A	83	HIS	2.3
1	B	80	PHE	2.3
1	B	144	ILE	2.2
1	A	295	GLY	2.2
1	B	56	GLY	2.2
1	A	252	ALA	2.2
1	A	2	SER	2.2
1	B	99	TRP	2.2
1	A	75	ASP	2.2
1	A	136	HIS	2.2
1	A	244	HIS	2.2
1	B	120	LEU	2.2
1	A	89	VAL	2.2
1	B	296	HIS	2.1
1	B	2	SER	2.1
1	A	151	GLU	2.1
1	A	51	LEU	2.1
1	A	296	HIS	2.0
1	B	138	ASP	2.0
1	B	63	ALA	2.0
1	B	140	ASN	2.0

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

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	MPD	A	328	8/8	0.73	0.23	49,56,57,60	0
2	SO4	B	1	5/5	0.87	0.16	89,90,90,92	0
2	SO4	B	327	5/5	0.89	0.18	66,66,69,71	0
2	SO4	B	326	5/5	0.91	0.23	73,73,75,76	0
4	TRS	B	328	8/8	0.91	0.11	53,53,55,56	0
2	SO4	A	326	5/5	0.92	0.10	82,82,83,85	0
4	TRS	B	329	8/8	0.93	0.17	62,63,64,64	0
2	SO4	A	327	5/5	0.94	0.08	60,63,66,66	0

6.5 Other polymers [i](#)

There are no such residues in this entry.