



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 08:28 AM GMT

PDB ID : 2R4J
Title : Crystal structure of Escherichia coli SeMet substituted Glycerol-3-phosphate
Dehydrogenase in complex with DHAP
Authors : Yeh, J.I.; Du, S.; Chinte, U.
Deposited on : 2007-08-31
Resolution : 1.96 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

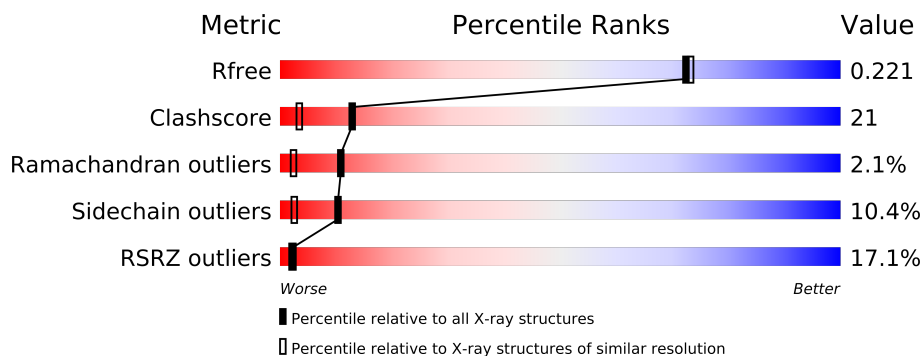
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.96 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}	66092	1321 (1.96-1.96)
Clashscore	79885	1488 (1.96-1.96)
Ramachandran outliers	78287	1475 (1.96-1.96)
Sidechain outliers	78261	1475 (1.96-1.96)
RSRZ outliers	66119	1321 (1.96-1.96)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	501	
1	B	501	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	BOG	A	1949	-	X
2	BOG	A	1950	-	X
2	BOG	A	800	-	X
2	BOG	B	700	-	X
2	BOG	B	800	-	X
3	PO4	B	801	-	X

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Mol	Type	Chain	Res	Geometry	Electron density
5	EDO	A	1952	-	X
5	EDO	A	1953	-	X
5	EDO	A	1955	-	X
5	EDO	A	1956	-	X
5	EDO	A	1957	-	X
5	EDO	A	1958	-	X
5	EDO	A	1965	-	X
5	EDO	B	802	-	X
5	EDO	B	803	-	X
5	EDO	B	806	-	X
5	EDO	B	808	-	X
5	EDO	B	810	-	X
5	EDO	B	811	-	X
5	EDO	B	814	-	X
5	EDO	B	817	-	X
6	IMD	A	1960	-	X
6	IMD	A	1966	-	X
7	TAM	B	812	-	X
9	BCN	A	1969	-	X
9	BCN	B	820	-	X
9	BCN	B	821	-	X

2 Entry composition i

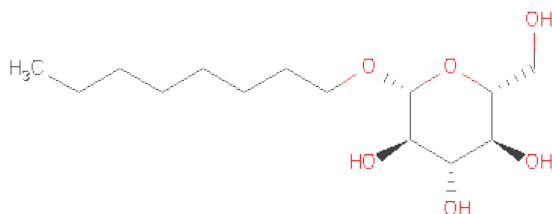
There are 10 unique types of molecules in this entry. The entry contains 8645 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Aerobic glycerol-3-phosphate dehydrogenase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	494	Total	C	N	O	S	Se	0	0	0
			3953	2510	703	727	6	7			
1	B	497	Total	C	N	O	S	Se	0	0	0
			3981	2527	710	731	6	7			

- Molecule 2 is SUGAR (B-OCTYLGLUCOSIDE) (three-letter code: BOG) (formula: C₁₄H₂₈O₆).



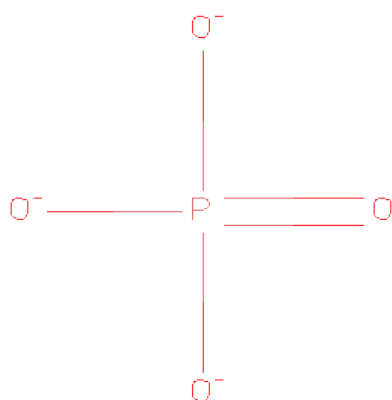
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			20	14	6		
2	A	1	Total	C	O	0	0
			20	14	6		
2	B	1	Total	C	O	0	0
			20	14	6		
2	B	1	Total	C	O	0	0
			20	14	6		

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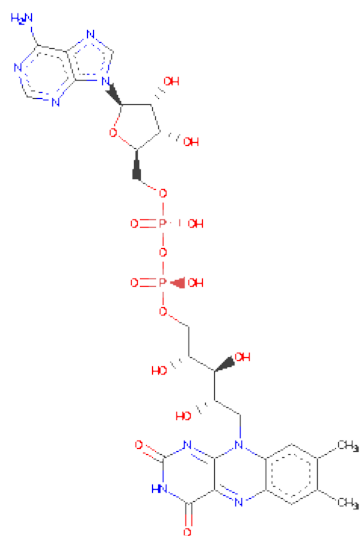
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			20	14	6		
2	A	1	Total	C	O	0	0
			20	14	6		

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



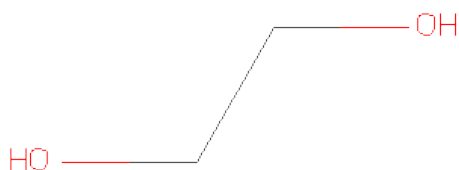
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	O	P	0	0
			5	4	1		
3	A	1	Total	O	P	0	0
			5	4	1		

- Molecule 4 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
4	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



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

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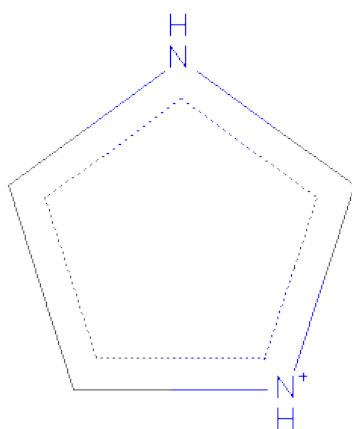
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	A	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	A	1	Total 4	C 2	O 2	0	0
5	A	1	Total 4	C 2	O 2	0	0
5	A	1	Total 4	C 2	O 2	0	0
5	A	1	Total 4	C 2	O 2	0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

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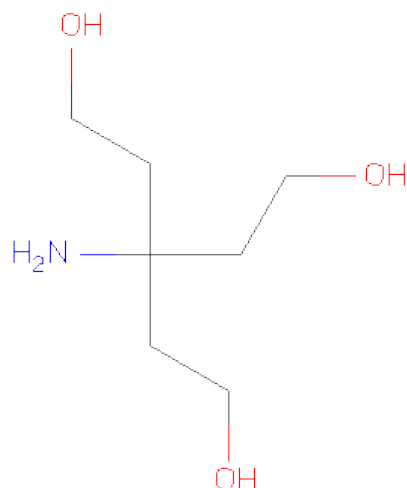
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 6 is IMIDAZOLE (three-letter code: IMD) (formula: $C_3H_5N_2$).



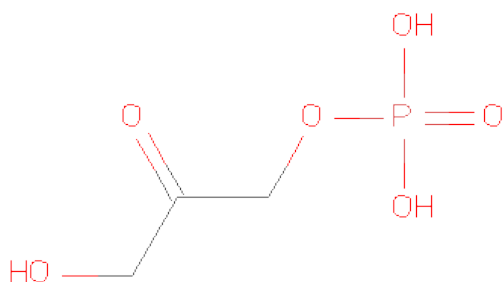
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	N	0	0
			5	3	2		
6	A	1	Total	C	N	0	0
			5	3	2		
6	A	1	Total	C	N	0	0
			5	3	2		
6	A	1	Total	C	N	0	0
			5	3	2		
6	A	1	Total	C	N	0	0
			5	3	2		
6	B	1	Total	C	N	0	0
			5	3	2		

- Molecule 7 is TRIS(HYDROXYETHYL)AMINOMETHANE (three-letter code: TAM) (formula: $C_7H_{17}NO_3$).



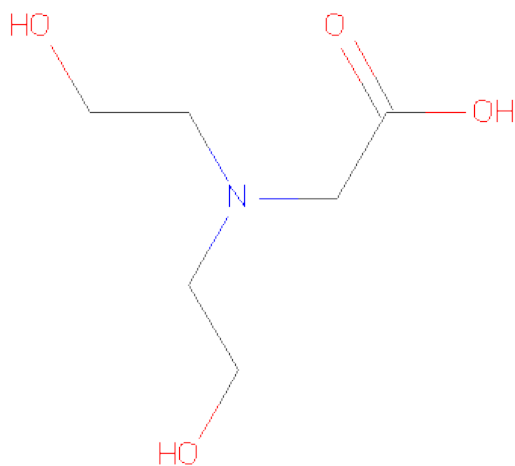
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	B	1	Total	C	N	O	0	0
			11	7	1	3		

- Molecule 8 is 1,3-DIHYDROXYACETONEPHOSPHATE (three-letter code: 13P) (formula: $C_3H_7O_6P$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	A	1	Total	C	O	P	0	0
			10	3	6	1		
8	B	1	Total	C	O	P	0	0
			10	3	6	1		

- Molecule 9 is BICINE (three-letter code: BCN) (formula: $C_6H_{13}NO_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
9	B	1	Total	C	N	O	0	0
			11	6	1	4		
9	B	1	Total	C	N	O	0	0
			11	6	1	4		
9	A	1	Total	C	N	O	0	0
			11	6	1	4		

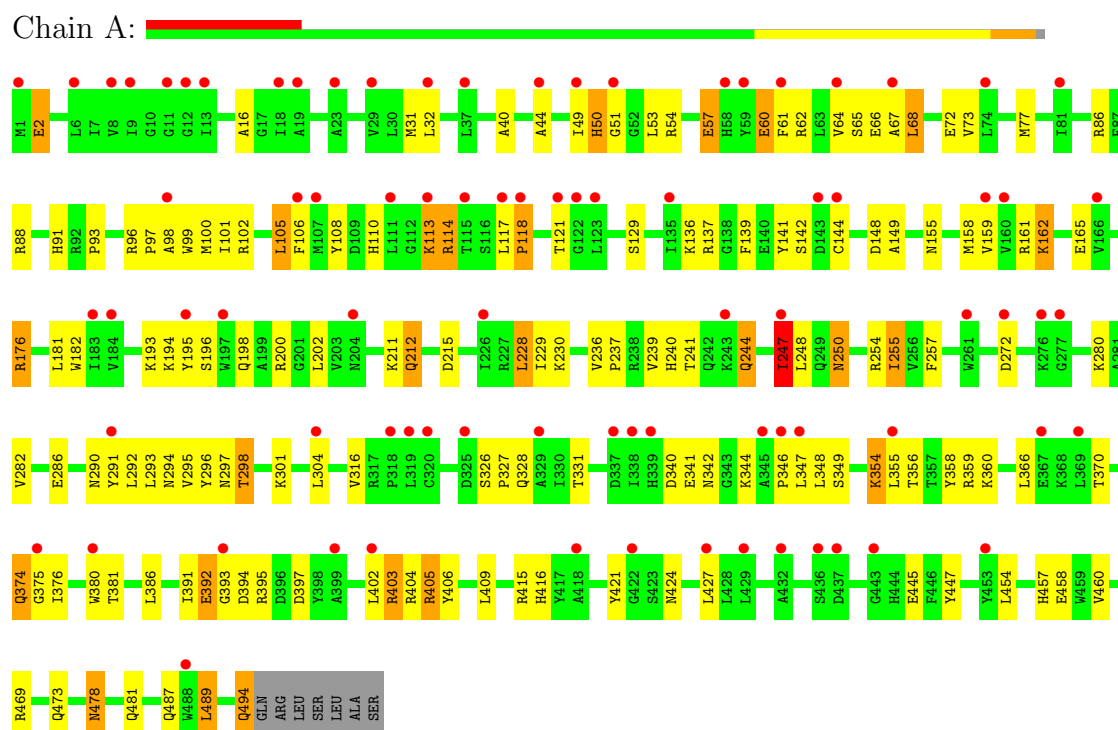
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	140	Total	O	0	0
			140	140		
10	B	137	Total	O	0	0
			137	137		

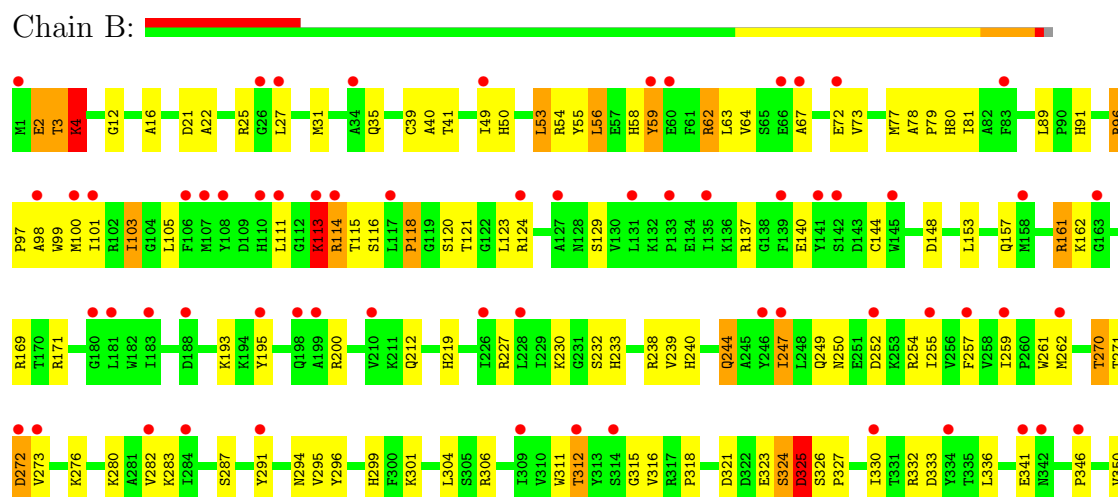
3 Residue-property plots

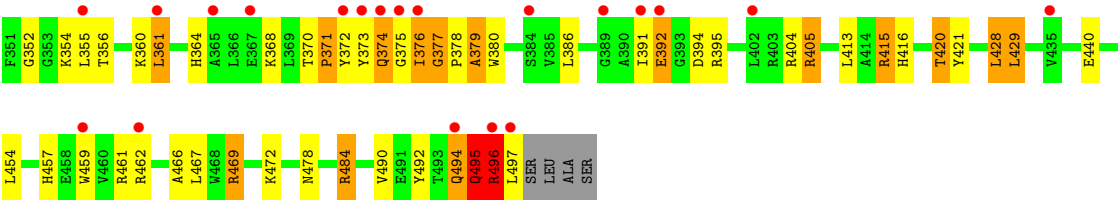
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according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

- Molecule 1: Aerobic glycerol-3-phosphate dehydrogenase



- Molecule 1: Aerobic glycerol-3-phosphate dehydrogenase





4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	113.92Å 114.14Å 193.59Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 1.96 40.32 – 1.96	Depositor EDS
% Data completeness (in resolution range)	100.0 (10.00-1.96) 91.0 (40.32-1.96)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	133.85 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.211 , 0.263 0.206 , 0.221	Depositor DCC
R_{free} test set	4126 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	33.3	Xtriage
Anisotropy	0.369	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 30.8	EDS
Estimated twinning fraction	0.450 for k,h,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.25$	Xtriage
Outliers	0 of 82520 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8645	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.15% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: BCN, IMD, PO4, EDO, 13P, TAM, FAD, BOG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	1.10	4/4041 (0.1%)	1.03	6/5461 (0.1%)
1	B	1.12	2/4069 (0.0%)	1.05	12/5498 (0.2%)
All	All	1.11	6/8110 (0.1%)	1.04	18/10959 (0.2%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	295	VAL	C-O	5.60	1.33	1.23
1	A	447	TYR	CD2-CE2	5.58	1.47	1.39
1	A	165	GLU	CB-CG	5.50	1.62	1.52
1	B	12	GLY	N-CA	-5.27	1.38	1.46
1	A	236	VAL	CB-CG2	5.19	1.63	1.52
1	B	466	ALA	CA-CB	5.07	1.63	1.52

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	169	ARG	NE-CZ-NH2	-5.88	117.36	120.30
1	A	247	ILE	CB-CA-C	-5.74	100.11	111.60
1	A	355	LEU	CB-CG-CD1	-5.71	101.28	111.00
1	B	148	ASP	CB-CG-OD1	5.68	123.41	118.30
1	A	355	LEU	CA-CB-CG	5.60	128.17	115.30
1	B	304	LEU	CB-CG-CD1	-5.58	101.52	111.00
1	B	429	LEU	CA-CB-CG	5.55	128.06	115.30
1	A	86	ARG	NE-CZ-NH1	-5.51	117.54	120.30
1	B	153	LEU	CB-CG-CD1	-5.42	101.79	111.00
1	B	113	LYS	N-CA-C	5.40	125.58	111.00
1	B	96	ARG	NE-CZ-NH2	-5.33	117.64	120.30
1	B	469	ARG	NE-CZ-NH2	-5.30	117.65	120.30
1	A	228	LEU	CA-CB-CG	-5.28	103.16	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	489	LEU	CA-CB-CG	5.28	127.44	115.30
1	B	53	LEU	CB-CG-CD1	-5.19	102.17	111.00
1	B	96	ARG	NE-CZ-NH1	5.13	122.86	120.30
1	B	283	LYS	CB-CA-C	-5.13	100.15	110.40
1	B	462	ARG	NE-CZ-NH1	5.08	122.84	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3953	0	3903	144	0
1	B	3981	0	3935	177	0
2	A	80	0	112	7	0
2	B	40	0	56	9	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
4	A	53	0	31	1	0
4	B	53	0	31	0	0
5	A	44	0	66	28	0
5	B	60	0	90	24	0
6	A	25	0	25	21	0
6	B	5	0	5	1	0
7	B	11	0	17	7	0
8	A	10	0	5	0	0
8	B	10	0	5	4	0
9	A	11	0	12	1	0
9	B	22	0	24	5	0
10	A	140	0	0	9	0
10	B	137	0	0	13	0
All	All	8645	0	8317	347	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including

hydrogens) of the entry. The overall clashscore for this entry is 21.

All (347) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:27:LEU:HD21	1:B:376:ILE:CD1	1.54	1.37
1:B:374:GLN:CB	1:B:375:GLY:HA3	1.59	1.27
1:B:467:LEU:HA	10:B:938:HOH:O	1.37	1.22
5:B:813:EDO:H22	10:B:869:HOH:O	1.37	1.20
1:B:428:LEU:HD12	1:B:428:LEU:O	1.36	1.20
5:A:1959:EDO:H21	6:A:1963:IMD:H5	1.22	1.17
1:A:237:PRO:HG3	6:A:1966:IMD:N3	1.60	1.16
1:B:361:LEU:O	1:B:361:LEU:HD12	1.44	1.14
1:A:51:GLY:HA3	1:A:68:LEU:HD13	1.29	1.12
1:A:374:GLN:CB	1:A:375:GLY:HA3	1.78	1.12
1:A:237:PRO:HG3	6:A:1966:IMD:HN3	1.02	1.11
1:A:341:GLU:CG	1:A:342:ASN:H	1.63	1.11
1:B:374:GLN:HB2	1:B:375:GLY:CA	1.79	1.10
1:A:374:GLN:HB2	1:A:375:GLY:HA3	1.28	1.10
5:A:1957:EDO:H11	6:A:1960:IMD:H5	1.29	1.10
1:A:237:PRO:CG	6:A:1966:IMD:HN3	1.66	1.07
5:B:807:EDO:H11	5:B:808:EDO:O1	1.55	1.06
1:B:50:HIS:CE1	1:B:354:LYS:NZ	2.24	1.06
1:A:341:GLU:HG3	1:A:342:ASN:N	1.65	1.05
1:B:27:LEU:HD21	1:B:376:ILE:HD12	1.13	1.05
1:B:361:LEU:HD12	1:B:361:LEU:C	1.77	1.04
5:A:1959:EDO:H21	6:A:1963:IMD:C5	1.86	1.04
1:B:50:HIS:CE1	1:B:354:LYS:HZ3	1.75	1.03
1:A:158:MSE:CE	1:A:161:ARG:HD2	1.87	1.03
1:B:113:LYS:HB3	1:B:114:ARG:HA	1.41	1.02
1:B:50:HIS:HE1	1:B:354:LYS:NZ	1.58	1.01
1:A:158:MSE:HE1	1:A:161:ARG:HD2	1.02	0.99
1:B:428:LEU:HD12	1:B:428:LEU:C	1.82	0.98
1:A:294:ASN:O	1:A:298:THR:HG23	1.62	0.97
5:A:1959:EDO:C2	6:A:1963:IMD:H5	1.95	0.97
1:A:176:ARG:HH11	1:A:176:ARG:HG2	1.29	0.96
1:A:341:GLU:HG3	1:A:342:ASN:H	0.81	0.95
1:B:62:ARG:HD3	1:B:333:ASP:OD2	1.66	0.95
1:A:158:MSE:HE1	1:A:161:ARG:CD	1.96	0.95
5:A:1957:EDO:C1	6:A:1960:IMD:H5	1.97	0.93
1:B:73:VAL:HG12	1:B:77:MSE:CE	1.96	0.93
1:A:297:ASN:HD21	1:A:304:LEU:H	1.05	0.93
1:A:457:HIS:C	6:A:1963:IMD:H4	1.89	0.93
1:B:467:LEU:HD23	10:B:938:HOH:O	1.68	0.92
1:B:50:HIS:CE1	1:B:354:LYS:CE	2.53	0.91

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:27:LEU:CD2	1:B:376:ILE:HD12	2.00	0.91
1:A:141:TYR:OH	1:A:247:ILE:CD1	2.21	0.88
1:B:376:ILE:HG23	1:B:376:ILE:O	1.70	0.88
1:B:374:GLN:HB2	1:B:375:GLY:HA3	0.88	0.88
1:B:315:GLY:HA2	5:B:813:EDO:H11	1.56	0.88
1:B:73:VAL:HG12	1:B:77:MSE:HE1	1.55	0.87
1:B:27:LEU:CD2	1:B:376:ILE:CD1	2.49	0.85
1:A:51:GLY:HA3	1:A:68:LEU:CD1	2.06	0.84
1:A:176:ARG:CG	1:A:176:ARG:HH11	1.91	0.83
1:B:315:GLY:CA	5:B:813:EDO:H11	2.09	0.83
1:B:212:GLN:HG2	5:B:810:EDO:H22	1.59	0.82
1:A:445:GLU:HB2	6:A:1960:IMD:H2	1.61	0.82
1:A:158:MSE:HA	1:A:158:MSE:HE3	1.61	0.82
1:B:323:GLU:O	1:B:324:SER:HB3	1.80	0.82
1:B:27:LEU:HD11	1:B:376:ILE:HD11	1.60	0.82
1:B:374:GLN:CB	1:B:375:GLY:CA	2.43	0.82
1:A:478:ASN:ND2	1:A:481:GLN:H	1.78	0.80
1:A:487:GLN:HG2	10:A:2079:HOH:O	1.80	0.80
5:B:807:EDO:H11	5:B:808:EDO:HO1	1.47	0.80
1:A:374:GLN:CB	1:A:375:GLY:CA	2.59	0.80
1:B:161:ARG:HB2	7:B:812:TAM:O5	1.81	0.79
1:A:141:TYR:OH	1:A:247:ILE:HD11	1.82	0.79
1:B:50:HIS:HE1	1:B:354:LYS:CE	1.93	0.79
1:B:59:TYR:CE2	2:B:800:BOG:H5'2	2.17	0.79
1:A:392:GLU:CA	1:A:392:GLU:OE1	2.30	0.78
1:B:416:HIS:O	1:B:420:THR:HG23	1.83	0.78
1:A:280:LYS:HA	1:A:280:LYS:HE2	1.65	0.78
1:B:56:LEU:HD13	1:B:64:VAL:HG21	1.66	0.78
1:B:361:LEU:C	1:B:361:LEU:CD1	2.52	0.77
1:B:262:MSE:HE2	5:B:807:EDO:H12	1.67	0.77
1:B:113:LYS:CB	1:B:114:ARG:HA	2.12	0.75
1:A:392:GLU:HA	1:A:392:GLU:OE1	1.84	0.75
1:B:364:HIS:HB2	10:B:885:HOH:O	1.86	0.75
1:A:254:ARG:HD2	1:A:328:GLN:HB3	1.68	0.74
1:B:316:VAL:H	5:B:813:EDO:C1	2.00	0.74
1:B:361:LEU:O	1:B:361:LEU:CD1	2.30	0.74
1:B:157:GLN:HG2	7:B:812:TAM:N	2.01	0.74
1:B:27:LEU:HD21	1:B:376:ILE:HD11	1.67	0.74
1:A:96:ARG:NE	2:A:800:BOG:H5	2.02	0.73
1:B:114:ARG:HG2	1:B:114:ARG:O	1.88	0.73
1:A:374:GLN:HB3	1:A:375:GLY:HA3	1.71	0.72
1:B:324:SER:O	1:B:325:ASP:HB2	1.87	0.72

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:336:LEU:HD12	1:B:364:HIS:HD2	1.52	0.72
1:B:324:SER:OG	1:B:325:ASP:N	2.23	0.72
1:B:376:ILE:CG2	1:B:376:ILE:O	2.38	0.71
1:B:99:TRP:O	1:B:103:ILE:HG12	1.90	0.71
1:A:202:LEU:HD23	1:A:347:LEU:HD13	1.71	0.71
1:A:424:ASN:ND2	6:A:1961:IMD:H4	2.06	0.70
1:B:129:SER:O	1:B:301:LYS:HE2	1.91	0.69
1:A:176:ARG:NH1	1:A:176:ARG:HG2	2.06	0.69
1:B:299:HIS:HE1	10:B:882:HOH:O	1.75	0.68
1:B:50:HIS:CE1	1:B:354:LYS:HE2	2.28	0.68
1:B:113:LYS:HB3	1:B:114:ARG:CA	2.20	0.68
1:B:114:ARG:NH1	1:B:118:PRO:O	2.26	0.68
1:A:53:LEU:HG	2:A:800:BOG:H2	1.76	0.68
1:B:72:GLU:OE2	1:B:114:ARG:HB2	1.93	0.68
1:A:370:THR:HG23	1:A:376:ILE:HG21	1.76	0.68
1:B:378:PRO:O	1:B:379:ALA:O	2.12	0.67
1:A:406:TYR:HE1	10:A:2045:HOH:O	1.76	0.67
5:B:806:EDO:H21	10:B:958:HOH:O	1.95	0.67
1:B:250:ASN:ND2	1:B:291:TYR:CZ	2.63	0.67
1:B:405:ARG:HG3	10:B:894:HOH:O	1.93	0.67
1:B:461:ARG:NH1	10:B:942:HOH:O	2.29	0.66
5:A:1958:EDO:C2	10:A:2097:HOH:O	2.44	0.66
1:B:496:ARG:O	1:B:497:LEU:HG	1.96	0.66
1:B:336:LEU:HD12	1:B:364:HIS:CD2	2.30	0.66
1:A:254:ARG:NE	1:A:272:ASP:OD2	2.27	0.65
1:A:392:GLU:N	1:A:392:GLU:OE1	2.30	0.65
1:B:492:TYR:HB2	6:B:819:IMD:H4	1.79	0.65
1:A:51:GLY:CA	1:A:68:LEU:HD13	2.17	0.65
1:B:2:GLU:O	1:B:4:LYS:HB2	1.97	0.65
1:A:290:ASN:HA	5:A:1956:EDO:H12	1.79	0.64
1:B:301:LYS:NZ	10:B:933:HOH:O	2.29	0.64
1:A:181:LEU:HD23	10:A:2063:HOH:O	1.98	0.64
1:B:428:LEU:CD1	1:B:428:LEU:C	2.58	0.64
1:B:161:ARG:HB2	7:B:812:TAM:HO5	1.63	0.63
1:A:50:HIS:NE2	1:A:354:LYS:NZ	2.44	0.62
1:B:273:VAL:HG13	9:B:820:BCN:H32	1.80	0.62
1:B:428:LEU:CD1	1:B:428:LEU:O	2.30	0.62
1:A:403:ARG:NH1	1:A:409:LEU:O	2.28	0.62
1:A:117:LEU:O	1:A:118:PRO:O	2.17	0.62
1:A:93:PRO:HB3	2:A:1950:BOG:H2	1.82	0.61
1:B:373:TYR:HB2	1:B:376:ILE:HD13	1.82	0.61
1:A:54:ARG:HA	2:A:800:BOG:O6	2.01	0.61

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:287:SER:CB	9:B:821:BCN:H62	2.30	0.61
1:B:73:VAL:CG1	1:B:77:MSE:HE1	2.30	0.61
1:B:239:VAL:HG12	1:B:240:HIS:CD2	2.36	0.60
1:B:440:GLU:OE2	1:B:484:ARG:NH1	2.34	0.60
1:B:421:TYR:OH	1:B:454:LEU:HD21	2.00	0.60
1:B:16:ALA:HA	1:B:31:MSE:HE2	1.82	0.60
1:B:59:TYR:CE2	2:B:800:BOG:C5'	2.84	0.60
1:B:323:GLU:O	1:B:324:SER:CB	2.48	0.59
1:B:316:VAL:H	5:B:813:EDO:H11	1.66	0.59
1:B:370:THR:HG23	1:B:376:ILE:HG21	1.84	0.59
1:A:297:ASN:ND2	1:A:304:LEU:H	1.89	0.59
1:B:262:MSE:HE2	5:B:807:EDO:C1	2.33	0.59
1:A:391:ILE:O	1:A:393:GLY:N	2.30	0.59
1:A:239:VAL:HG23	1:A:240:HIS:CD2	2.38	0.59
1:A:394:ASP:HB3	1:A:397:ASP:HB2	1.84	0.58
1:A:478:ASN:C	1:A:478:ASN:HD22	2.06	0.58
1:A:148:ASP:OD1	1:A:149:ALA:N	2.35	0.58
1:A:290:ASN:CA	5:A:1956:EDO:H12	2.34	0.58
1:A:250:ASN:OD1	1:A:291:TYR:CZ	2.56	0.58
1:A:137:ARG:HH22	2:A:700:BOG:H5	1.68	0.57
1:A:406:TYR:CE1	10:A:2045:HOH:O	2.52	0.57
1:A:141:TYR:CZ	1:A:247:ILE:HD11	2.40	0.57
1:B:316:VAL:N	5:B:813:EDO:H11	2.20	0.57
1:A:341:GLU:CG	1:A:342:ASN:N	2.36	0.57
1:B:378:PRO:O	1:B:379:ALA:C	2.43	0.56
1:B:254:ARG:HH21	8:B:816:13P:H12	1.70	0.56
1:B:73:VAL:CG1	1:B:77:MSE:CE	2.77	0.56
5:A:1958:EDO:H21	10:A:2097:HOH:O	2.04	0.56
1:B:50:HIS:HE1	1:B:354:LYS:HZ1	1.52	0.56
1:B:62:ARG:HG3	1:B:63:LEU:N	2.20	0.56
1:B:405:ARG:NH2	10:B:947:HOH:O	2.37	0.56
1:A:211:LYS:NZ	1:A:215:ASP:OD2	2.37	0.56
1:A:370:THR:HG23	1:A:376:ILE:CG2	2.36	0.55
1:A:454:LEU:HB3	1:A:460:VAL:HG21	1.87	0.55
5:B:804:EDO:H21	10:B:924:HOH:O	2.04	0.55
1:B:262:MSE:HE1	5:B:807:EDO:O2	2.07	0.55
1:B:103:ILE:CD1	2:B:700:BOG:H6'2	2.37	0.55
1:A:101:ILE:HG22	1:A:105:LEU:HD22	1.88	0.55
1:A:473:GLN:NE2	5:A:1957:EDO:H21	2.22	0.55
1:B:262:MSE:CE	5:B:807:EDO:H12	2.36	0.55
5:B:811:EDO:H11	7:B:812:TAM:H62	1.88	0.54
1:A:72:GLU:OE2	1:A:114:ARG:HB2	2.07	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:373:TYR:CB	1:B:376:ILE:HD13	2.36	0.54
1:A:57:GLU:HB2	2:A:800:BOG:O6	2.08	0.54
1:A:97:PRO:HG2	1:A:100:MET:HB2	1.90	0.54
1:A:469:ARG:NH1	5:A:1953:EDO:O2	2.40	0.54
1:A:286:GLU:HA	1:A:286:GLU:OE1	2.07	0.54
1:A:230:LYS:HD3	1:A:282:VAL:HG23	1.87	0.54
1:A:494:GLN:HA	1:A:494:GLN:OE1	2.06	0.54
1:B:25:ARG:HD3	1:B:377:GLY:O	2.08	0.54
1:B:157:GLN:HG2	7:B:812:TAM:HN2	1.70	0.54
1:A:193:LYS:HD2	1:A:195:TYR:CZ	2.43	0.53
1:A:155:ASN:O	1:A:159:VAL:HG23	2.09	0.53
1:B:376:ILE:O	1:B:377:GLY:O	2.28	0.52
1:A:374:GLN:HB2	1:A:375:GLY:CA	2.19	0.52
1:B:101:ILE:HG23	2:B:800:BOG:H61	1.92	0.52
1:A:229:ILE:HD12	1:A:327:PRO:HB3	1.92	0.52
1:B:97:PRO:HD2	1:B:100:MET:SD	2.50	0.52
1:B:232:SER:O	1:B:270:THR:HG22	2.10	0.52
1:B:120:SER:HA	1:B:140:GLU:O	2.10	0.52
1:A:97:PRO:HB2	1:A:99:TRP:CD1	2.45	0.51
1:A:50:HIS:NE2	1:A:354:LYS:CE	2.74	0.51
1:A:391:ILE:C	1:A:393:GLY:H	2.13	0.51
1:B:27:LEU:CD1	1:B:376:ILE:HD11	2.38	0.51
1:A:54:ARG:HG2	10:A:2109:HOH:O	2.11	0.51
1:A:64:VAL:HG12	1:A:68:LEU:HD22	1.91	0.51
1:B:58:HIS:O	1:B:59:TYR:O	2.27	0.51
1:A:340:ASP:HB2	1:A:344:LYS:O	2.11	0.50
1:A:129:SER:O	1:A:301:LYS:HE2	2.12	0.50
1:A:114:ARG:HG2	1:A:114:ARG:O	2.09	0.50
1:A:40:ALA:O	5:A:1965:EDO:O1	2.28	0.50
1:B:97:PRO:HG2	1:B:100:MET:HB2	1.92	0.50
5:A:1958:EDO:H22	10:A:2097:HOH:O	2.07	0.50
1:B:272:ASP:HB2	5:B:802:EDO:H21	1.94	0.50
5:A:1953:EDO:H12	5:A:1965:EDO:O2	2.12	0.50
1:B:490:VAL:O	1:B:494:GLN:HB2	2.10	0.50
1:B:324:SER:O	1:B:325:ASP:CB	2.58	0.50
1:A:478:ASN:HD21	1:A:481:GLN:H	1.55	0.49
1:B:27:LEU:HD21	1:B:376:ILE:HD13	1.77	0.49
1:A:478:ASN:HD22	1:A:481:GLN:H	1.57	0.49
1:B:227:ARG:HG2	1:B:321:ASP:HB2	1.93	0.49
1:A:200:ARG:O	1:A:346:PRO:HD2	2.12	0.49
5:A:1959:EDO:H11	6:A:1961:IMD:H2	1.94	0.49
1:B:376:ILE:O	1:B:377:GLY:C	2.49	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:49:ILE:HB	1:B:144:CYS:HB2	1.94	0.49
1:B:252:ASP:OD1	1:B:254:ARG:HB2	2.12	0.49
1:A:294:ASN:ND2	5:A:1955:EDO:O2	2.46	0.49
1:A:193:LYS:HE3	1:A:195:TYR:OH	2.12	0.49
1:A:424:ASN:HD21	6:A:1961:IMD:H4	1.75	0.49
1:A:237:PRO:CG	6:A:1966:IMD:N3	2.43	0.49
1:B:212:GLN:CG	5:B:810:EDO:H22	2.37	0.49
1:B:391:ILE:C	1:B:392:GLU:HG2	2.32	0.49
1:A:96:ARG:HE	2:A:800:BOG:H5	1.78	0.49
1:A:316:VAL:HG22	5:A:1958:EDO:H11	1.95	0.49
1:B:287:SER:HB3	9:B:821:BCN:H62	1.95	0.49
1:B:39:CYS:HA	1:B:469:ARG:HD3	1.95	0.49
1:B:2:GLU:O	1:B:4:LYS:HD2	2.13	0.48
1:B:336:LEU:CD2	1:B:350:VAL:HG22	2.43	0.48
1:B:311:TRP:O	1:B:312:THR:HG22	2.12	0.48
1:B:67:ALA:HB1	1:B:356:THR:HG21	1.94	0.48
1:A:416:HIS:ND1	5:A:1957:EDO:H11	2.27	0.48
1:B:59:TYR:HE2	2:B:800:BOG:C5'	2.25	0.48
1:B:459:TRP:CE2	5:B:811:EDO:H12	2.49	0.48
1:A:457:HIS:CA	6:A:1963:IMD:H4	2.44	0.48
1:B:457:HIS:C	5:B:811:EDO:H21	2.34	0.48
1:A:16:ALA:HA	1:A:31:MSE:HE2	1.95	0.48
1:B:336:LEU:HD22	1:B:350:VAL:HG22	1.95	0.47
1:B:54:ARG:HA	2:B:800:BOG:O2	2.14	0.47
1:A:421:TYR:OH	1:A:454:LEU:HD21	2.14	0.47
1:A:469:ARG:HG3	1:A:469:ARG:NH1	2.28	0.47
1:A:254:ARG:NH1	1:A:254:ARG:HB3	2.30	0.47
1:B:3:THR:HA	1:B:4:LYS:HB2	1.97	0.47
1:A:106:PHE:HE2	5:A:1964:EDO:HO1	1.63	0.47
1:A:114:ARG:NH2	1:A:118:PRO:O	2.47	0.47
1:A:158:MSE:HE3	1:A:161:ARG:HB3	1.97	0.47
1:A:241:THR:O	1:A:241:THR:CG2	2.63	0.47
8:B:816:13P:O3P	8:B:816:13P:H31	2.15	0.47
1:B:193:LYS:HD3	1:B:195:TYR:CZ	2.49	0.47
1:B:478:ASN:C	1:B:478:ASN:OD1	2.52	0.47
1:A:88:ARG:NE	1:A:244:GLN:HG3	2.30	0.47
1:A:255:ILE:HG13	1:A:255:ILE:H	1.58	0.46
1:A:67:ALA:HB1	1:A:356:THR:HG21	1.95	0.46
1:B:457:HIS:HD2	7:B:812:TAM:O6	1.98	0.46
1:B:59:TYR:HE2	2:B:800:BOG:H5'2	1.73	0.46
1:A:458:GLU:N	6:A:1963:IMD:H4	2.29	0.46
5:A:1957:EDO:H12	6:A:1960:IMD:H5	1.93	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
9:A:1969:BCN:H62	9:A:1969:BCN:H11	1.55	0.46
1:B:415:ARG:HH11	1:B:415:ARG:HG2	1.80	0.46
1:A:298:THR:HG22	5:A:1955:EDO:O1	2.15	0.46
1:A:293:LEU:HB2	5:A:1956:EDO:H21	1.98	0.46
1:B:311:TRP:C	1:B:312:THR:CG2	2.84	0.46
1:B:296:TYR:C	1:B:296:TYR:CD2	2.89	0.46
1:B:415:ARG:NH1	1:B:415:ARG:HG2	2.31	0.45
1:A:296:TYR:CD1	1:A:296:TYR:C	2.89	0.45
1:B:96:ARG:HH11	1:B:249:GLN:HE21	1.65	0.45
1:A:416:HIS:ND1	5:A:1957:EDO:C1	2.80	0.45
1:A:106:PHE:O	1:A:110:HIS:CD2	2.70	0.45
1:B:497:LEU:HA	1:B:497:LEU:HD23	1.85	0.45
1:B:200:ARG:O	1:B:346:PRO:HD2	2.16	0.45
4:A:600:FAD:H4B	4:A:600:FAD:O2A	2.17	0.45
9:B:821:BCN:O4	9:B:821:BCN:H12	2.17	0.45
1:A:469:ARG:HH11	1:A:469:ARG:HG3	1.82	0.45
1:B:370:THR:O	1:B:372:TYR:N	2.50	0.45
5:A:1959:EDO:C1	6:A:1961:IMD:H2	2.47	0.45
1:B:364:HIS:NE2	1:B:368:LYS:NZ	2.65	0.45
1:B:73:VAL:O	1:B:77:MSE:HE3	2.16	0.45
1:A:106:PHE:O	1:A:110:HIS:HD2	2.00	0.45
1:B:244:GLN:HE21	1:B:244:GLN:HB2	1.61	0.45
1:B:113:LYS:HE3	1:B:113:LYS:H	1.82	0.44
1:A:254:ARG:HD2	1:A:328:GLN:CB	2.42	0.44
1:B:232:SER:O	1:B:270:THR:CG2	2.66	0.44
1:B:21:ASP:O	1:B:22:ALA:C	2.56	0.44
1:A:359:ARG:NH1	1:A:381:THR:OG1	2.45	0.44
1:B:336:LEU:HD21	1:B:361:LEU:CD1	2.47	0.44
1:B:103:ILE:HD13	2:B:700:BOG:H6'2	1.99	0.44
1:B:77:MSE:HG2	1:B:386:LEU:HG	1.98	0.44
1:B:89:LEU:HD12	1:B:247:ILE:O	2.17	0.44
1:B:262:MSE:CE	5:B:807:EDO:C1	2.93	0.44
1:B:59:TYR:HE2	2:B:800:BOG:C6'	2.31	0.44
1:B:227:ARG:HG2	1:B:321:ASP:CB	2.48	0.44
1:B:250:ASN:ND2	1:B:291:TYR:CE1	2.86	0.44
1:B:261:TRP:CZ2	1:B:472:LYS:HD2	2.53	0.44
1:B:294:ASN:ND2	5:B:803:EDO:O1	2.51	0.44
1:B:254:ARG:HH21	8:B:816:13P:C1	2.30	0.43
1:A:77:MSE:HE2	1:A:386:LEU:HG	2.00	0.43
1:B:233:HIS:CE1	1:B:270:THR:HG23	2.53	0.43
1:A:117:LEU:HD22	1:A:142:SER:HB3	2.00	0.43
1:B:3:THR:HA	1:B:4:LYS:CB	2.49	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
5:B:803:EDO:H21	5:B:817:EDO:H11	2.01	0.43
1:B:323:GLU:CB	10:B:912:HOH:O	2.67	0.42
1:A:366:LEU:HA	1:A:366:LEU:HD13	1.92	0.42
1:A:91:HIS:CE1	1:A:98:ALA:HB2	2.54	0.42
1:A:247:ILE:HG23	1:A:257:PHE:CE1	2.54	0.42
1:B:295:VAL:O	1:B:299:HIS:HD2	2.02	0.42
9:B:820:BCN:H62	9:B:820:BCN:H31	1.64	0.42
1:B:55:TYR:OH	8:B:816:13P:O1P	2.24	0.42
1:A:44:ALA:HB2	5:A:1965:EDO:C2	2.49	0.42
1:A:158:MSE:CA	1:A:158:MSE:HE3	2.41	0.42
1:B:232:SER:N	1:B:270:THR:HG22	2.34	0.42
1:A:248:LEU:O	1:A:255:ILE:HA	2.19	0.42
1:A:139:PHE:N	1:A:139:PHE:CD2	2.87	0.42
1:B:259:ILE:HD13	1:B:259:ILE:HG21	1.83	0.42
1:B:370:THR:O	1:B:371:PRO:C	2.57	0.42
1:A:228:LEU:HD23	1:A:228:LEU:HA	1.79	0.42
1:A:457:HIS:HA	6:A:1963:IMD:C4	2.50	0.42
1:A:374:GLN:HB3	1:A:375:GLY:CA	2.39	0.42
1:A:415:ARG:HD3	10:A:2077:HOH:O	2.18	0.42
1:A:358:TYR:C	1:A:358:TYR:CD2	2.93	0.42
5:A:1959:EDO:O1	6:A:1961:IMD:N1	2.53	0.42
1:B:440:GLU:CD	1:B:484:ARG:NH1	2.73	0.42
1:A:61:PHE:O	1:A:62:ARG:C	2.57	0.42
1:A:212:GLN:HB3	1:A:212:GLN:HE21	1.71	0.42
1:B:80:HIS:CD2	1:B:81:ILE:HG23	2.54	0.42
1:B:157:GLN:NE2	7:B:812:TAM:H41	2.35	0.42
1:B:91:HIS:NE2	1:B:98:ALA:HB2	2.35	0.42
1:A:241:THR:HG22	1:A:241:THR:O	2.20	0.42
1:A:68:LEU:HD11	1:A:108:TYR:CE2	2.55	0.41
1:B:495:GLN:C	1:B:497:LEU:H	2.22	0.41
1:B:270:THR:HB	1:B:271:THR:H	1.40	0.41
1:A:292:LEU:HA	1:A:292:LEU:HD23	1.77	0.41
5:A:1959:EDO:H21	6:A:1963:IMD:C4	2.46	0.41
1:B:77:MSE:HE3	1:B:77:MSE:HB2	1.96	0.41
1:B:257:PHE:HB3	1:B:259:ILE:HG13	2.03	0.41
1:A:66:GLU:HG3	1:A:360:LYS:HD2	2.02	0.41
1:A:405:ARG:HD3	5:A:1962:EDO:O2	2.20	0.41
1:A:254:ARG:HB3	1:A:254:ARG:CZ	2.51	0.41
1:A:158:MSE:O	1:A:162:LYS:HG3	2.20	0.41
1:B:171:ARG:HD2	5:B:815:EDO:O2	2.21	0.41
1:B:58:HIS:O	1:B:59:TYR:C	2.59	0.41
1:B:378:PRO:C	1:B:379:ALA:O	2.59	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:40:ALA:O	1:B:41:THR:C	2.58	0.41
1:B:326:SER:HA	1:B:327:PRO:HD3	1.89	0.41
1:B:416:HIS:CE1	1:B:420:THR:HG21	2.56	0.41
1:B:299:HIS:CE1	10:B:882:HOH:O	2.61	0.41
1:A:73:VAL:O	1:A:77:MSE:HG3	2.21	0.41
1:B:230:LYS:HD3	1:B:282:VAL:HG23	2.03	0.41
1:B:327:PRO:HA	1:B:330:ILE:HD12	2.02	0.41
1:B:318:PRO:O	1:B:352:GLY:HA2	2.21	0.41
1:A:176:ARG:CG	1:A:176:ARG:NH1	2.61	0.40
1:A:60:GLU:OE1	1:A:331:THR:OG1	2.31	0.40
1:A:49:ILE:HB	1:A:144:CYS:HB2	2.03	0.40
1:B:78:ALA:N	1:B:79:PRO:CD	2.85	0.40
1:A:182:TRP:O	1:A:198:GLN:HA	2.21	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	492/501 (98%)	456 (93%)	30 (6%)	6 (1%)	19	6
1	B	495/501 (99%)	454 (92%)	26 (5%)	15 (3%)	7	1
All	All	987/1002 (98%)	910 (92%)	56 (6%)	21 (2%)	11	2

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	113	LYS
1	A	118	PRO
1	B	113	LYS
1	B	118	PRO
1	B	324	SER
1	B	325	ASP
1	B	379	ALA

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Mol	Chain	Res	Type
1	B	59	TYR
1	B	495	GLN
1	B	496	ARG
1	A	374	GLN
1	B	380	TRP
1	A	2	GLU
1	A	380	TRP
1	B	4	LYS
1	B	115	THR
1	B	374	GLN
1	A	354	LYS
1	B	371	PRO
1	B	377	GLY
1	B	376	ILE

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	413/412 (100%)	377 (91%)	36 (9%)	15 4
1	B	416/412 (101%)	366 (88%)	50 (12%)	7 1
All	All	829/824 (101%)	743 (90%)	86 (10%)	10 2

All (86) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	GLU
1	A	32	LEU
1	A	50	HIS
1	A	57	GLU
1	A	60	GLU
1	A	65	SER
1	A	68	LEU
1	A	102	ARG
1	A	105	LEU
1	A	113	LYS
1	A	114	ARG

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Mol	Chain	Res	Type
1	A	121	THR
1	A	136	LYS
1	A	162	LYS
1	A	176	ARG
1	A	194	LYS
1	A	196	SER
1	A	212	GLN
1	A	244	GLN
1	A	247	ILE
1	A	250	ASN
1	A	255	ILE
1	A	298	THR
1	A	326	SER
1	A	348	LEU
1	A	349	SER
1	A	392	GLU
1	A	395	ARG
1	A	402	LEU
1	A	403	ARG
1	A	404	ARG
1	A	405	ARG
1	A	427	LEU
1	A	478	ASN
1	A	489	LEU
1	A	494	GLN
1	B	2	GLU
1	B	3	THR
1	B	4	LYS
1	B	35	GLN
1	B	53	LEU
1	B	56	LEU
1	B	62	ARG
1	B	103	ILE
1	B	105	LEU
1	B	111	LEU
1	B	113	LYS
1	B	114	ARG
1	B	116	SER
1	B	121	THR
1	B	123	LEU
1	B	124	ARG
1	B	137	ARG

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Mol	Chain	Res	Type
1	B	161	ARG
1	B	162	LYS
1	B	219	HIS
1	B	238	ARG
1	B	244	GLN
1	B	247	ILE
1	B	255	ILE
1	B	270	THR
1	B	272	ASP
1	B	276	LYS
1	B	280	LYS
1	B	306	ARG
1	B	312	THR
1	B	325	ASP
1	B	332	ARG
1	B	341	GLU
1	B	355	LEU
1	B	360	LYS
1	B	361	LEU
1	B	392	GLU
1	B	394	ASP
1	B	395	ARG
1	B	404	ARG
1	B	405	ARG
1	B	413	LEU
1	B	415	ARG
1	B	420	THR
1	B	428	LEU
1	B	429	LEU
1	B	484	ARG
1	B	494	GLN
1	B	495	GLN
1	B	496	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	14	ASN
1	A	110	HIS
1	A	128	ASN
1	A	212	GLN
1	A	242	GLN

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Mol	Chain	Res	Type
1	A	244	GLN
1	A	250	ASN
1	A	290	ASN
1	A	297	ASN
1	A	364	HIS
1	A	424	ASN
1	A	444	HIS
1	A	473	GLN
1	A	478	ASN
1	A	482	GLN
1	B	50	HIS
1	B	244	GLN
1	B	249	GLN
1	B	294	ASN
1	B	299	HIS
1	B	424	ASN
1	B	457	HIS
1	B	494	GLN
1	B	495	GLN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

48 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	BOG	A	1949	-	20,20,20	0.94	1 (5%)	25,25,25	0.98	2 (8%)
2	BOG	A	1950	-	20,20,20	0.92	1 (5%)	25,25,25	1.72	5 (20%)
3	PO4	A	1951	-	4,4,4	0.12	0	6,6,6	0.32	0
5	EDO	A	1952	-	3,3,3	0.79	0	2,2,2	0.26	0
5	EDO	A	1953	-	3,3,3	0.69	0	2,2,2	0.96	0
5	EDO	A	1954	-	3,3,3	0.65	0	2,2,2	0.28	0
5	EDO	A	1955	-	3,3,3	0.77	0	2,2,2	0.39	0
5	EDO	A	1956	-	3,3,3	0.58	0	2,2,2	0.24	0
5	EDO	A	1957	-	3,3,3	1.17	0	2,2,2	0.88	0
5	EDO	A	1958	-	3,3,3	0.54	0	2,2,2	0.24	0
5	EDO	A	1959	-	3,3,3	0.76	0	2,2,2	0.52	0
6	IMD	A	1960	-	5,5,5	1.63	2 (40%)	5,5,5	1.47	0
6	IMD	A	1961	-	5,5,5	1.60	2 (40%)	5,5,5	0.59	0
5	EDO	A	1962	-	3,3,3	0.63	0	2,2,2	0.36	0
6	IMD	A	1963	-	5,5,5	1.52	2 (40%)	5,5,5	0.73	0
5	EDO	A	1964	-	3,3,3	0.65	0	2,2,2	0.47	0
5	EDO	A	1965	-	3,3,3	0.73	0	2,2,2	0.68	0
6	IMD	A	1966	-	5,5,5	1.59	2 (40%)	5,5,5	1.33	1 (20%)
6	IMD	A	1967	-	5,5,5	1.50	2 (40%)	5,5,5	1.21	0
8	13P	A	1968	-	9,9,9	5.23	6 (66%)	12,12,12	4.47	6 (50%)
9	BCN	A	1969	-	10,10,10	0.93	0	11,11,11	1.98	4 (36%)
4	FAD	A	600	-	58,58,58	1.76	19 (32%)	85,89,89	1.87	16 (18%)
2	BOG	A	700	-	20,20,20	0.74	1 (5%)	25,25,25	1.18	2 (8%)
2	BOG	A	800	-	20,20,20	0.61	0	25,25,25	1.56	5 (20%)
4	FAD	B	600	-	58,58,58	1.82	12 (20%)	85,89,89	2.18	20 (23%)
2	BOG	B	700	-	20,20,20	0.77	1 (5%)	25,25,25	1.10	2 (8%)
2	BOG	B	800	-	20,20,20	0.55	0	25,25,25	1.27	3 (12%)
3	PO4	B	801	-	4,4,4	0.32	0	6,6,6	0.34	0
5	EDO	B	802	-	3,3,3	1.05	0	2,2,2	0.90	0
5	EDO	B	803	-	3,3,3	0.89	0	2,2,2	0.82	0
5	EDO	B	804	-	3,3,3	0.60	0	2,2,2	0.91	0
5	EDO	B	805	-	3,3,3	0.61	0	2,2,2	0.69	0
5	EDO	B	806	-	3,3,3	1.14	0	2,2,2	0.37	0
5	EDO	B	807	-	3,3,3	1.08	0	2,2,2	0.63	0
5	EDO	B	808	-	3,3,3	0.91	0	2,2,2	0.36	0
5	EDO	B	809	-	3,3,3	0.67	0	2,2,2	0.42	0
5	EDO	B	810	-	3,3,3	0.66	0	2,2,2	0.44	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	EDO	B	811	-	3,3,3	0.42	0	2,2,2	0.81	0
7	TAM	B	812	-	10,10,10	0.63	0	12,12,12	2.56	6 (50%)
5	EDO	B	813	-	3,3,3	0.31	0	2,2,2	0.88	0
5	EDO	B	814	-	3,3,3	0.79	0	2,2,2	0.21	0
5	EDO	B	815	-	3,3,3	0.47	0	2,2,2	0.92	0
8	13P	B	816	-	9,9,9	5.23	6 (66%)	12,12,12	4.22	6 (50%)
5	EDO	B	817	-	3,3,3	0.73	0	2,2,2	0.33	0
5	EDO	B	818	-	3,3,3	0.64	0	2,2,2	0.30	0
6	IMD	B	819	-	5,5,5	1.51	2 (40%)	5,5,5	0.48	0
9	BCN	B	820	-	10,10,10	0.90	0	11,11,11	2.04	3 (27%)
9	BCN	B	821	-	10,10,10	1.04	0	11,11,11	2.07	4 (36%)

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
2	BOG	A	1949	-	-	0/11/31/31	0/1/1/1
2	BOG	A	1950	-	-	0/11/31/31	0/1/1/1
3	PO4	A	1951	-	-	0/0/0/0	0/0/0/0
5	EDO	A	1952	-	-	0/1/1/1	0/0/0/0
5	EDO	A	1953	-	-	0/1/1/1	0/0/0/0
5	EDO	A	1954	-	-	0/1/1/1	0/0/0/0
5	EDO	A	1955	-	-	0/1/1/1	0/0/0/0
5	EDO	A	1956	-	-	0/1/1/1	0/0/0/0
5	EDO	A	1957	-	-	0/1/1/1	0/0/0/0
5	EDO	A	1958	-	-	0/1/1/1	0/0/0/0
5	EDO	A	1959	-	-	0/1/1/1	0/0/0/0
6	IMD	A	1960	-	-	0/0/0/0	0/1/1/1
6	IMD	A	1961	-	-	0/0/0/0	0/1/1/1
5	EDO	A	1962	-	-	0/1/1/1	0/0/0/0
6	IMD	A	1963	-	-	0/0/0/0	0/1/1/1
5	EDO	A	1964	-	-	0/1/1/1	0/0/0/0
5	EDO	A	1965	-	-	0/1/1/1	0/0/0/0
6	IMD	A	1966	-	-	0/0/0/0	0/1/1/1
6	IMD	A	1967	-	-	0/0/0/0	0/1/1/1
8	13P	A	1968	-	-	0/8/8/8	0/0/0/0
9	BCN	A	1969	-	-	0/10/10/10	0/0/0/0
4	FAD	A	600	-	-	0/34/50/50	0/1/6/6
2	BOG	A	700	-	-	0/11/31/31	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	BOG	A	800	-	-	0/11/31/31	0/1/1/1
4	FAD	B	600	-	-	0/34/50/50	0/1/6/6
2	BOG	B	700	-	-	0/11/31/31	0/1/1/1
2	BOG	B	800	-	-	0/11/31/31	0/1/1/1
3	PO4	B	801	-	-	0/0/0/0	0/0/0/0
5	EDO	B	802	-	-	0/1/1/1	0/0/0/0
5	EDO	B	803	-	-	0/1/1/1	0/0/0/0
5	EDO	B	804	-	-	0/1/1/1	0/0/0/0
5	EDO	B	805	-	-	0/1/1/1	0/0/0/0
5	EDO	B	806	-	-	0/1/1/1	0/0/0/0
5	EDO	B	807	-	-	0/1/1/1	0/0/0/0
5	EDO	B	808	-	-	0/1/1/1	0/0/0/0
5	EDO	B	809	-	-	0/1/1/1	0/0/0/0
5	EDO	B	810	-	-	0/1/1/1	0/0/0/0
5	EDO	B	811	-	-	0/1/1/1	0/0/0/0
7	TAM	B	812	-	-	0/12/12/12	0/0/0/0
5	EDO	B	813	-	-	0/1/1/1	0/0/0/0
5	EDO	B	814	-	-	0/1/1/1	0/0/0/0
5	EDO	B	815	-	-	0/1/1/1	0/0/0/0
8	13P	B	816	-	-	0/8/8/8	0/0/0/0
5	EDO	B	817	-	-	0/1/1/1	0/0/0/0
5	EDO	B	818	-	-	0/1/1/1	0/0/0/0
6	IMD	B	819	-	-	0/0/0/0	0/1/1/1
9	BCN	B	820	-	-	0/10/10/10	0/0/0/0
9	BCN	B	821	-	-	0/10/10/10	0/0/0/0

All (59) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	B	816	13P	O1-C1	12.46	1.51	1.43
8	A	1968	13P	O1-C1	11.86	1.51	1.43
4	B	600	FAD	C1'-N10	6.77	1.55	1.48
8	A	1968	13P	O2-C2	5.98	1.33	1.21
8	B	816	13P	C1-C2	5.74	1.61	1.50
8	A	1968	13P	C1-C2	5.48	1.61	1.50
8	B	816	13P	O2-C2	5.31	1.31	1.21
4	A	600	FAD	C6-C5X	-4.50	1.36	1.41
4	B	600	FAD	C6-C5X	-3.93	1.37	1.41
8	A	1968	13P	C3-C2	3.91	1.61	1.50
4	B	600	FAD	C10-N1	3.71	1.42	1.35
4	B	600	FAD	C4'-C3'	-3.56	1.46	1.53
8	A	1968	13P	O3-C3	3.55	1.53	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	B	816	13P	C3-C2	3.47	1.59	1.50
4	A	600	FAD	PA-O2A	-3.08	1.41	1.55
8	B	816	13P	O3-C3	2.99	1.51	1.41
2	A	1949	BOG	O1-C1	2.88	1.45	1.40
2	A	1950	BOG	O1-C1	2.84	1.45	1.40
4	A	600	FAD	P-O2P	-2.83	1.42	1.55
4	A	600	FAD	C9A-C5X	-2.76	1.36	1.42
4	A	600	FAD	O4'-C4'	-2.73	1.37	1.43
6	A	1961	IMD	C2-N1	2.69	1.37	1.31
4	B	600	FAD	O4B-C4B	-2.68	1.38	1.45
4	A	600	FAD	P-O1P	-2.67	1.41	1.51
2	B	700	BOG	O1-C1	2.65	1.45	1.40
6	A	1966	IMD	C2-N3	2.63	1.36	1.31
4	A	600	FAD	C2B-C1B	-2.59	1.49	1.53
4	B	600	FAD	C9A-C5X	-2.58	1.37	1.42
4	A	600	FAD	PA-O3P	-2.56	1.55	1.59
4	A	600	FAD	O2'-C2'	-2.54	1.37	1.43
6	A	1960	IMD	C2-N3	2.54	1.36	1.31
8	B	816	13P	P-O1	2.53	1.68	1.60
6	A	1963	IMD	C2-N1	2.44	1.36	1.31
4	A	600	FAD	C4-C4X	-2.43	1.37	1.41
4	A	600	FAD	O3'-C3'	-2.43	1.37	1.43
4	B	600	FAD	C9-C9A	-2.41	1.36	1.40
4	B	600	FAD	PA-O5B	-2.36	1.48	1.59
4	A	600	FAD	O4B-C4B	-2.36	1.39	1.45
6	B	819	IMD	C2-N3	2.34	1.36	1.31
8	A	1968	13P	P-O1	2.34	1.68	1.60
4	A	600	FAD	C5A-N7A	-2.33	1.31	1.40
4	B	600	FAD	C4-N3	-2.30	1.33	1.37
4	A	600	FAD	O4-C4	-2.29	1.20	1.24
6	A	1967	IMD	C2-N1	2.24	1.36	1.31
6	A	1960	IMD	C2-N1	2.22	1.36	1.31
6	B	819	IMD	C2-N1	2.20	1.36	1.31
6	A	1967	IMD	C2-N3	2.18	1.36	1.31
2	A	700	BOG	O1-C1	2.18	1.44	1.40
4	B	600	FAD	P-O2P	-2.17	1.45	1.55
6	A	1966	IMD	C2-N1	2.16	1.36	1.31
4	B	600	FAD	C4X-N5	2.15	1.40	1.36
4	A	600	FAD	C9-C9A	-2.15	1.36	1.40
6	A	1961	IMD	C2-N3	2.13	1.35	1.31
4	A	600	FAD	C4X-C10	-2.13	1.37	1.40
6	A	1963	IMD	C2-N3	2.09	1.35	1.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	600	FAD	O3B-C3B	-2.08	1.37	1.43
4	A	600	FAD	O2-C2	-2.05	1.19	1.23
4	A	600	FAD	PA-O5B	-2.02	1.50	1.59
4	B	600	FAD	PA-O2A	-2.02	1.46	1.55

All (85) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	600	FAD	N3A-C2A-N1A	-8.93	121.25	128.71
8	A	1968	13P	O1-C1-C2	8.56	124.14	110.53
8	B	816	13P	O1-C1-C2	8.44	123.95	110.53
8	A	1968	13P	O3P-P-O1	-8.40	83.48	106.65
4	A	600	FAD	N3A-C2A-N1A	-8.08	121.96	128.71
8	B	816	13P	O3P-P-O1	-7.56	85.77	106.65
4	B	600	FAD	C4X-N5-C5X	7.19	124.77	116.69
8	B	816	13P	O2P-P-O1	6.09	123.44	106.65
4	A	600	FAD	C2-N1-C10	5.96	120.99	114.98
8	A	1968	13P	O2P-P-O1	5.94	123.05	106.65
4	B	600	FAD	C2-N1-C10	5.65	120.67	114.98
8	B	816	13P	O3P-P-O1P	5.44	128.24	110.44
2	A	1950	BOG	C3-C4-C5	5.18	119.44	110.20
8	A	1968	13P	O3P-P-O1P	5.13	127.21	110.44
4	B	600	FAD	C4X-C10-N10	-4.79	118.12	120.51
9	A	1969	BCN	C5-N1-C3	4.62	122.72	111.45
9	B	821	BCN	C1-N1-C5	4.59	123.85	112.34
9	B	820	BCN	C1-N1-C3	4.50	123.63	112.34
4	A	600	FAD	N3A-C4A-N9A	4.35	133.29	125.43
7	B	812	TAM	C3-C-N	-4.35	98.18	107.92
7	B	812	TAM	C3-C-C2	-4.33	103.02	110.49
8	A	1968	13P	O1-P-O1P	-4.28	94.16	106.71
4	A	600	FAD	O2A-PA-O5B	-4.26	87.04	108.51
4	B	600	FAD	C4X-C10-N1	-4.01	118.72	122.73
7	B	812	TAM	C2-C-N	3.87	116.59	107.92
4	B	600	FAD	O2'-C2'-C3'	-3.83	99.51	109.05
2	B	800	BOG	C1-C2-C3	3.74	117.28	110.00
8	A	1968	13P	P-O1-C1	3.49	133.82	121.22
2	A	800	BOG	C3-C4-C5	3.43	116.33	110.20
9	B	820	BCN	C5-N1-C3	3.40	119.77	111.45
4	A	600	FAD	C4X-N5-C5X	3.36	120.46	116.69
2	A	800	BOG	O5-C5-C4	3.24	115.76	109.76
4	A	600	FAD	C5A-C4A-N3A	-3.22	118.68	125.70
2	A	1950	BOG	O1-C1-C2	3.13	112.16	108.18

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	600	FAD	C10-C4X-N5	-3.05	116.75	120.45
4	B	600	FAD	C9A-N10-C10	3.01	124.72	121.77
2	B	700	BOG	O5-C5-C6	3.00	113.72	106.34
2	A	800	BOG	O3-C3-C4	-3.00	103.63	110.35
4	B	600	FAD	O2'-C2'-C1'	3.00	117.16	109.71
2	A	1950	BOG	C4-C3-C2	2.98	116.33	110.82
9	B	821	BCN	C2-C1-N1	2.97	124.73	113.73
2	A	800	BOG	C4-C3-C2	2.96	116.29	110.82
4	A	600	FAD	C4B-O4B-C1B	2.94	112.95	109.75
4	A	600	FAD	C4X-C10-N1	-2.94	119.79	122.73
4	A	600	FAD	O3P-PA-O5B	2.92	116.47	103.41
2	A	1949	BOG	O1-C1-C2	2.91	111.88	108.18
2	A	1950	BOG	O5-C5-C4	2.88	115.08	109.76
4	B	600	FAD	N3A-C4A-N9A	2.87	130.61	125.43
4	B	600	FAD	C9A-C5X-N5	-2.87	117.97	122.37
8	B	816	13P	P-O1-C1	2.81	131.36	121.22
4	A	600	FAD	C5X-C9A-N10	2.78	119.54	116.80
9	B	821	BCN	C1-N1-C3	2.74	119.22	112.34
4	B	600	FAD	N1-C10-N10	2.71	123.08	115.97
4	B	600	FAD	O3'-C3'-C2'	-2.70	101.91	108.74
4	B	600	FAD	O4B-C1B-N9A	-2.69	105.93	108.44
4	A	600	FAD	O5B-C5B-C4B	2.68	118.77	108.94
4	B	600	FAD	C1'-C2'-C3'	-2.68	102.16	109.82
7	B	812	TAM	C5-C2-C	2.63	120.10	115.81
9	A	1969	BCN	C1-N1-C5	2.59	118.83	112.34
4	B	600	FAD	O5'-P-O1P	-2.55	99.38	109.37
7	B	812	TAM	C2-C-C1	2.47	114.75	110.49
2	A	700	BOG	O2-C2-C1	2.47	115.42	110.04
4	A	600	FAD	O4B-C1B-N9A	2.45	110.72	108.44
9	A	1969	BCN	C1-N1-C3	2.43	118.44	112.34
2	A	700	BOG	C1-O5-C5	2.42	118.44	113.73
2	A	1950	BOG	O4-C4-C3	-2.42	104.93	110.35
8	B	816	13P	O1-P-O1P	-2.41	99.64	106.71
4	A	600	FAD	C4A-C5A-N7A	-2.40	107.47	109.52
4	A	600	FAD	O5B-PA-O1A	2.35	118.60	109.37
4	B	600	FAD	C4B-O4B-C1B	-2.34	107.21	109.75
2	A	800	BOG	O5-C1-C2	-2.32	105.55	110.31
9	A	1969	BCN	C6-C5-N1	-2.31	106.03	113.62
4	B	600	FAD	C2'-C1'-N10	-2.29	109.41	112.45
4	A	600	FAD	C2A-N3A-C4A	2.25	120.40	114.01
7	B	812	TAM	C3-C-C1	2.24	114.36	110.49
9	B	821	BCN	C5-N1-C3	2.23	116.90	111.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	800	BOG	C4-C3-C2	2.22	114.92	110.82
6	A	1966	IMD	C4-N3-C2	2.18	111.33	108.20
4	A	600	FAD	O4'-C4'-C3'	2.11	114.31	109.05
2	B	700	BOG	O5-C1-O1	2.11	114.93	109.98
9	B	820	BCN	C2-C1-N1	2.11	121.52	113.73
4	B	600	FAD	C2B-C3B-C4B	2.10	106.84	102.65
4	B	600	FAD	C4A-C5A-N7A	-2.08	107.74	109.52
2	B	800	BOG	O3-C3-C4	-2.04	105.79	110.35
2	A	1949	BOG	O5-C5-C6	2.02	111.30	106.34

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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	494/501 (98%)	1.28	83 (16%) 2 2	20, 41, 65, 84	0
1	B	497/501 (99%)	1.34	83 (16%) 2 2	19, 41, 66, 98	0
All	All	991/1002 (98%)	1.31	166 (16%) 2 2	19, 41, 66, 98	0

All (166) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	113	LYS	10.4
1	B	494	GLN	8.6
1	B	59	TYR	7.6
1	A	1	MET	6.7
1	B	106	PHE	6.3
1	B	1	MET	5.7
1	A	115	THR	5.5
1	A	325	ASP	5.2
1	B	376	ILE	5.1
1	B	141	TYR	4.3
1	B	135	ILE	4.3
1	A	23	ALA	4.1
1	A	197	TRP	4.1
1	B	111	LEU	3.9
1	B	497	LEU	3.9
1	A	58	HIS	3.8
1	B	110	HIS	3.8
1	B	26	GLY	3.8
1	A	369	LEU	3.7
1	B	374	GLN	3.7
1	B	391	ILE	3.7
1	B	163	GLY	3.7
1	B	49	ILE	3.6
1	A	318	PRO	3.5

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Mol	Chain	Res	Type	RSRZ
1	B	389	GLY	3.5
1	A	320	CYS	3.4
1	A	113	LYS	3.4
1	A	98	ALA	3.4
1	B	100	MET	3.3
1	A	453	TYR	3.2
1	A	345	ALA	3.2
1	B	375	GLY	3.1
1	A	111	LEU	3.1
1	B	402	LEU	3.1
1	A	81	ILE	3.1
1	B	435	VAL	3.0
1	A	19	ALA	3.0
1	B	117	LEU	3.0
1	A	117	LEU	3.0
1	B	139	PHE	3.0
1	A	319	LEU	3.0
1	A	122	GLY	2.9
1	B	72	GLU	2.9
1	B	247	ILE	2.9
1	A	106	PHE	2.9
1	A	59	TYR	2.9
1	A	121	THR	2.8
1	A	272	ASP	2.8
1	B	262	MSE	2.8
1	A	64	VAL	2.8
1	B	124	ARG	2.8
1	B	346	PRO	2.8
1	A	8	VAL	2.7
1	A	367	GLU	2.7
1	B	367	GLU	2.7
1	A	380	TRP	2.7
1	B	181	LEU	2.7
1	B	133	PRO	2.7
1	A	437	ASP	2.7
1	B	66	GLU	2.7
1	B	291	TYR	2.6
1	B	330	ILE	2.6
1	A	12	GLY	2.6
1	B	365	ALA	2.6
1	A	393	GLY	2.6
1	A	337	ASP	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	131	LEU	2.6
1	B	142	SER	2.6
1	B	273	VAL	2.6
1	A	243	LYS	2.6
1	A	51	GLY	2.6
1	B	309	ILE	2.6
1	A	195	TYR	2.6
1	B	83	PHE	2.5
1	A	304	LEU	2.5
1	A	427	LEU	2.5
1	B	198	GLN	2.5
1	A	443	GLY	2.5
1	A	107	MET	2.5
1	A	11	GLY	2.5
1	A	355	LEU	2.5
1	A	123	LEU	2.5
1	A	160	VAL	2.5
1	B	108	TYR	2.5
1	A	204	ASN	2.5
1	A	329	ALA	2.5
1	A	402	LEU	2.4
1	B	259	ILE	2.4
1	B	180	GLY	2.4
1	B	98	ALA	2.4
1	B	361	LEU	2.4
1	B	312	THR	2.4
1	A	432	ALA	2.4
1	B	183	ILE	2.4
1	A	29	VAL	2.4
1	A	74	LEU	2.4
1	B	67	ALA	2.4
1	A	143	ASP	2.4
1	B	210	VAL	2.4
1	A	6	LEU	2.4
1	A	37	LEU	2.4
1	B	355	LEU	2.4
1	A	488	TRP	2.4
1	A	61	PHE	2.3
1	B	226	ILE	2.3
1	B	341	GLU	2.3
1	A	436	SER	2.3
1	B	145	TRP	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	399	ALA	2.3
1	B	127	ALA	2.3
1	A	291	TYR	2.3
1	B	252	ASP	2.3
1	B	282	VAL	2.3
1	A	429	LEU	2.3
1	B	228	LEU	2.3
1	A	13	ILE	2.3
1	A	247	ILE	2.3
1	A	277	GLY	2.3
1	A	375	GLY	2.3
1	B	255	ILE	2.3
1	B	342	ASN	2.3
1	B	392	GLU	2.2
1	A	159	VAL	2.2
1	A	184	VAL	2.2
1	A	418	ALA	2.2
1	B	60	GLU	2.2
1	A	9	ILE	2.2
1	A	49	ILE	2.2
1	B	101	ILE	2.2
1	A	118	PRO	2.2
1	B	27	LEU	2.2
1	B	107	MET	2.2
1	B	462	ARG	2.2
1	B	257	PHE	2.2
1	B	158	MSE	2.2
1	A	44	ALA	2.2
1	B	195	TYR	2.2
1	B	246	TYR	2.2
1	A	135	ILE	2.1
1	A	226	ILE	2.1
1	B	314	SER	2.1
1	B	373	TYR	2.1
1	B	459	TRP	2.1
1	A	347	LEU	2.1
1	A	144	CYS	2.1
1	A	18	ILE	2.1
1	A	276	LYS	2.1
1	B	496	ARG	2.1
1	A	339	HIS	2.1
1	A	338	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	272	ASP	2.1
1	A	261	TRP	2.1
1	B	34	ALA	2.1
1	B	384	SER	2.1
1	A	166	VAL	2.1
1	B	334	TYR	2.1
1	A	183	ILE	2.0
1	B	284	ILE	2.0
1	A	32	LEU	2.0
1	B	372	TYR	2.0
1	B	199	ALA	2.0
1	A	346	PRO	2.0
1	B	114	ARG	2.0
1	B	188	ASP	2.0
1	A	67	ALA	2.0
1	A	422	GLY	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 ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	BOG	A	800	20/20	1.70	33.06	131,136,138,138	0
5	EDO	B	817	4/4	0.74	23.45	55,62,65,66	0
5	EDO	B	806	4/4	0.63	21.18	34,45,50,55	0
6	IMD	A	1960	5/5	0.37	9.65	58,58,60,60	0
5	EDO	A	1956	4/4	0.49	9.45	41,44,50,54	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	PO4	B	801	5/5	0.52	9.06	68,68,70,71	0
6	IMD	A	1966	5/5	0.24	8.68	70,70,72,72	0
2	BOG	A	1950	20/20	0.67	8.50	90,93,94,95	0
5	EDO	A	1958	4/4	0.36	7.98	32,34,40,46	0
5	EDO	A	1955	4/4	0.31	7.94	38,50,50,52	0
7	TAM	B	812	11/11	0.49	7.76	56,63,66,70	0
9	BCN	A	1969	11/11	0.59	6.66	63,76,79,79	0
9	BCN	B	820	11/11	0.47	6.23	64,69,72,74	0
2	BOG	B	800	20/20	0.66	5.88	133,136,139,140	0
9	BCN	B	821	11/11	0.44	5.21	58,64,70,70	0
5	EDO	A	1953	4/4	0.22	5.06	32,45,46,47	0
5	EDO	B	814	4/4	0.23	4.79	54,56,56,56	0
5	EDO	B	802	4/4	0.37	3.84	38,44,48,50	0
5	EDO	A	1952	4/4	0.23	3.06	63,63,64,64	0
5	EDO	B	811	4/4	0.28	3.05	36,43,44,49	0
2	BOG	B	700	20/20	0.30	2.98	63,71,77,79	0
5	EDO	A	1965	4/4	0.24	2.83	37,37,42,42	0
5	EDO	B	808	4/4	0.29	2.77	52,56,57,60	0
5	EDO	A	1957	4/4	0.18	2.62	43,48,48,49	0
5	EDO	B	803	4/4	0.23	2.35	47,49,52,53	0
5	EDO	B	810	4/4	0.26	2.21	53,53,54,58	0
2	BOG	A	1949	20/20	0.55	2.00	92,95,97,97	0
5	EDO	B	805	4/4	0.20	1.86	63,64,64,68	0
5	EDO	B	813	4/4	0.25	1.60	36,39,41,43	0
6	IMD	A	1961	5/5	0.25	1.50	73,73,74,74	0
5	EDO	B	807	4/4	0.22	1.37	46,47,50,52	0
8	13P	B	816	10/10	0.24	1.19	28,40,45,45	0
8	13P	A	1968	10/10	0.23	0.84	28,38,43,44	0
2	BOG	A	700	20/20	0.31	0.73	86,91,94,94	0
5	EDO	B	804	4/4	0.23	0.73	49,49,51,52	0
6	IMD	A	1967	5/5	0.19	0.50	63,64,65,67	0
6	IMD	A	1963	5/5	0.19	0.29	42,47,49,50	0
5	EDO	B	815	4/4	0.17	-0.22	54,57,59,62	0
4	FAD	A	600	53/53	0.16	-0.39	18,26,29,36	0
4	FAD	B	600	53/53	0.15	-0.40	19,25,29,32	0
6	IMD	B	819	5/5	0.18	-0.58	78,78,79,81	0
5	EDO	A	1954	4/4	0.15	-0.97	64,65,65,66	0
5	EDO	A	1962	4/4	0.16	-1.10	70,71,72,74	0
3	PO4	A	1951	5/5	0.15	-1.11	84,84,84,85	0
5	EDO	B	809	4/4	0.19	-1.55	65,66,67,68	0
5	EDO	A	1959	4/4	0.14	-1.69	43,48,49,52	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	EDO	A	1964	4/4	0.19	-2.05	64,66,66,67	0
5	EDO	B	818	4/4	0.21	-	72,75,75,75	0

6.5 Other polymers ⓘ

There are no such residues in this entry.