



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 19, 2026 – 11:31 AM UTC

PDB ID : 1JVN / pdb_00001jvn
Title : CRYSTAL STRUCTURE OF IMIDAZOLE GLYCEROL PHOSPHATE SYNTHASE: A TUNNEL THROUGH A (BETA/ALPHA)₈ BARREL JOINS TWO ACTIVE SITES
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Deposited on : 2001-08-30
Resolution : 2.10 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

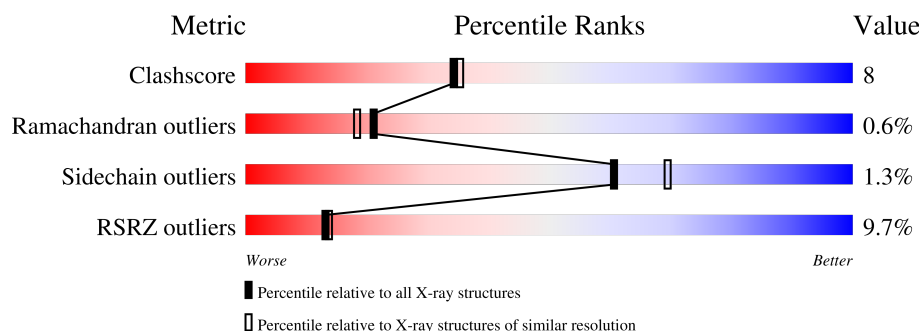
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	555	9% 80% 15% • •
1	B	555	9% 79% 16% • •

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
1	143	A	83	-	-	X	-
4	POP	B	999	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8818 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 BIFUNCTIONAL HISTIDINE BIOSYNTHESIS PROTEIN HISHF.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	537	Total	C	N	O	S	0	0	0
			4185	2659	699	812	15			
1	B	537	Total	C	N	O	S	0	0	0
			4185	2659	699	812	15			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	expression tag	UNP P33734
A	-2	SER	-	expression tag	UNP P33734
A	-1	HIS	-	expression tag	UNP P33734
A	83	143	CYS	modified residue	UNP P33734
B	-3	GLY	-	expression tag	UNP P33734
B	-2	SER	-	expression tag	UNP P33734
B	-1	HIS	-	expression tag	UNP P33734
B	83	143	CYS	modified residue	UNP P33734

- Molecule 2 is NICKEL (II) ION (CCD ID: NI) (formula: Ni).

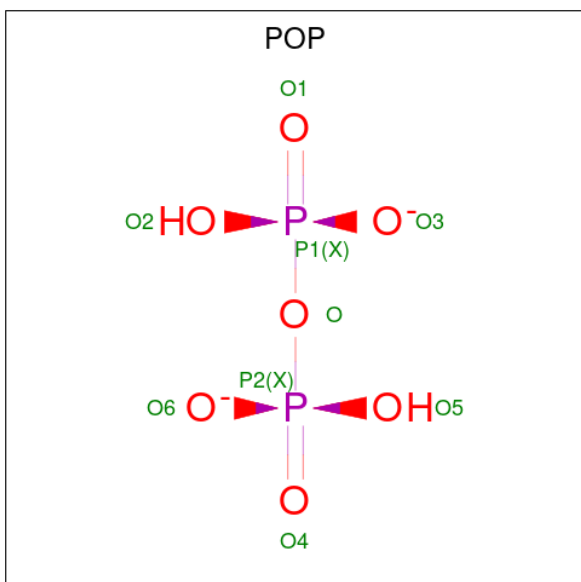
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Ni	0	0
			1	1		
2	B	1	Total	Ni	0	0
			1	1		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 4 is PYROPHOSPHATE 2- (CCD ID: POP) (formula: $\text{H}_2\text{O}_7\text{P}_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	O	P	0	0
			9	7	2		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	212	Total	O	2	0
			212	212		
5	B	200	Total	O	0	0
			200	200		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	98.37Å 112.17Å 116.68Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	75.21 – 2.10 75.21 – 2.10	Depositor EDS
% Data completeness (in resolution range)	97.6 (75.21-2.10) 97.7 (75.21-2.10)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.69 (at 2.10Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.224 , 0.242 0.213 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	28.8	Xtriage
Anisotropy	0.284	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 36.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.010 for -h,l,k	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8818	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.27% 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: 143, POP, NI, 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	0.84	7/4255 (0.2%)	1.05	15/5753 (0.3%)
1	B	0.79	0/4255	1.04	15/5753 (0.3%)
All	All	0.82	7/8510 (0.1%)	1.04	30/11506 (0.3%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	343	LYS	C-N	7.19	1.40	1.33
1	A	344	ILE	N-CA	6.82	1.51	1.46
1	A	404	ASP	C-O	-6.38	1.17	1.24
1	A	305	CYS	C-N	6.08	1.41	1.33
1	A	306	PRO	C-N	-5.49	1.27	1.33
1	A	470	CYS	C-N	-5.34	1.26	1.33
1	A	121	GLU	N-CA	-5.03	1.40	1.46

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	289	GLN	OE1-CD-NE2	-10.60	112.00	122.60
1	A	437	GLN	OE1-CD-NE2	-9.79	112.81	122.60
1	A	288	GLN	OE1-CD-NE2	-9.33	113.27	122.60
1	B	437	GLN	OE1-CD-NE2	-9.13	113.47	122.60
1	B	289	GLN	OE1-CD-NE2	-9.02	113.58	122.60
1	A	344	ILE	N-CA-C	8.26	116.44	107.60
1	B	214	PRO	N-CA-CB	7.22	107.23	103.19
1	A	288	GLN	CG-CD-NE2	6.94	126.81	116.40
1	A	289	GLN	CG-CD-NE2	6.89	126.73	116.40
1	B	289	GLN	CG-CD-NE2	6.87	126.70	116.40
1	A	437	GLN	CG-CD-NE2	6.69	126.43	116.40
1	B	492	VAL	N-CA-C	6.42	118.08	108.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	147	HIS	N-CA-C	6.40	118.80	109.07
1	B	404	ASP	CA-C-N	6.37	126.83	120.52
1	B	404	ASP	C-N-CA	6.37	126.83	120.52
1	B	437	GLN	CG-CD-NE2	6.32	125.88	116.40
1	B	104	TYR	N-CA-C	-6.18	104.65	111.82
1	B	256	VAL	N-CA-C	-5.98	106.98	112.96
1	B	216	ILE	O-C-N	5.83	127.44	121.14
1	A	51	GLY	O-C-N	5.82	128.68	122.86
1	A	119	VAL	CB-CA-C	5.76	115.72	110.13
1	A	305	CYS	CA-C-N	5.71	125.66	119.78
1	A	305	CYS	C-N-CA	5.71	125.66	119.78
1	A	344	ILE	O-C-N	5.65	126.12	120.92
1	A	405	PRO	CB-CA-C	-5.54	103.69	110.95
1	B	491	ALA	N-CA-C	5.32	116.88	111.14
1	A	492	VAL	N-CA-C	5.30	116.48	108.96
1	B	482	LEU	N-CA-C	5.24	116.99	111.28
1	A	148	SER	N-CA-C	-5.09	107.62	113.88
1	B	14	LEU	N-CA-C	-5.07	107.14	113.72

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	4185	0	4113	65	0
1	B	4185	0	4113	65	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	15	0	0	0	0
3	B	10	0	0	0	0
4	B	9	0	0	5	0
5	A	212	0	0	1	0
5	B	200	0	0	3	0
All	All	8818	0	8226	127	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 8.

All (127) 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:B:404:ASP:O	1:B:471:ILE:HD11	1.70	0.91
1:B:541:LEU:HD13	1:B:547:VAL:HG22	1.57	0.87
1:A:403:VAL:HG12	1:A:405:PRO:HD3	1.59	0.84
1:A:541:LEU:HD13	1:A:547:VAL:HG22	1.64	0.79
1:B:475:GLY:HA2	4:B:999:POP:O3	1.83	0.78
1:B:1:MET:HE1	1:B:25:GLY:O	1.89	0.72
1:B:245:ASP:HB3	1:B:299:ILE:HD11	1.72	0.72
1:A:404:ASP:C	1:A:471:ILE:HD11	2.15	0.71
1:B:146:VAL:HG22	1:B:397:GLN:HB2	1.70	0.71
1:A:194:PRO:O	1:A:201:GLY:HA3	1.91	0.71
1:B:403:VAL:HG11	1:B:451:VAL:HG23	1.74	0.68
1:B:524:GLY:HA3	4:B:999:POP:O1	1.94	0.67
1:A:245:ASP:HB3	1:A:299:ILE:HD11	1.76	0.66
1:A:464:GLY:O	1:A:495:PRO:HD2	1.96	0.65
1:B:65:PHE:C	1:B:68:PRO:HD2	2.22	0.65
1:A:403:VAL:HG11	1:A:451:VAL:HG23	1.78	0.64
1:A:307:LEU:HD13	1:A:336:ILE:CD1	2.28	0.64
1:A:403:VAL:HG12	1:A:405:PRO:CD	2.28	0.63
1:A:404:ASP:O	1:A:471:ILE:HD11	1.99	0.62
1:B:121:GLU:OE1	1:B:147:HIS:NE2	2.28	0.60
1:A:2:PRO:HB3	1:B:476:SER:HB2	1.82	0.60
1:B:459:GLU:HG3	1:B:494:ILE:CD1	2.32	0.60
1:B:541:LEU:HD13	1:B:547:VAL:CG2	2.32	0.59
1:B:146:VAL:HG13	1:B:146:VAL:O	2.03	0.59
1:B:1:HIS:CD2	1:B:1:MET:HG3	2.38	0.59
1:B:355:ARG:HD3	5:B:1192:HOH:O	2.02	0.59
1:B:464:GLY:O	1:B:495:PRO:HD2	2.04	0.58
1:B:293:GLU:HG3	1:B:326:PRO:HB2	1.85	0.58
1:B:459:GLU:HG3	1:B:494:ILE:HD13	1.85	0.58
1:A:405:PRO:HB3	1:A:436:TYR:HB3	1.87	0.56
1:A:405:PRO:HD2	1:A:469:ASN:O	2.05	0.56
1:A:80:MET:HA	1:A:188:PHE:O	2.06	0.56
1:A:83:143:OJ1	1:A:149:PHE:N	2.30	0.56
1:B:168:ILE:CD1	1:B:182:VAL:HG12	2.37	0.55
1:A:146:VAL:HG22	1:A:397:GLN:HB2	1.90	0.54
1:A:121:GLU:HB2	1:A:173:TYR:CE2	2.42	0.54
1:A:312:MET:HE3	1:A:315:VAL:HB	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:540:LEU:HD22	1:A:545:LEU:HD12	1.89	0.53
1:B:67:LYS:HB2	1:B:68:PRO:HD3	1.90	0.53
1:B:194:PRO:O	1:B:201:GLY:HA3	2.09	0.53
1:A:83:143:HI	1:A:87:GLN:HE22	1.73	0.53
1:B:121:GLU:HB2	1:B:173:TYR:CE2	2.44	0.53
1:B:524:GLY:CA	4:B:999:POP:O1	2.57	0.52
1:B:524:GLY:HA3	4:B:999:POP:P1	2.50	0.52
1:B:257:THR:HB	1:B:261:ASP:OD2	2.10	0.52
1:A:67:LYS:N	1:A:68:PRO:HD2	2.24	0.52
1:B:168:ILE:HD13	1:B:182:VAL:HG12	1.92	0.51
1:A:470:CYS:H	1:A:500:SER:HB2	1.76	0.51
1:A:521:LEU:C	1:A:521:LEU:HD23	2.35	0.51
1:A:345:PRO:O	1:A:349:VAL:HG23	2.11	0.51
1:B:254:LEU:HD21	1:B:311:PRO:HB2	1.93	0.50
1:A:146:VAL:O	1:A:146:VAL:HG13	2.12	0.50
1:A:247:ARG:NH1	1:A:299:ILE:HD12	2.27	0.50
1:B:521:LEU:C	1:B:521:LEU:HD23	2.36	0.50
1:B:247:ARG:NH1	1:B:299:ILE:HD12	2.26	0.49
1:A:87:GLN:HB3	1:A:109:LEU:HD13	1.95	0.49
1:B:403:VAL:HG12	1:B:405:PRO:HD3	1.94	0.49
1:B:510:GLU:HG3	1:B:514:LYS:HD2	1.93	0.49
1:A:124:TRP:CE3	1:A:548:ARG:HD2	2.48	0.49
1:B:307:LEU:HD22	1:B:336:ILE:HD11	1.95	0.49
1:A:-3:GLY:HA3	1:B:425:TYR:CZ	2.48	0.49
1:A:118:PRO:HA	1:A:493:LYS:O	2.13	0.48
1:B:46:ILE:HG22	1:B:48:PRO:HD3	1.95	0.48
1:B:90:PHE:O	1:B:101:GLY:HA3	2.13	0.48
1:B:420:VAL:HG12	1:B:435:TRP:HB3	1.95	0.48
1:B:146:VAL:CG2	1:B:397:GLN:HB2	2.40	0.48
1:B:527:HIS:HB2	5:B:1198:HOH:O	2.14	0.47
1:A:307:LEU:HD13	1:A:336:ILE:HD13	1.95	0.47
1:A:323:VAL:CG1	1:A:325:VAL:HG22	2.44	0.47
1:A:534:ASN:O	1:A:538:GLU:HG2	2.15	0.47
1:A:276:LEU:C	1:A:276:LEU:HD12	2.40	0.47
1:B:80:MET:HA	1:B:188:PHE:O	2.14	0.47
1:A:83:143:HI	1:A:87:GLN:NE2	2.30	0.47
1:A:2:PRO:HD3	1:B:472:ASP:O	2.16	0.46
1:A:541:LEU:HD13	1:A:547:VAL:CG2	2.42	0.45
1:B:81:GLY:O	1:B:189:ALA:HA	2.16	0.45
1:B:336:ILE:HG13	1:B:344:ILE:HB	1.99	0.45
1:A:83:143:CE	1:A:191:GLN:HE22	2.30	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:459:GLU:HG3	1:B:494:ILE:HG12	1.97	0.45
1:A:506:GLU:HG3	1:A:539:TYR:OH	2.16	0.45
1:A:124:TRP:HB3	1:A:142:ARG:HB3	1.98	0.45
1:B:308:LYS:HG3	5:B:1169:HOH:O	2.17	0.44
1:B:169:ALA:HB3	1:B:181:ALA:HB3	2.00	0.44
1:A:65:PHE:O	1:A:68:PRO:HG2	2.17	0.44
1:B:459:GLU:HG3	1:B:494:ILE:CG1	2.47	0.44
1:A:83:143:SG	1:A:84:VAL:HG23	2.57	0.44
1:B:124:TRP:HB3	1:B:142:ARG:HB3	1.99	0.44
1:B:400:VAL:HG22	1:B:465:GLU:HB3	2.00	0.44
1:A:403:VAL:C	1:A:405:PRO:HD3	2.42	0.44
1:A:257:THR:HG22	1:A:261:ASP:OD2	2.16	0.43
1:A:336:ILE:HG13	1:A:344:ILE:HB	2.00	0.43
1:B:67:LYS:N	1:B:68:PRO:CD	2.80	0.43
1:B:249:ASN:HD22	1:B:261:ASP:C	2.27	0.43
1:B:289:GLN:O	1:B:534:ASN:CG	2.61	0.43
1:A:513:LEU:HD21	1:A:545:LEU:HD21	2.00	0.43
1:A:50:VAL:HG22	5:A:1027:HOH:O	2.17	0.43
1:B:276:LEU:C	1:B:276:LEU:HD12	2.44	0.43
1:A:46:ILE:HG22	1:A:48:PRO:HD3	2.01	0.43
1:A:199:LYS:HE2	1:A:203:ASN:OD1	2.19	0.43
1:A:255:VAL:HG11	1:A:261:ASP:C	2.43	0.43
1:B:124:TRP:CE3	1:B:548:ARG:HD2	2.54	0.43
1:B:524:GLY:HA3	4:B:999:POP:O2	2.19	0.43
1:A:146:VAL:CG2	1:A:397:GLN:HB2	2.49	0.42
1:B:307:LEU:HD13	1:B:336:ILE:CD1	2.49	0.42
1:B:534:ASN:O	1:B:538:GLU:HG2	2.19	0.42
1:A:83:143:NI	1:A:87:GLN:OE1	2.52	0.42
1:B:344:ILE:HA	1:B:345:PRO:HD2	1.91	0.42
1:A:57:VAL:HG11	1:A:100:THR:O	2.20	0.42
1:B:467:LEU:HD11	1:B:521:LEU:HD13	2.02	0.42
1:B:540:LEU:HD22	1:B:545:LEU:HD12	2.01	0.42
1:A:237:THR:HB	1:A:547:VAL:HA	2.02	0.42
1:A:413:GLN:OE1	1:A:420:VAL:HG22	2.19	0.42
1:A:312:MET:HE3	1:A:312:MET:O	2.20	0.41
1:B:413:GLN:OE1	1:B:420:VAL:HG22	2.20	0.41
1:A:65:PHE:HA	1:A:68:PRO:HG2	2.03	0.41
1:A:154:ASN:O	1:A:158:LYS:HG2	2.21	0.41
1:B:307:LEU:CD2	1:B:344:ILE:HD13	2.51	0.41
1:A:403:VAL:HG12	1:A:405:PRO:CG	2.51	0.41
1:B:305:CYS:HA	1:B:306:PRO:HD2	1.77	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:471:ILE:H	1:A:471:ILE:HG13	1.48	0.41
1:B:49:GLY:O	1:B:83:143:HB2	2.21	0.41
1:A:83:143:O	1:A:86:LEU:HB3	2.21	0.40
1:A:90:PHE:O	1:A:101:GLY:HA3	2.21	0.40
1:A:306:PRO:O	1:A:307:LEU:C	2.58	0.40
1:A:504:VAL:HG22	1:A:507:HIS:CE1	2.57	0.40
1:B:311:PRO:O	1:B:315:VAL:HG23	2.21	0.40
1:A:425:TYR:HA	1:A:426:PRO:HD2	1.87	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	530/555 (96%)	516 (97%)	11 (2%)	3 (1%)	21	18
1	B	530/555 (96%)	516 (97%)	11 (2%)	3 (1%)	21	18
All	All	1060/1110 (96%)	1032 (97%)	22 (2%)	6 (1%)	21	18

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	259	LYS
1	A	-2	SER
1	B	-2	SER
1	A	500	SER
1	B	500	SER
1	B	146	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	450/466 (97%)	449 (100%)	1 (0%)	87	93
1	B	450/466 (97%)	439 (98%)	11 (2%)	43	49
All	All	900/932 (97%)	888 (99%)	12 (1%)	61	69

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	248	THR
1	B	50	VAL
1	B	94	VAL
1	B	155	SER
1	B	216	ILE
1	B	247	ARG
1	B	248	THR
1	B	250	ASP
1	B	257	THR
1	B	327	LEU
1	B	411	ASN
1	B	541	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	183	ASN
1	A	289	GLN
1	A	298	ASN
1	A	527	HIS
1	B	55	HIS
1	B	183	ASN
1	B	211	GLN
1	B	212	GLN
1	B	298	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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
1	143	A	83	1	12,16,17	1.45	1 (8%)	5,21,23	2.60	2 (40%)
1	143	B	83	1	12,16,17	1.44	1 (8%)	5,21,23	2.62	2 (40%)

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
1	143	A	83	1	-	4/9/24/26	0/1/1/1
1	143	B	83	1	-	4/9/24/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	83	143	CD-SG	-2.62	1.78	1.82
1	B	83	143	CD-SG	-2.60	1.78	1.82

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	83	143	CI-CF-CE	4.19	138.97	133.64
1	A	83	143	CI-CF-CE	4.16	138.94	133.64
1	B	83	143	OF-CF-CI	-3.63	111.78	116.06
1	A	83	143	OF-CF-CI	-3.62	111.79	116.06

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	83	143	NI-CI-CJ-OJ1
1	A	83	143	OF-CF-CI-CJ
1	A	83	143	CE-CF-CI-CJ
1	B	83	143	CA-CB-SG-CD
1	B	83	143	NI-CI-CJ-OJ1
1	B	83	143	OF-CF-CI-CJ
1	B	83	143	CE-CF-CI-CJ
1	A	83	143	CA-CB-SG-CD

There are no ring outliers.

2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	83	143	7	0
1	B	83	143	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 2 are monoatomic - leaving 6 for Mogul analysis.

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	$\# Z > 2$
4	POP	B	999	-	6,8,8	1.18	0	12,13,13	1.87	5 (41%)
3	SO4	A	806	-	4,4,4	0.39	0	6,6,6	0.09	0
3	SO4	B	803	-	4,4,4	0.32	0	6,6,6	0.09	0
3	SO4	A	801	-	4,4,4	0.39	0	6,6,6	0.12	0
3	SO4	B	804	-	4,4,4	0.35	0	6,6,6	0.12	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SO4	A	802	-	4,4,4	0.38	0	6,6,6	0.17	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
4	POP	B	999	-	-	0/6/6/6	-

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	999	POP	O5-P2-O	-3.54	92.76	104.64
4	B	999	POP	O5-P2-O4	-2.88	99.61	110.83
4	B	999	POP	O-P2-O4	2.88	126.18	111.04
4	B	999	POP	O6-P2-O4	2.49	120.52	110.83
4	B	999	POP	O3-P1-O	2.09	111.64	104.64

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	999	POP	5	0

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	536/555 (96%)	0.59	52 (9%) 13 14	16, 28, 50, 72	0
1	B	536/555 (96%)	0.65	52 (9%) 13 14	17, 29, 49, 72	0
All	All	1072/1110 (96%)	0.62	104 (9%) 13 14	16, 29, 49, 72	0

All (104) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	259	LYS	5.5
1	A	252	GLY	5.5
1	A	261	ASP	5.3
1	B	254	LEU	4.6
1	B	98	LYS	4.5
1	B	216	ILE	4.5
1	B	100	THR	4.1
1	A	214	PRO	4.0
1	B	304	ASP	3.9
1	B	276	LEU	3.9
1	A	247	ARG	3.7
1	B	300	THR	3.7
1	B	261	ASP	3.6
1	A	414	ALA	3.6
1	B	259	LYS	3.5
1	B	248	THR	3.5
1	B	218	ASN	3.5
1	A	304	ASP	3.4
1	B	249	ASN	3.4
1	A	300	THR	3.4
1	A	250	ASP	3.4
1	A	257	THR	3.3
1	A	260	GLY	3.3
1	A	299	ILE	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	471	ILE	3.2
1	A	470	CYS	3.2
1	B	62	ASN	3.2
1	A	339	VAL	3.2
1	B	40	SER	3.1
1	B	257	THR	3.1
1	A	448	ASP	3.1
1	B	342	THR	3.0
1	B	251	GLN	3.0
1	B	529	GLY	3.0
1	B	527	HIS	3.0
1	B	299	ILE	3.0
1	B	308	LYS	3.0
1	A	251	GLN	2.9
1	A	344	ILE	2.9
1	A	341	GLY	2.9
1	A	255	VAL	2.9
1	A	256	VAL	2.9
1	A	336	ILE	2.9
1	B	250	ASP	2.9
1	B	215	PRO	2.9
1	A	248	THR	2.8
1	B	339	VAL	2.8
1	A	100	THR	2.8
1	B	471	ILE	2.8
1	A	306	PRO	2.8
1	B	260	GLY	2.7
1	A	440	ILE	2.7
1	A	254	LEU	2.7
1	B	217	PRO	2.7
1	A	67	LYS	2.6
1	B	306	PRO	2.6
1	A	50	VAL	2.6
1	B	336	ILE	2.6
1	B	380	ARG	2.6
1	A	342	THR	2.6
1	A	222	GLU	2.5
1	A	551	GLU	2.5
1	A	411	ASN	2.5
1	A	343	LYS	2.5
1	B	247	ARG	2.5
1	B	253	ASP	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	222	GLU	2.5
1	B	343	LYS	2.5
1	B	551	GLU	2.5
1	B	214	PRO	2.5
1	B	428	PRO	2.5
1	A	337	VAL	2.4
1	B	337	VAL	2.4
1	A	377	LEU	2.4
1	A	253	ASP	2.4
1	A	216	ILE	2.3
1	A	338	ASP	2.3
1	B	252	GLY	2.3
1	A	276	LEU	2.3
1	A	55	HIS	2.2
1	A	552	GLU	2.2
1	B	256	VAL	2.2
1	B	1	MET	2.2
1	B	541	LEU	2.2
1	A	35	LYS	2.2
1	A	221	ALA	2.2
1	B	307	LEU	2.2
1	B	246	VAL	2.1
1	A	446	SER	2.1
1	A	308	LYS	2.1
1	B	530	GLU	2.1
1	B	48	PRO	2.1
1	B	411	ASN	2.1
1	A	380	ARG	2.1
1	A	475	GLY	2.1
1	B	528	ARG	2.0
1	B	432	LYS	2.0
1	A	249	ASN	2.0
1	A	444	ARG	2.0
1	A	305	CYS	2.0
1	B	438	CYS	2.0
1	B	55	HIS	2.0
1	A	310	THR	2.0
1	B	277	GLY	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	143	A	83	16/17	0.74	0.25	20,25,49,51	0
1	143	B	83	16/17	0.77	0.18	20,25,49,51	0

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
4	POP	B	999	9/9	0.76	0.20	52,55,57,57	0
3	SO4	A	806	5/5	0.77	0.21	74,75,76,76	0
3	SO4	A	802	5/5	0.80	0.15	57,57,58,58	0
3	SO4	A	801	5/5	0.86	0.11	59,60,60,60	0
3	SO4	B	804	5/5	0.90	0.12	51,51,51,51	0
3	SO4	B	803	5/5	0.93	0.13	49,52,52,52	0
2	NI	A	902	1/1	0.99	0.03	26,26,26,26	0
2	NI	B	901	1/1	0.99	0.04	27,27,27,27	0

6.5 Other polymers [i](#)

There are no such residues in this entry.