



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 02:14 PM UTC

PDB ID : 2W40 / pdb_00002w40
Title : Crystal structure of Plasmodium falciparum glycerol kinase with bound glycerol
Authors : Schnick, C.; Polley, S.D.; Fivelman, Q.L.; Ranford-Cartwright, L.; Wilkinson, S.R.; Brannigan, J.A.; Wilkinson, A.J.; Baker, D.A.
Deposited on : 2008-11-18
Resolution : 1.49 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

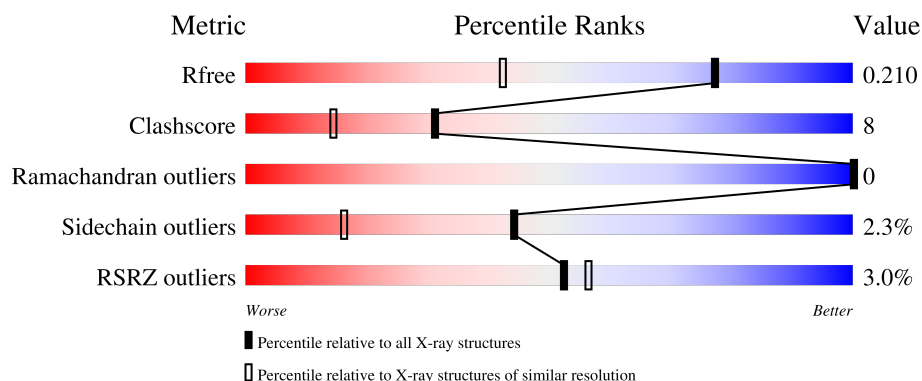
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.49 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4037 (1.50-1.50)
Clashscore	190562	4235 (1.50-1.50)
Ramachandran outliers	187476	4153 (1.50-1.50)
Sidechain outliers	187428	4150 (1.50-1.50)
RSRZ outliers	180081	4039 (1.50-1.50)

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	503	4% 84% 14% .
1	B	503	3% 84% 14% .
1	C	503	3% 83% 15% .
1	D	503	3% 84% 14% .

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
2	EDO	A	1507	-	-	X	-
2	EDO	A	1509	-	-	X	-
2	EDO	B	1514	-	-	X	-
2	EDO	B	1518	-	-	X	-
2	EDO	C	1510	-	-	X	-
2	EDO	C	1514	-	X	X	-
2	EDO	C	1520	-	-	X	-
2	EDO	D	1510	-	-	X	-
2	EDO	D	1515	-	X	X	-
3	GOL	C	1527	-	-	X	-

2 Entry composition [i](#)

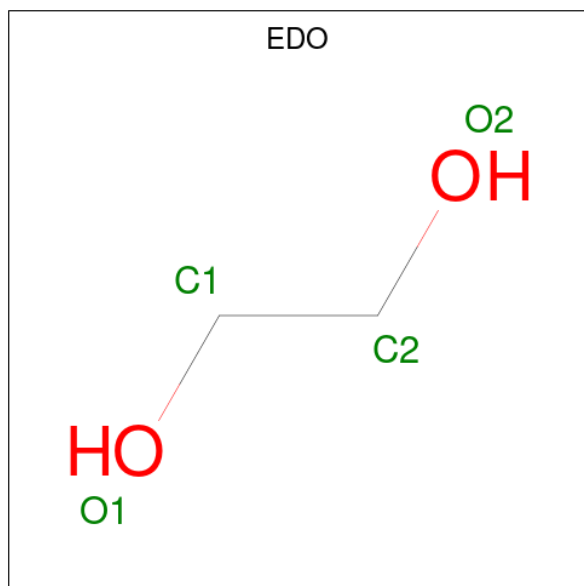
There are 4 unique types of molecules in this entry. The entry contains 18365 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 GLYCEROL KINASE, PUTATIVE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	501	Total	C	N	O	S	11	10	0
			4034	2568	668	771	27			
1	B	503	Total	C	N	O	S	5	7	0
			4030	2565	667	771	27			
1	C	502	Total	C	N	O	S	21	9	0
			4037	2569	669	773	26			
1	D	503	Total	C	N	O	S	3	7	0
			4015	2555	662	772	26			

- Molecule 2 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



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

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

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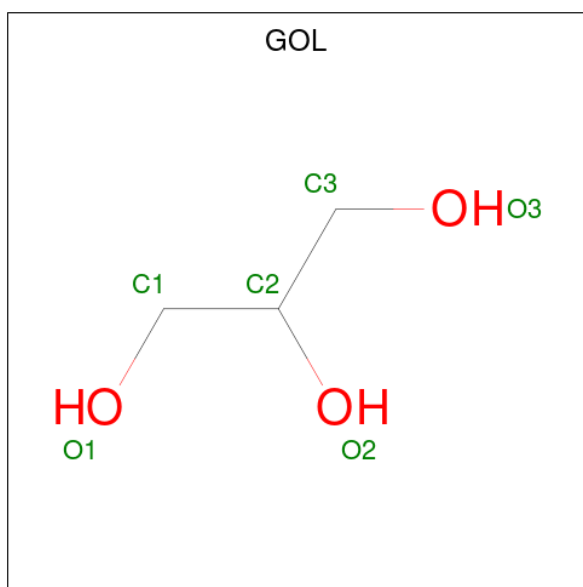
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total 4	C 2	O 2	0	0
2	B	1	Total 4	C 2	O 2	0	0
2	B	1	Total 4	C 2	O 2	0	0
2	B	1	Total 4	C 2	O 2	0	0
2	B	1	Total 4	C 2	O 2	0	0
2	B	1	Total 4	C 2	O 2	0	0
2	C	1	Total 4	C 2	O 2	0	0
2	C	1	Total 4	C 2	O 2	0	0
2	C	1	Total 4	C 2	O 2	0	0
2	C	1	Total 4	C 2	O 2	0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

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

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		

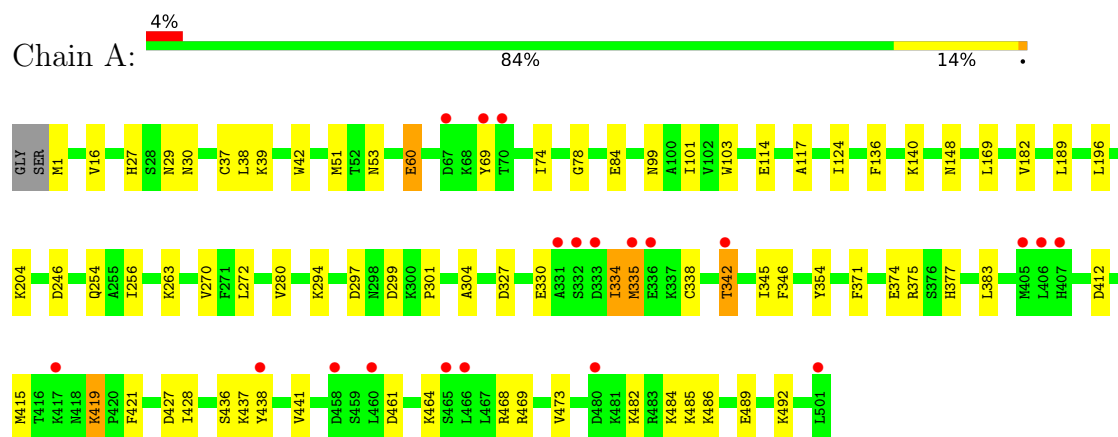
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	463	Total 463	O 463	0	0
4	B	492	Total 492	O 492	0	0
4	C	499	Total 499	O 499	0	0
4	D	475	Total 475	O 475	0	0

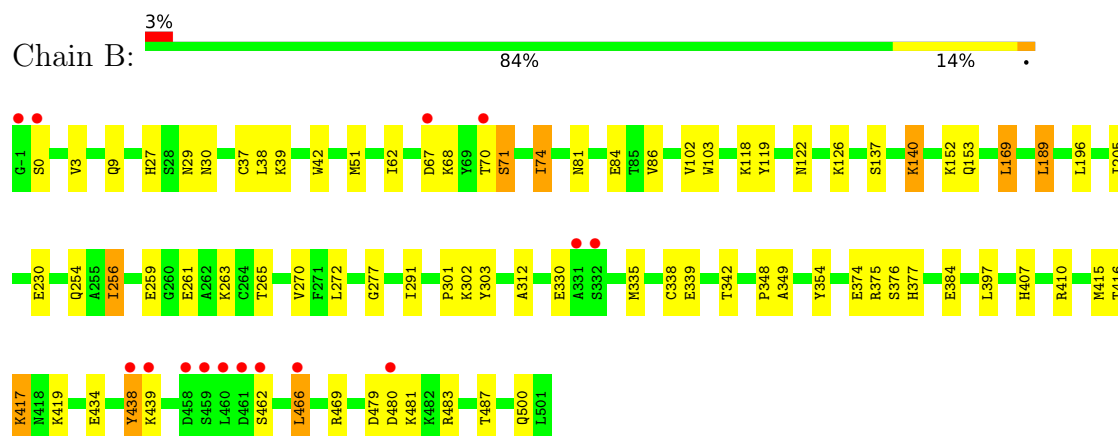
3 Residue-property plots [i](#)

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.

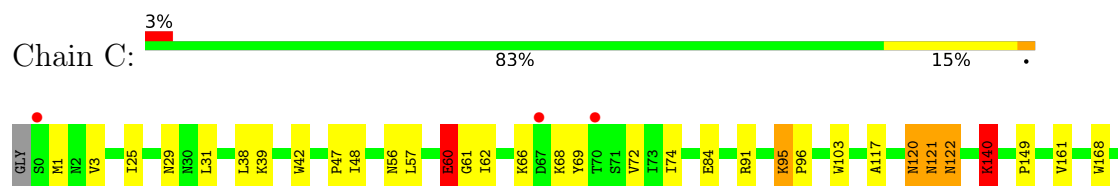
• Molecule 1: GLYCEROL KINASE, PUTATIVE

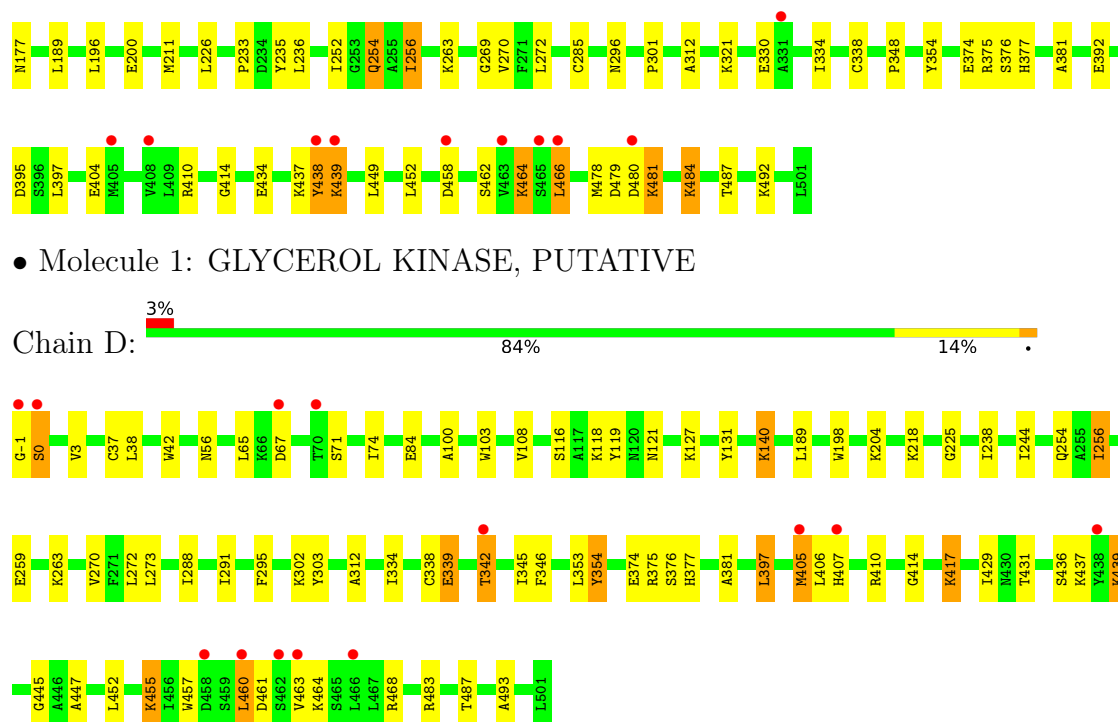


• Molecule 1: GLYCEROL KINASE, PUTATIVE



• Molecule 1: GLYCEROL KINASE, PUTATIVE





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	101.92Å 100.72Å 107.81Å 90.00° 92.56° 90.00°	Depositor
Resolution (Å)	107.83 – 1.49 107.70 – 1.49	Depositor EDS
% Data completeness (in resolution range)	73.5 (107.83-1.49) 73.6 (107.70-1.49)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.66 (at 1.49Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.168 , 0.210 0.161 , 0.210	Depositor DCC
R_{free} test set	13131 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	13.0	Xtriage
Anisotropy	0.258	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 51.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.000 for k,h,-l 0.000 for -k,-h,-l 0.000 for h,-k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	18365	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 49.76 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 7.1291e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: EDO, GOL

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	1.51	16/4133 (0.4%)	1.21	10/5592 (0.2%)
1	B	1.56	19/4113 (0.5%)	1.23	8/5565 (0.1%)
1	C	1.58	22/4132 (0.5%)	1.24	10/5590 (0.2%)
1	D	1.55	25/4111 (0.6%)	1.20	5/5563 (0.1%)
All	All	1.55	82/16489 (0.5%)	1.22	33/22310 (0.1%)

All (82) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	95	LYS	CD-CE	-12.66	1.14	1.52
1	C	68	LYS	CG-CD	-11.83	1.17	1.52
1	C	439	LYS	C-O	11.13	1.37	1.24
1	C	439	LYS	N-CA	10.54	1.59	1.46
1	B	439	LYS	CA-CB	10.06	1.69	1.53
1	B	439	LYS	CA-C	9.88	1.65	1.52
1	C	434	GLU	CB-CG	-8.94	1.25	1.52
1	C	285	CYS	CA-CB	7.35	1.64	1.53
1	D	375	ARG	CB-CG	-7.18	1.30	1.52
1	D	140	LYS	N-CA	7.08	1.54	1.46
1	C	439	LYS	CB-CG	7.00	1.73	1.52
1	A	346	PHE	N-CA	6.95	1.54	1.46
1	C	121	ASN	N-CA	6.89	1.54	1.46
1	C	438	TYR	C-O	-6.75	1.16	1.24
1	B	480	ASP	CG-OD1	6.71	1.38	1.25
1	D	127	LYS	C-O	-6.63	1.16	1.24
1	B	348	PRO	C-O	6.52	1.32	1.24
1	B	375	ARG	CB-CG	-6.30	1.33	1.52
1	D	342	THR	CA-CB	6.26	1.63	1.53
1	A	375	ARG	CB-CG	-6.22	1.33	1.52
1	A	427	ASP	N-CA	6.22	1.53	1.46
1	D	225	GLY	C-O	-6.20	1.17	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	244	ILE	CA-CB	6.19	1.61	1.55
1	B	384	GLU	N-CA	-6.18	1.38	1.46
1	C	25	ILE	CA-CB	6.13	1.61	1.53
1	A	124	ILE	CA-CB	6.12	1.62	1.54
1	C	480	ASP	CG-OD2	6.08	1.36	1.25
1	C	437	LYS	CA-C	6.07	1.60	1.52
1	A	78	GLY	N-CA	6.05	1.52	1.45
1	D	259	GLU	C-O	5.91	1.30	1.23
1	A	342	THR	N-CA	5.87	1.54	1.46
1	D	288	ILE	C-O	-5.86	1.17	1.24
1	D	447	ALA	CA-CB	-5.85	1.44	1.53
1	D	445	GLY	C-O	5.85	1.30	1.23
1	D	342	THR	N-CA	5.83	1.54	1.46
1	B	74	ILE	CA-CB	5.78	1.60	1.53
1	B	205	ILE	CA-CB	5.73	1.60	1.54
1	D	218	LYS	CD-CE	-5.71	1.35	1.52
1	D	67	ASP	CB-CG	5.70	1.66	1.52
1	B	480	ASP	CB-CG	5.68	1.66	1.52
1	C	381	ALA	CA-CB	5.64	1.62	1.53
1	B	102	VAL	CA-CB	5.63	1.62	1.54
1	B	439	LYS	C-O	5.61	1.30	1.24
1	C	168	TRP	CA-C	5.61	1.59	1.52
1	A	280	VAL	C-O	5.56	1.29	1.24
1	D	140	LYS	CE-NZ	-5.52	1.32	1.49
1	B	349	ALA	C-O	-5.52	1.17	1.24
1	D	354[A]	TYR	C-O	5.51	1.30	1.24
1	D	354[B]	TYR	C-O	5.51	1.30	1.24
1	A	16	VAL	CA-C	5.50	1.59	1.52
1	B	254	GLN	N-CA	5.46	1.53	1.46
1	B	354	TYR	CA-C	5.46	1.56	1.53
1	A	38	LEU	CA-C	-5.44	1.45	1.52
1	D	256	ILE	CA-CB	5.42	1.61	1.54
1	C	161	VAL	C-O	5.41	1.29	1.24
1	B	407	HIS	N-CA	5.39	1.52	1.46
1	D	381	ALA	N-CA	-5.37	1.39	1.46
1	A	375	ARG	CA-C	-5.37	1.45	1.52
1	D	463	VAL	CA-CB	5.34	1.61	1.54
1	C	348	PRO	C-O	5.32	1.30	1.24
1	C	269	GLY	N-CA	5.31	1.48	1.44
1	D	204	LYS	C-O	5.31	1.30	1.24
1	C	117	ALA	C-O	5.28	1.30	1.24
1	B	265	THR	CA-CB	5.24	1.61	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	346	PHE	N-CA	5.23	1.53	1.46
1	A	304	ALA	C-O	-5.22	1.17	1.23
1	A	148[A]	ASN	C-O	-5.22	1.20	1.24
1	A	148[B]	ASN	C-O	-5.22	1.20	1.24
1	B	342	THR	CA-CB	-5.20	1.45	1.53
1	D	334	ILE	CA-CB	5.19	1.60	1.54
1	A	419	LYS	C-O	-5.17	1.19	1.24
1	D	295	PHE	CE1-CZ	-5.17	1.23	1.38
1	C	149	PRO	C-O	-5.16	1.17	1.24
1	B	438	TYR	C-O	-5.15	1.17	1.23
1	D	100	ALA	CA-CB	5.13	1.61	1.53
1	C	321	LYS	CD-CE	5.11	1.67	1.52
1	A	270	VAL	N-CA	5.10	1.52	1.46
1	C	375	ARG	CB-CG	-5.09	1.37	1.52
1	C	60	GLU	CD-OE2	5.06	1.34	1.25
1	A	53	ASN	C-O	-5.02	1.18	1.24
1	D	353	LEU	C-N	5.02	1.37	1.33
1	B	189	LEU	CA-C	5.01	1.58	1.52

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	439	LYS	N-CA-CB	-12.01	92.65	110.07
1	C	95	LYS	CG-CD-CE	8.51	130.87	111.30
1	A	246	ASP	N-CA-C	7.53	119.49	111.28
1	B	439	LYS	N-CA-C	7.23	118.94	111.14
1	C	481	LYS	CA-CB-CG	-6.39	101.32	114.10
1	D	437	LYS	CB-CG-CD	-6.23	96.97	111.30
1	B	335	MET	N-CA-C	6.15	118.78	111.71
1	A	335	MET	CG-SD-CE	-6.03	87.62	100.90
1	C	464	LYS	N-CA-C	5.93	117.75	111.28
1	D	218	LYS	CG-CD-CE	5.86	124.77	111.30
1	B	71	SER	N-CA-C	-5.83	99.55	108.76
1	C	414	GLY	N-CA-C	5.82	119.69	112.64
1	D	218	LYS	CB-CG-CD	5.65	124.29	111.30
1	A	182	VAL	N-CA-C	5.61	117.01	110.62
1	A	297	ASP	CA-C-N	5.57	128.77	120.31
1	A	297	ASP	C-N-CA	5.57	128.77	120.31
1	B	480	ASP	N-CA-CB	5.54	118.76	110.22
1	C	438	TYR	CA-C-O	-5.50	114.20	120.58
1	C	60	GLU	CB-CG-CD	5.46	121.88	112.60
1	A	437	LYS	CB-CG-CD	-5.41	98.86	111.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	51	MET	CG-SD-CE	5.39	112.75	100.90
1	A	335	MET	N-CA-C	5.30	117.96	111.82
1	C	256	ILE	CA-C-O	5.24	127.33	120.78
1	B	439	LYS	CA-C-O	5.22	126.08	120.70
1	B	439	LYS	CB-CA-C	5.14	119.02	110.90
1	B	81	ASN	N-CA-C	5.07	116.77	108.76
1	C	478	MET	CA-C-N	5.07	128.41	120.75
1	C	478	MET	C-N-CA	5.07	128.41	120.75
1	A	428	ILE	N-CA-C	5.06	115.83	110.72
1	D	198	TRP	N-CA-C	-5.06	102.89	110.23
1	C	140	LYS	CA-CB-CG	-5.04	104.03	114.10
1	A	334	ILE	N-CA-CB	5.02	116.42	110.55
1	D	375	ARG	CB-CA-C	-5.02	102.15	110.68

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	4034	0	4060	48	0
1	B	4030	0	4051	63	0
1	C	4037	0	4064	76	0
1	D	4015	0	4033	59	0
2	A	36	0	53	9	0
2	B	80	0	120	16	0
2	C	76	0	114	27	0
2	D	56	0	81	14	0
3	A	12	0	16	3	0
3	B	12	0	15	1	0
3	C	42	0	56	14	0
3	D	6	0	8	0	0
4	A	463	0	0	10	1
4	B	492	0	0	22	2
4	C	499	0	0	24	0
4	D	475	0	0	14	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	18365	0	16671	270	2

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 (270) 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)
2:C:1510:EDO:C1	2:C:1510:EDO:O1	1.64	1.44
1:D:116:SER:O	2:D:1515:EDO:H12	1.31	1.27
1:D:116:SER:O	2:D:1515:EDO:C1	1.98	1.10
3:C:1522:GOL:H2	4:C:2429:HOH:O	1.53	1.08
1:D:452:LEU:O	1:D:455:LYS:HE2	1.51	1.07
1:C:226:LEU:H	2:C:1510:EDO:H12	1.19	1.05
1:B:126:LYS:NZ	4:B:2173:HOH:O	1.95	0.99
2:C:1514:EDO:H21	4:C:2356:HOH:O	1.62	0.96
1:B:500:GLN:HG2	4:B:2467:HOH:O	1.64	0.96
1:C:140:LYS:HE2	4:C:2110:HOH:O	1.65	0.95
3:B:1523:GOL:H2	4:B:2432:HOH:O	1.65	0.95
1:A:335:MET:HE1	1:A:421:PHE:HB2	1.48	0.94
1:B:374:GLU:H	1:B:377:HIS:HD2	1.16	0.93
1:D:374:GLU:H	1:D:377:HIS:HD2	1.17	0.93
1:C:91[B]:ARG:NH1	4:C:2119:HOH:O	1.88	0.91
1:D:121:ASN:ND2	2:D:1515:EDO:H21	1.84	0.91
1:B:118:LYS:HE3	4:B:2156:HOH:O	1.70	0.91
1:C:374:GLU:H	1:C:377:HIS:HD2	1.21	0.88
1:A:374:GLU:H	1:A:377:HIS:HD2	1.20	0.88
3:C:1527:GOL:O1	4:C:2499:HOH:O	1.86	0.86
1:A:330:GLU:OE2	4:A:2327:HOH:O	1.95	0.84
1:D:452:LEU:O	1:D:455:LYS:CE	2.26	0.84
1:D:118:LYS:NZ	4:D:2162:HOH:O	2.12	0.83
1:D:354[A]:TYR:HD1	4:D:2322:HOH:O	1.63	0.82
1:D:116:SER:C	2:D:1515:EDO:H12	2.03	0.82
1:C:140:LYS:CE	4:C:2110:HOH:O	2.24	0.80
1:C:479:ASP:OD2	1:C:481:LYS:HB2	1.83	0.79
1:C:200:GLU:CG	3:C:1527:GOL:O2	2.31	0.79
1:A:117:ALA:HB3	3:A:1512:GOL:H32	1.64	0.77
1:C:458:ASP:OD2	4:C:2433:HOH:O	2.02	0.77
1:D:338[A]:CYS:SG	1:D:376[A]:SER:HB2	2.26	0.75
1:D:457:TRP:O	4:D:2429:HOH:O	2.04	0.75
2:D:1515:EDO:H22	4:D:2472:HOH:O	1.85	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:226:LEU:N	2:C:1510:EDO:H12	2.00	0.75
1:B:37:CYS:HB3	2:B:1514:EDO:H11	1.67	0.74
1:A:485:LYS:NZ	1:A:489:GLU:OE2	2.21	0.73
1:A:117:ALA:CB	3:A:1512:GOL:H32	2.17	0.73
1:A:335:MET:HE2	1:A:383:LEU:HD13	1.70	0.73
1:C:200:GLU:HG2	3:C:1527:GOL:O2	1.88	0.73
1:C:438:TYR:OH	1:C:466:LEU:CD2	2.38	0.71
1:C:236:LEU:HD22	2:C:1520:EDO:H12	1.71	0.71
1:C:330:GLU:OE1	4:C:2341:HOH:O	2.08	0.71
1:A:327:ASP:O	1:A:330:GLU:HG2	1.91	0.70
2:A:1507:EDO:H22	4:A:2341:HOH:O	1.92	0.70
2:B:1518:EDO:O2	4:B:2488:HOH:O	1.90	0.69
1:C:200:GLU:HG3	3:C:1527:GOL:O2	1.92	0.69
1:C:57:LEU:O	1:C:60:GLU:HG3	1.94	0.68
1:D:436:SER:O	4:D:2417:HOH:O	2.13	0.66
1:B:410:ARG:NH1	1:B:434:GLU:OE2	2.28	0.66
1:C:177:ASN:HD21	2:C:1510:EDO:H11	1.61	0.66
1:B:417:LYS:HE2	4:B:2442:HOH:O	1.94	0.66
1:C:62:ILE:HG13	1:C:66:LYS:HE2	1.78	0.66
1:A:196:LEU:HD11	1:A:301:PRO:HG3	1.78	0.65
1:B:0:SER:CB	1:B:71:SER:OG	2.44	0.65
1:B:27:HIS:HD2	4:B:2010:HOH:O	1.78	0.65
1:A:114[A]:GLU:HG3	4:A:2139:HOH:O	1.96	0.65
1:C:31:LEU:HD21	3:C:1523:GOL:H32	1.79	0.64
2:C:1510:EDO:C1	2:C:1510:EDO:HO1	2.08	0.64
1:B:330:GLU:OE1	4:B:2341:HOH:O	2.14	0.64
1:D:3:VAL:CG2	1:D:74:ILE:HD13	2.27	0.64
1:D:397:LEU:C	1:D:397:LEU:HD23	2.22	0.64
1:A:473:VAL:H	1:B:153:GLN:HE22	1.45	0.64
2:C:1514:EDO:C2	4:C:2356:HOH:O	2.33	0.64
1:D:56:ASN:HB3	2:D:1514:EDO:O1	1.97	0.64
1:A:204:LYS:HG2	2:A:1509:EDO:H11	1.78	0.63
1:C:42:TRP:O	2:C:1514:EDO:H12	1.98	0.63
1:C:233:PRO:O	2:C:1520:EDO:H21	1.98	0.63
1:A:464:LYS:HB3	1:A:468:ARG:NH1	2.14	0.62
1:D:439:LYS:H	1:D:439:LYS:CD	2.10	0.62
1:C:392:GLU:O	2:C:1516:EDO:H21	2.00	0.62
1:D:119:TYR:O	2:D:1515:EDO:O1	2.17	0.62
1:D:342:THR:HG23	1:D:377:HIS:CE1	2.34	0.62
1:A:37:CYS:HB3	2:A:1507:EDO:H11	1.82	0.61
1:A:371:PHE:HD2	2:C:1514:EDO:H21	1.65	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:338[B]:CYS:SG	1:C:376:SER:HB2	2.40	0.61
1:C:376:SER:HB3	4:C:2387:HOH:O	1.99	0.61
1:A:204:LYS:HG2	2:A:1509:EDO:C1	2.30	0.60
1:D:272:LEU:C	1:D:272:LEU:HD23	2.27	0.60
1:D:121:ASN:HD22	2:D:1515:EDO:H21	1.65	0.60
1:A:99:ASN:HD22	2:A:1508:EDO:H22	1.65	0.60
1:C:121:ASN:HB2	4:C:2166:HOH:O	2.02	0.60
1:D:338[A]:CYS:SG	1:D:376[A]:SER:CB	2.90	0.60
1:C:438:TYR:HH	1:C:466:LEU:HD23	1.67	0.59
1:C:397:LEU:C	1:C:397:LEU:HD23	2.27	0.59
1:A:354[A]:TYR:HD2	4:A:2345:HOH:O	1.85	0.59
1:B:118:LYS:CE	4:B:2156:HOH:O	2.40	0.59
1:A:334:ILE:O	1:A:338[B]:CYS:HB2	2.03	0.59
1:B:29:ASN:C	2:B:1521:EDO:H22	2.27	0.58
1:C:438:TYR:OH	1:C:466:LEU:HD23	2.03	0.58
1:C:29:ASN:ND2	1:C:60:GLU:OE1	2.27	0.57
2:B:1517:EDO:H12	4:B:2070:HOH:O	2.04	0.57
1:A:464:LYS:NZ	4:A:2423:HOH:O	2.38	0.56
1:B:119:TYR:HA	2:B:1519:EDO:H12	1.85	0.56
1:C:438:TYR:OH	1:C:466:LEU:HD22	2.04	0.56
1:A:272:LEU:HD23	1:A:272:LEU:C	2.31	0.56
1:C:296:ASN:HD21	2:C:1507:EDO:H22	1.71	0.56
1:D:339:GLU:O	1:D:339:GLU:HG2	2.05	0.56
1:B:302:LYS:HZ2	2:B:1502:EDO:H22	1.70	0.56
2:D:1510:EDO:H22	4:D:2347:HOH:O	2.06	0.55
1:B:42:TRP:O	2:B:1514:EDO:C2	2.55	0.55
1:C:338[B]:CYS:SG	1:C:376:SER:OG	2.65	0.55
1:A:30:ASN:ND2	4:A:2029:HOH:O	2.39	0.55
1:C:177:ASN:ND2	2:C:1510:EDO:H11	2.20	0.55
1:A:436:SER:O	4:A:2411:HOH:O	2.18	0.55
2:D:1510:EDO:C2	4:D:2138:HOH:O	2.55	0.55
1:B:277:GLY:HA3	2:B:1518:EDO:H12	1.88	0.54
1:A:204:LYS:NZ	2:A:1509:EDO:H11	2.22	0.54
1:C:121:ASN:ND2	4:C:2157:HOH:O	2.41	0.54
1:D:0:SER:HA	1:D:71:SER:OG	2.06	0.54
1:B:419:LYS:HE3	4:B:2181:HOH:O	2.07	0.53
1:B:263:LYS:HD2	1:B:263:LYS:C	2.33	0.53
1:B:397:LEU:HD23	1:B:397:LEU:C	2.33	0.53
1:C:354:TYR:HD2	4:C:2360:HOH:O	1.90	0.53
1:D:405:MET:HE2	1:D:407:HIS:HA	1.90	0.53
2:B:1514:EDO:H22	4:B:2356:HOH:O	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:1516:EDO:H11	4:C:2487:HOH:O	2.08	0.53
1:C:338[B]:CYS:SG	1:C:376:SER:CB	2.96	0.53
1:A:136:PHE:O	1:A:140[B]:LYS:HE2	2.07	0.53
1:B:3:VAL:CG2	1:B:74:ILE:HD13	2.39	0.53
1:D:439:LYS:H	1:D:439:LYS:HD3	1.72	0.52
1:B:137:SER:HA	1:B:140[B]:LYS:HE2	1.92	0.52
1:D:460:LEU:HD13	1:D:464:LYS:CE	2.40	0.52
1:A:42:TRP:O	2:A:1507:EDO:H12	2.10	0.52
1:D:338[A]:CYS:SG	1:D:376[A]:SER:OG	2.68	0.52
1:D:483:ARG:O	1:D:487:THR:HG23	2.10	0.52
1:B:30:ASN:OD1	2:B:1521:EDO:H21	2.10	0.52
1:B:376:SER:HB3	4:B:2384:HOH:O	2.08	0.52
1:C:120:ASN:C	1:C:120:ASN:HD22	2.19	0.51
1:B:256[B]:ILE:HG23	1:B:261:GLU:HB3	1.92	0.51
1:C:452:LEU:HB3	2:C:1507:EDO:H11	1.93	0.51
1:D:256:ILE:HG22	1:D:256:ILE:O	2.09	0.51
1:B:302:LYS:NZ	2:B:1502:EDO:H22	2.27	0.50
1:B:67:ASP:HB2	4:B:2097:HOH:O	2.09	0.50
1:C:96:PRO:HB2	2:C:1511:EDO:H11	1.92	0.50
1:A:374:GLU:H	1:A:377:HIS:CD2	2.12	0.50
1:B:42:TRP:O	2:B:1514:EDO:C1	2.60	0.50
1:C:91[B]:ARG:NE	4:C:2116:HOH:O	2.43	0.50
1:A:1:MET:HE3	1:A:69:TYR:CE1	2.47	0.50
1:B:338[A]:CYS:SG	1:B:376:SER:OG	2.66	0.50
1:C:31:LEU:HD11	3:C:1523:GOL:H11	1.94	0.50
1:D:460:LEU:HD13	1:D:464:LYS:HE2	1.93	0.50
1:B:438:TYR:HE2	1:B:466:LEU:O	1.95	0.49
1:B:68:LYS:NZ	4:B:2101:HOH:O	2.35	0.49
1:A:101:ILE:HG21	1:A:140[A]:LYS:HG2	1.93	0.49
1:D:342:THR:HG23	1:D:377:HIS:ND1	2.28	0.49
1:B:42:TRP:O	2:B:1514:EDO:H12	2.12	0.49
1:D:354[A]:TYR:CD1	4:D:2322:HOH:O	2.49	0.49
1:C:236:LEU:HD22	2:C:1520:EDO:C1	2.41	0.49
1:B:0:SER:HA	1:B:71:SER:OG	2.13	0.49
1:B:374:GLU:H	1:B:377:HIS:CD2	2.09	0.49
1:D:468:ARG:NH2	4:D:2436:HOH:O	2.46	0.49
1:B:27:HIS:HE1	4:B:2096:HOH:O	1.96	0.48
1:D:291:ILE:HD13	1:D:303:TYR:CE1	2.48	0.48
1:D:272:LEU:HD23	1:D:273:LEU:N	2.29	0.48
1:A:342:THR:HG23	1:A:377:HIS:CE1	2.49	0.48
1:D:37:CYS:HB3	2:D:1510:EDO:H11	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:256:ILE:HG13	1:C:256:ILE:O	2.14	0.48
2:C:1510:EDO:H22	4:C:2481:HOH:O	2.13	0.47
1:D:42:TRP:O	2:D:1510:EDO:H12	2.13	0.47
1:D:464:LYS:HB3	1:D:468:ARG:HH11	1.79	0.47
1:A:415:MET:HG3	4:A:2284:HOH:O	2.13	0.47
1:D:84:GLU:HB2	1:D:103:TRP:HB3	1.96	0.47
1:A:482:LYS:HE3	1:A:486:LYS:HE2	1.97	0.47
1:A:492:LYS:HD3	4:A:2450:HOH:O	2.15	0.47
1:B:338[A]:CYS:SG	1:B:376:SER:HB2	2.54	0.47
1:C:48:ILE:HD12	2:C:1511:EDO:H22	1.95	0.47
1:A:42:TRP:O	2:A:1507:EDO:C1	2.63	0.47
1:D:414:GLY:O	1:D:417:LYS:HD3	2.15	0.47
1:B:51:MET:SD	1:B:169[B]:LEU:HD12	2.55	0.47
1:A:374:GLU:N	1:A:377:HIS:HD2	1.99	0.46
1:C:438:TYR:HH	1:C:466:LEU:CD2	2.25	0.46
3:C:1522:GOL:C2	4:C:2429:HOH:O	2.33	0.46
3:C:1525:GOL:C3	4:C:2495:HOH:O	2.63	0.46
1:B:483:ARG:O	1:B:487:THR:HG23	2.16	0.46
1:C:1:MET:HE2	1:C:69:TYR:CE1	2.50	0.46
2:D:1510:EDO:H21	4:D:2138:HOH:O	2.14	0.46
1:C:57:LEU:HA	1:C:60:GLU:HG2	1.96	0.46
1:D:-1:GLY:HA3	1:D:71:SER:H	1.81	0.46
1:C:296:ASN:ND2	2:C:1507:EDO:H22	2.31	0.45
1:B:270:VAL:HG23	1:B:312:ALA:CB	2.46	0.45
1:D:263:LYS:HD2	1:D:263:LYS:C	2.42	0.45
1:D:429:ILE:HG13	1:D:431:THR:HB	1.99	0.45
1:C:334:ILE:O	1:C:338[B]:CYS:HB2	2.16	0.45
1:C:484:LYS:HD2	1:C:484:LYS:N	2.30	0.45
1:B:152:LYS:HE3	4:B:2199:HOH:O	2.17	0.45
1:B:272:LEU:C	1:B:272:LEU:HD23	2.42	0.44
1:C:211:MET:HG3	3:C:1527:GOL:H31	2.00	0.44
1:C:122:ASN:HD22	1:C:122:ASN:H	1.66	0.44
1:C:66:LYS:HZ3	1:C:72:VAL:HB	1.82	0.44
1:C:374:GLU:N	1:C:377:HIS:HD2	2.03	0.44
1:D:140:LYS:HD3	4:D:2112:HOH:O	2.17	0.44
1:B:0:SER:CA	1:B:71:SER:OG	2.66	0.44
1:B:0:SER:HB3	1:B:71:SER:OG	2.17	0.44
1:C:270:VAL:HG23	1:C:312:ALA:CB	2.48	0.44
1:C:272:LEU:C	1:C:272:LEU:HD23	2.43	0.44
1:D:464:LYS:HB3	1:D:468:ARG:NH1	2.33	0.44
1:B:62:ILE:HD13	1:B:74:ILE:HD11	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:415:MET:HG3	4:B:2295:HOH:O	2.17	0.44
1:C:395[B]:ASP:HB3	2:C:1516:EDO:H21	1.99	0.44
1:A:39:LYS:HD2	1:A:42:TRP:CE2	2.53	0.43
1:B:338[A]:CYS:SG	1:B:376:SER:CB	3.06	0.43
1:C:263:LYS:HA	1:C:410:ARG:O	2.17	0.43
1:D:342:THR:CG2	1:D:377:HIS:CE1	3.00	0.43
1:B:259:GLU:OE1	2:B:1518:EDO:C1	2.67	0.43
1:C:42:TRP:O	2:C:1514:EDO:C1	2.66	0.43
1:A:117:ALA:HB1	3:A:1512:GOL:H32	1.98	0.43
1:C:47:PRO:HD2	2:C:1511:EDO:H12	2.01	0.43
1:D:302:LYS:HE3	4:D:2293:HOH:O	2.19	0.43
1:A:263:LYS:C	1:A:263:LYS:HD2	2.44	0.43
1:C:254:GLN:NE2	1:C:410:ARG:HE	2.17	0.43
1:B:86:VAL:HG12	1:B:140[B]:LYS:HE3	2.01	0.43
1:D:270:VAL:HG23	1:D:312:ALA:HB2	1.99	0.43
1:C:60:GLU:HG3	1:C:61:GLY:N	2.34	0.42
1:C:270:VAL:HG23	1:C:312:ALA:HB2	2.01	0.42
3:C:1526:GOL:H32	4:C:2498:HOH:O	2.19	0.42
1:C:62:ILE:HD13	1:C:74:ILE:HD11	2.01	0.42
1:C:74:ILE:HD13	1:C:235:TYR:CE1	2.54	0.42
1:B:259:GLU:OE1	2:B:1518:EDO:H11	2.19	0.42
1:B:291:ILE:HD13	1:B:303:TYR:CE1	2.53	0.42
1:D:345:ILE:HG21	1:D:493:ALA:CB	2.48	0.42
1:C:140:LYS:HE3	4:C:2110:HOH:O	2.05	0.42
1:D:121:ASN:HB3	1:D:131:TYR:CD1	2.54	0.42
1:A:84:GLU:HB2	1:A:103:TRP:HB3	2.02	0.42
1:A:294:LYS:HE3	1:A:299:ASP:O	2.19	0.42
1:C:252[A]:ILE:HD13	1:C:449:LEU:HD21	2.01	0.42
1:A:438:TYR:OH	1:A:469:ARG:HD2	2.19	0.42
1:D:108:VAL:HG22	4:D:2139:HOH:O	2.19	0.42
1:D:374:GLU:H	1:D:377:HIS:CD2	2.10	0.42
1:A:371:PHE:HD2	2:C:1514:EDO:C2	2.32	0.42
1:B:84:GLU:HB2	1:B:103:TRP:HB3	2.01	0.42
1:D:405:MET:HE2	1:D:406:LEU:O	2.20	0.42
1:A:345:ILE:HD11	1:A:489:GLU:HB3	2.02	0.42
1:B:417:LYS:NZ	4:B:2416:HOH:O	2.52	0.42
1:C:196:LEU:HD11	1:C:301:PRO:HG3	2.01	0.42
1:D:270:VAL:HG23	1:D:312:ALA:CB	2.50	0.42
1:C:56:ASN:HB3	3:C:1523:GOL:O1	2.20	0.41
1:D:74:ILE:O	1:D:238:ILE:HD12	2.20	0.41
1:B:466:LEU:HD12	1:B:469:ARG:NH2	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:479:ASP:OD2	1:B:481:LYS:HB2	2.20	0.41
2:B:1514:EDO:C2	4:B:2132:HOH:O	2.67	0.41
1:C:84:GLU:HB2	1:C:103:TRP:HB3	2.03	0.41
1:C:233:PRO:O	2:C:1520:EDO:C2	2.68	0.41
1:C:464:LYS:NZ	4:C:2437:HOH:O	2.52	0.41
1:B:196:LEU:HD11	1:B:301:PRO:HG3	2.02	0.41
1:B:339:GLU:HB3	4:B:2350:HOH:O	2.20	0.41
1:C:1:MET:HG3	1:C:3:VAL:HG13	2.02	0.41
1:A:204:LYS:HZ2	2:A:1509:EDO:H11	1.86	0.41
1:A:29:ASN:ND2	1:A:60:GLU:OE2	2.47	0.41
1:A:256:ILE:HG22	1:A:256:ILE:O	2.20	0.41
2:C:1510:EDO:H22	4:C:2220:HOH:O	2.20	0.41
3:C:1525:GOL:H31	4:C:2495:HOH:O	2.19	0.41
3:C:1527:GOL:H12	4:C:2242:HOH:O	2.19	0.41
2:D:1510:EDO:H22	4:D:2138:HOH:O	2.19	0.41
1:C:38[A]:LEU:HG	1:C:39:LYS:HG3	2.02	0.41
1:D:439:LYS:HD3	1:D:439:LYS:N	2.35	0.41
1:C:374:GLU:H	1:C:377:HIS:CD2	2.14	0.41
1:D:263:LYS:HA	1:D:410:ARG:O	2.21	0.41
1:A:27:HIS:HE1	4:A:2081:HOH:O	2.04	0.41
1:B:416:THR:C	1:B:417:LYS:HE3	2.46	0.41
1:B:9:GLN:HB2	1:B:169[A]:LEU:HD21	2.03	0.40
1:B:122:ASN:ND2	4:B:2167:HOH:O	2.53	0.40
1:D:65:LEU:HD23	1:D:65:LEU:HA	1.91	0.40
1:D:342:THR:HG23	1:D:377:HIS:CG	2.56	0.40
1:B:38:LEU:C	1:B:39:LYS:HG3	2.47	0.40
1:B:230:GLU:H	1:B:230:GLU:CD	2.29	0.40
1:A:412:ASP:CG	1:A:441:VAL:HG21	2.46	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:2181:HOH:O	4:B:2203:HOH:O[1_556]	2.00	0.20
4:B:2266:HOH:O	4:D:2117:HOH:O[1_565]	2.16	0.04

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	509/503 (101%)	496 (97%)	13 (3%)	0	100	100
1	B	508/503 (101%)	495 (97%)	13 (3%)	0	100	100
1	C	509/503 (101%)	497 (98%)	12 (2%)	0	100	100
1	D	508/503 (101%)	496 (98%)	12 (2%)	0	100	100
All	All	2034/2012 (101%)	1984 (98%)	50 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	459/450 (102%)	451 (98%)	8 (2%)	53	26
1	B	457/450 (102%)	446 (98%)	11 (2%)	43	15
1	C	459/450 (102%)	446 (97%)	13 (3%)	38	11
1	D	457/450 (102%)	445 (97%)	12 (3%)	40	13
All	All	1832/1800 (102%)	1788 (98%)	44 (2%)	44	15

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	60	GLU
1	A	74	ILE

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Mol	Chain	Res	Type
1	A	169	LEU
1	A	189	LEU
1	A	254	GLN
1	A	419	LYS
1	A	461	ASP
1	A	484	LYS
1	B	70	THR
1	B	140[A]	LYS
1	B	140[B]	LYS
1	B	169[A]	LEU
1	B	169[B]	LEU
1	B	189	LEU
1	B	256[A]	ILE
1	B	256[B]	ILE
1	B	417	LYS
1	B	462	SER
1	B	466	LEU
1	C	60	GLU
1	C	95	LYS
1	C	120	ASN
1	C	122	ASN
1	C	140	LYS
1	C	189	LEU
1	C	254	GLN
1	C	404	GLU
1	C	439	LYS
1	C	462	SER
1	C	466	LEU
1	C	484	LYS
1	C	492	LYS
1	D	0	SER
1	D	38	LEU
1	D	189	LEU
1	D	254	GLN
1	D	339	GLU
1	D	397	LEU
1	D	405	MET
1	D	417	LYS
1	D	439	LYS
1	D	455	LYS
1	D	460	LEU
1	D	461	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (47) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	12	GLN
1	A	24	ASN
1	A	27	HIS
1	A	30	ASN
1	A	81	ASN
1	A	99	ASN
1	A	125	GLN
1	A	172	ASN
1	A	184	ASN
1	A	254	GLN
1	A	322	ASN
1	A	372	ASN
1	A	377	HIS
1	B	24	ASN
1	B	27	HIS
1	B	99	ASN
1	B	122	ASN
1	B	153	GLN
1	B	172	ASN
1	B	322	ASN
1	B	372	ASN
1	B	377	HIS
1	B	423	GLN
1	B	500	GLN
1	C	24	ASN
1	C	27	HIS
1	C	35	GLN
1	C	81	ASN
1	C	82	GLN
1	C	120	ASN
1	C	122	ASN
1	C	172	ASN
1	C	177	ASN
1	C	184	ASN
1	C	254	GLN
1	C	322	ASN
1	C	377	HIS
1	C	475	HIS
1	C	500	GLN
1	D	24	ASN
1	D	27	HIS

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Mol	Chain	Res	Type
1	D	125	GLN
1	D	172	ASN
1	D	254	GLN
1	D	322	ASN
1	D	372	ASN
1	D	475	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](#)

74 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	EDO	C	1511	-	3,3,3	0.59	0	2,2,2	0.45	0
2	EDO	A	1509	-	3,3,3	0.60	0	2,2,2	0.42	0
2	EDO	A	1510	-	3,3,3	0.18	0	2,2,2	0.23	0
2	EDO	B	1508	-	3,3,3	0.56	0	2,2,2	0.74	0
2	EDO	B	1512	-	3,3,3	0.36	0	2,2,2	0.24	0
2	EDO	D	1512	-	3,3,3	0.73	0	2,2,2	0.05	0
2	EDO	A	1504	-	3,3,3	0.08	0	2,2,2	0.86	0
2	EDO	C	1519	-	3,3,3	0.67	0	2,2,2	0.44	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	EDO	A	1503	-	3,3,3	0.43	0	2,2,2	0.61	0
2	EDO	C	1505	-	3,3,3	0.40	0	2,2,2	1.08	0
3	GOL	C	1523	-	5,5,5	0.44	0	5,5,5	0.21	0
2	EDO	A	1505	-	3,3,3	0.35	0	2,2,2	0.09	0
2	EDO	B	1511	-	3,3,3	0.38	0	2,2,2	0.27	0
2	EDO	B	1510	-	3,3,3	0.84	0	2,2,2	0.66	0
2	EDO	D	1502	-	3,3,3	0.62	0	2,2,2	1.04	0
2	EDO	C	1506	-	3,3,3	0.59	0	2,2,2	0.68	0
2	EDO	A	1508	-	3,3,3	0.78	0	2,2,2	1.51	1 (50%)
2	EDO	A	1506	-	3,3,3	0.65	0	2,2,2	0.55	0
2	EDO	D	1504	-	3,3,3	0.51	0	2,2,2	0.82	0
2	EDO	D	1513	-	3,3,3	0.38	0	2,2,2	0.69	0
2	EDO	C	1512	-	3,3,3	0.37	0	2,2,2	0.22	0
2	EDO	C	1520	-	3,3,3	0.38	0	2,2,2	0.15	0
2	EDO	D	1514	-	3,3,3	0.29	0	2,2,2	0.88	0
2	EDO	C	1503	-	3,3,3	0.47	0	2,2,2	1.09	0
2	EDO	A	1507	-	3,3,3	1.21	0	2,2,2	1.52	0
2	EDO	B	1516	-	3,3,3	1.06	0	2,2,2	0.19	0
3	GOL	C	1525	-	5,5,5	0.41	0	5,5,5	1.11	0
3	GOL	A	1511	-	5,5,5	1.04	0	5,5,5	0.87	0
2	EDO	B	1509	-	3,3,3	0.27	0	2,2,2	1.68	1 (50%)
2	EDO	B	1513	-	3,3,3	0.34	0	2,2,2	0.33	0
2	EDO	B	1517	-	3,3,3	0.51	0	2,2,2	0.54	0
2	EDO	B	1506	-	3,3,3	1.01	0	2,2,2	0.75	0
2	EDO	B	1515	-	3,3,3	0.46	0	2,2,2	0.21	0
2	EDO	C	1518	-	3,3,3	0.22	0	2,2,2	0.72	0
2	EDO	B	1503	-	3,3,3	0.41	0	2,2,2	0.61	0
2	EDO	D	1505	-	3,3,3	0.17	0	2,2,2	1.25	0
2	EDO	D	1507	-	3,3,3	0.84	0	2,2,2	0.92	0
3	GOL	D	1516	-	5,5,5	0.91	0	5,5,5	0.89	0
2	EDO	B	1514	-	3,3,3	0.80	0	2,2,2	1.04	0
2	EDO	C	1502	-	3,3,3	0.66	0	2,2,2	0.74	0
3	GOL	C	1527	-	5,5,5	0.38	0	5,5,5	1.17	0
3	GOL	C	1521	-	5,5,5	0.98	0	5,5,5	0.53	0
2	EDO	B	1521	-	3,3,3	0.53	0	2,2,2	0.35	0
2	EDO	B	1507	-	3,3,3	0.46	0	2,2,2	0.35	0
2	EDO	C	1509	-	3,3,3	1.89	1 (33%)	2,2,2	1.84	1 (50%)
3	GOL	C	1524	-	5,5,5	0.69	0	5,5,5	1.99	2 (40%)
3	GOL	C	1526	-	5,5,5	0.67	0	5,5,5	0.44	0
3	GOL	B	1523	-	5,5,5	1.64	1 (20%)	5,5,5	1.19	0
2	EDO	B	1504	-	3,3,3	0.69	0	2,2,2	1.01	0
2	EDO	C	1510	-	3,3,3	2.63	1 (33%)	2,2,2	1.38	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	EDO	C	1516	-	3,3,3	1.05	0	2,2,2	1.26	0
2	EDO	B	1519	-	3,3,3	0.30	0	2,2,2	0.87	0
2	EDO	D	1509	-	3,3,3	0.37	0	2,2,2	0.03	0
2	EDO	C	1504	-	3,3,3	0.66	0	2,2,2	0.99	0
2	EDO	C	1517	-	3,3,3	0.44	0	2,2,2	0.43	0
2	EDO	C	1508	-	3,3,3	0.72	0	2,2,2	0.17	0
2	EDO	C	1513	-	3,3,3	0.30	0	2,2,2	0.32	0
2	EDO	C	1515	-	3,3,3	0.80	0	2,2,2	0.33	0
3	GOL	C	1522	-	5,5,5	1.18	1 (20%)	5,5,5	1.29	0
2	EDO	B	1502	-	3,3,3	0.56	0	2,2,2	1.02	0
2	EDO	C	1514	-	3,3,3	2.21	1 (33%)	2,2,2	2.45	2 (100%)
2	EDO	A	1502	-	3,3,3	0.39	0	2,2,2	1.06	0
2	EDO	D	1510	-	3,3,3	0.90	0	2,2,2	1.22	0
2	EDO	B	1518	-	3,3,3	0.62	0	2,2,2	0.44	0
2	EDO	D	1506	-	3,3,3	0.21	0	2,2,2	0.86	0
2	EDO	B	1520	-	3,3,3	1.13	0	2,2,2	0.30	0
3	GOL	A	1512	-	5,5,5	0.47	0	5,5,5	0.64	0
2	EDO	D	1508	-	3,3,3	0.72	0	2,2,2	0.37	0
2	EDO	B	1505	-	3,3,3	0.14	0	2,2,2	1.14	0
3	GOL	B	1522	-	5,5,5	0.75	0	5,5,5	0.52	0
2	EDO	D	1515	-	3,3,3	3.93	2 (66%)	2,2,2	2.08	1 (50%)
2	EDO	D	1503	-	3,3,3	0.46	0	2,2,2	0.73	0
2	EDO	C	1507	-	3,3,3	0.37	0	2,2,2	0.41	0
2	EDO	D	1511	-	3,3,3	0.37	0	2,2,2	0.28	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
2	EDO	C	1511	-	-	1/1/1/1	-
2	EDO	A	1509	-	-	1/1/1/1	-
2	EDO	A	1510	-	-	0/1/1/1	-
2	EDO	B	1508	-	-	1/1/1/1	-
2	EDO	B	1512	-	-	1/1/1/1	-
2	EDO	D	1512	-	-	1/1/1/1	-
2	EDO	A	1504	-	-	0/1/1/1	-
2	EDO	C	1519	-	-	0/1/1/1	-
2	EDO	A	1503	-	-	0/1/1/1	-
2	EDO	C	1505	-	-	0/1/1/1	-
3	GOL	C	1523	-	-	3/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	EDO	A	1505	-	-	0/1/1/1	-
2	EDO	B	1511	-	-	0/1/1/1	-
2	EDO	B	1510	-	-	0/1/1/1	-
2	EDO	D	1502	-	-	0/1/1/1	-
2	EDO	C	1506	-	-	1/1/1/1	-
2	EDO	A	1508	-	-	1/1/1/1	-
2	EDO	A	1506	-	-	0/1/1/1	-
2	EDO	D	1504	-	-	0/1/1/1	-
2	EDO	D	1513	-	-	1/1/1/1	-
2	EDO	C	1512	-	-	0/1/1/1	-
2	EDO	C	1520	-	-	0/1/1/1	-
2	EDO	D	1514	-	-	0/1/1/1	-
2	EDO	C	1503	-	-	0/1/1/1	-
2	EDO	A	1507	-	-	1/1/1/1	-
2	EDO	B	1516	-	-	0/1/1/1	-
3	GOL	C	1525	-	-	2/4/4/4	-
3	GOL	A	1511	-	-	0/4/4/4	-
2	EDO	B	1509	-	-	1/1/1/1	-
2	EDO	B	1513	-	-	1/1/1/1	-
2	EDO	B	1517	-	-	1/1/1/1	-
2	EDO	B	1506	-	-	0/1/1/1	-
2	EDO	B	1515	-	-	0/1/1/1	-
2	EDO	C	1518	-	-	0/1/1/1	-
2	EDO	B	1503	-	-	0/1/1/1	-
2	EDO	D	1505	-	-	0/1/1/1	-
2	EDO	D	1507	-	-	0/1/1/1	-
3	GOL	D	1516	-	-	0/4/4/4	-
2	EDO	B	1514	-	-	1/1/1/1	-
2	EDO	C	1502	-	-	0/1/1/1	-
3	GOL	C	1527	-	-	2/4/4/4	-
3	GOL	C	1521	-	-	0/4/4/4	-
2	EDO	B	1521	-	-	0/1/1/1	-
2	EDO	B	1507	-	-	1/1/1/1	-
2	EDO	C	1509	-	-	0/1/1/1	-
3	GOL	C	1524	-	-	0/4/4/4	-
3	GOL	C	1526	-	-	2/4/4/4	-
3	GOL	B	1523	-	-	4/4/4/4	-
2	EDO	B	1504	-	-	0/1/1/1	-
2	EDO	C	1510	-	-	1/1/1/1	-
2	EDO	C	1516	-	-	1/1/1/1	-
2	EDO	B	1519	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	EDO	D	1509	-	-	0/1/1/1	-
2	EDO	C	1504	-	-	1/1/1/1	-
2	EDO	C	1517	-	-	1/1/1/1	-
2	EDO	C	1508	-	-	0/1/1/1	-
2	EDO	C	1513	-	-	0/1/1/1	-
2	EDO	C	1515	-	-	0/1/1/1	-
3	GOL	C	1522	-	-	2/4/4/4	-
2	EDO	B	1502	-	-	0/1/1/1	-
2	EDO	C	1514	-	-	1/1/1/1	-
2	EDO	A	1502	-	-	0/1/1/1	-
2	EDO	D	1510	-	-	1/1/1/1	-
2	EDO	B	1518	-	-	1/1/1/1	-
2	EDO	D	1506	-	-	0/1/1/1	-
2	EDO	B	1520	-	-	1/1/1/1	-
3	GOL	A	1512	-	-	4/4/4/4	-
2	EDO	D	1508	-	-	1/1/1/1	-
2	EDO	B	1505	-	-	0/1/1/1	-
3	GOL	B	1522	-	-	0/4/4/4	-
2	EDO	D	1515	-	-	1/1/1/1	-
2	EDO	D	1503	-	-	0/1/1/1	-
2	EDO	C	1507	-	-	1/1/1/1	-
2	EDO	D	1511	-	-	0/1/1/1	-

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1515	EDO	O1-C1	-5.75	1.12	1.42
2	C	1510	EDO	O1-C1	4.38	1.64	1.42
2	D	1515	EDO	O2-C2	-3.64	1.23	1.42
2	C	1514	EDO	O1-C1	3.08	1.57	1.42
3	B	1523	GOL	O2-C2	2.40	1.50	1.43
3	C	1522	GOL	O2-C2	2.25	1.49	1.43
2	C	1509	EDO	O1-C1	2.21	1.53	1.42

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1524	GOL	O2-C2-C1	-3.05	96.55	109.18
2	C	1514	EDO	O2-C2-C1	2.68	132.76	112.39
2	D	1515	EDO	O2-C2-C1	-2.64	92.28	112.39
2	C	1509	EDO	O1-C1-C2	2.40	130.67	112.39
2	C	1514	EDO	O1-C1-C2	2.21	129.18	112.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1524	GOL	O2-C2-C3	2.15	118.09	109.18
2	A	1508	EDO	O2-C2-C1	-2.13	96.19	112.39
2	B	1509	EDO	O1-C1-C2	-2.03	96.92	112.39

There are no chirality outliers.

All (45) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1512	GOL	O1-C1-C2-C3
3	A	1512	GOL	C1-C2-C3-O3
3	B	1523	GOL	O1-C1-C2-C3
3	B	1523	GOL	C1-C2-C3-O3
3	C	1523	GOL	O1-C1-C2-C3
3	C	1525	GOL	C1-C2-C3-O3
3	C	1526	GOL	C1-C2-C3-O3
3	C	1522	GOL	O1-C1-C2-C3
3	C	1527	GOL	O1-C1-C2-C3
3	A	1512	GOL	O1-C1-C2-O2
3	A	1512	GOL	O2-C2-C3-O3
3	C	1525	GOL	O2-C2-C3-O3
3	C	1526	GOL	O2-C2-C3-O3
2	C	1510	EDO	O1-C1-C2-O2
2	C	1516	EDO	O1-C1-C2-O2
2	D	1512	EDO	O1-C1-C2-O2
3	B	1523	GOL	O1-C1-C2-O2
3	B	1523	GOL	O2-C2-C3-O3
3	C	1523	GOL	O1-C1-C2-O2
2	A	1507	EDO	O1-C1-C2-O2
2	B	1518	EDO	O1-C1-C2-O2
2	B	1519	EDO	O1-C1-C2-O2
2	C	1507	EDO	O1-C1-C2-O2
3	C	1522	GOL	O2-C2-C3-O3
2	A	1509	EDO	O1-C1-C2-O2
3	C	1523	GOL	C1-C2-C3-O3
2	B	1514	EDO	O1-C1-C2-O2
2	D	1513	EDO	O1-C1-C2-O2
3	C	1527	GOL	O1-C1-C2-O2
2	D	1515	EDO	O1-C1-C2-O2
2	B	1512	EDO	O1-C1-C2-O2
2	B	1517	EDO	O1-C1-C2-O2
2	C	1511	EDO	O1-C1-C2-O2
2	A	1508	EDO	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
2	B	1507	EDO	O1-C1-C2-O2
2	B	1513	EDO	O1-C1-C2-O2
2	C	1504	EDO	O1-C1-C2-O2
2	D	1508	EDO	O1-C1-C2-O2
2	D	1510	EDO	O1-C1-C2-O2
2	B	1508	EDO	O1-C1-C2-O2
2	B	1520	EDO	O1-C1-C2-O2
2	C	1517	EDO	O1-C1-C2-O2
2	B	1509	EDO	O1-C1-C2-O2
2	C	1506	EDO	O1-C1-C2-O2
2	C	1514	EDO	O1-C1-C2-O2

There are no ring outliers.

25 monomers are involved in 84 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1511	EDO	3	0
2	A	1509	EDO	4	0
3	C	1523	GOL	3	0
2	A	1508	EDO	1	0
2	C	1520	EDO	4	0
2	D	1514	EDO	1	0
2	A	1507	EDO	4	0
3	C	1525	GOL	2	0
2	B	1517	EDO	1	0
2	B	1514	EDO	6	0
3	C	1527	GOL	6	0
2	B	1521	EDO	2	0
3	C	1526	GOL	1	0
3	B	1523	GOL	1	0
2	C	1510	EDO	8	0
2	C	1516	EDO	3	0
2	B	1519	EDO	1	0
3	C	1522	GOL	2	0
2	B	1502	EDO	2	0
2	C	1514	EDO	6	0
2	D	1510	EDO	6	0
2	B	1518	EDO	4	0
3	A	1512	GOL	3	0
2	D	1515	EDO	7	0
2	C	1507	EDO	3	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	501/503 (99%)	-0.19	20 (3%)	42 46	5, 13, 33, 44	14 (2%)
1	B	503/503 (100%)	-0.29	15 (2%)	52 56	4, 12, 30, 40	10 (1%)
1	C	502/503 (99%)	-0.31	13 (2%)	57 61	6, 12, 28, 40	17 (3%)
1	D	503/503 (100%)	-0.22	13 (2%)	57 61	5, 13, 31, 44	10 (1%)
All	All	2009/2012 (99%)	-0.25	61 (3%)	52 56	4, 13, 32, 44	51 (2%)

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	-1	GLY	8.3
1	C	0	SER	6.7
1	D	-1	GLY	6.3
1	A	417	LYS	5.8
1	D	0	SER	5.5
1	C	438	TYR	4.6
1	C	70	THR	4.5
1	B	70	THR	4.2
1	B	0	SER	4.2
1	D	70	THR	3.8
1	B	331	ALA	3.5
1	B	438	TYR	3.3
1	A	466	LEU	3.2
1	B	466	LEU	3.2
1	A	70	THR	3.1
1	A	342	THR	3.1
1	D	460	LEU	3.1
1	C	466	LEU	2.9
1	A	438	TYR	2.9
1	D	67	ASP	2.8
1	D	463	VAL	2.8

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Mol	Chain	Res	Type	RSRZ
1	C	463	VAL	2.7
1	B	458	ASP	2.7
1	D	342	THR	2.6
1	D	438	TYR	2.6
1	C	439	LYS	2.5
1	C	331	ALA	2.5
1	D	407	HIS	2.5
1	C	67	ASP	2.4
1	A	465	SER	2.3
1	C	480	ASP	2.3
1	A	331	ALA	2.3
1	A	332	SER	2.3
1	A	336	GLU	2.3
1	B	462	SER	2.3
1	A	333	ASP	2.2
1	D	458	ASP	2.2
1	C	408	VAL	2.2
1	D	462	SER	2.2
1	B	67	ASP	2.2
1	A	67	ASP	2.2
1	B	480	ASP	2.2
1	A	407	HIS	2.2
1	C	465	SER	2.2
1	A	480	ASP	2.2
1	C	458	ASP	2.2
1	A	406	LEU	2.1
1	B	460	LEU	2.1
1	A	458	ASP	2.1
1	B	439	LYS	2.1
1	D	405	MET	2.1
1	A	405	MET	2.1
1	A	501	LEU	2.1
1	A	460	LEU	2.0
1	D	466	LEU	2.0
1	A	69	TYR	2.0
1	B	332	SER	2.0
1	B	459	SER	2.0
1	A	335	MET	2.0
1	C	405	MET	2.0
1	B	461	ASP	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	GOL	C	1527	6/6	0.62	0.18	31,36,39,40	0
3	GOL	C	1526	6/6	0.70	0.15	32,34,36,41	0
2	EDO	C	1517	4/4	0.73	0.21	54,54,55,55	0
2	EDO	C	1506	4/4	0.75	0.15	38,43,44,44	0
3	GOL	C	1525	6/6	0.75	0.18	42,46,48,51	0
3	GOL	A	1512	6/6	0.77	0.17	48,49,51,53	0
3	GOL	C	1523	6/6	0.79	0.17	58,60,61,61	0
2	EDO	D	1513	4/4	0.79	0.18	37,38,40,47	0
2	EDO	C	1509	4/4	0.80	0.17	24,31,31,32	0
2	EDO	D	1511	4/4	0.80	0.15	40,40,43,50	0
2	EDO	B	1517	4/4	0.81	0.15	30,31,32,33	0
2	EDO	D	1514	4/4	0.82	0.17	45,45,45,51	0
2	EDO	C	1519	4/4	0.82	0.16	30,32,37,41	0
2	EDO	D	1508	4/4	0.82	0.15	33,34,36,38	0
2	EDO	A	1509	4/4	0.82	0.15	39,45,46,46	0
2	EDO	D	1512	4/4	0.82	0.14	34,36,36,38	0
2	EDO	B	1512	4/4	0.82	0.16	43,43,44,44	0
2	EDO	B	1516	4/4	0.83	0.15	29,30,33,34	0
2	EDO	C	1510	4/4	0.84	0.16	11,18,24,28	0
3	GOL	C	1524	6/6	0.85	0.13	20,34,39,43	0
2	EDO	A	1504	4/4	0.85	0.13	25,30,30,35	0
2	EDO	A	1508	4/4	0.85	0.12	18,31,35,36	0
2	EDO	B	1521	4/4	0.85	0.12	24,27,38,38	0
2	EDO	C	1503	4/4	0.86	0.12	21,24,24,25	0
2	EDO	C	1513	4/4	0.87	0.12	39,40,40,40	0
2	EDO	C	1516	4/4	0.87	0.17	22,26,27,36	0
2	EDO	A	1507	4/4	0.87	0.37	19,25,27,40	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	EDO	B	1514	4/4	0.88	0.39	19,20,26,39	0
2	EDO	B	1515	4/4	0.88	0.13	36,36,37,39	0
2	EDO	D	1510	4/4	0.88	0.15	22,26,27,41	0
2	EDO	B	1507	4/4	0.88	0.11	29,32,37,40	0
3	GOL	C	1522	6/6	0.88	0.10	15,21,23,26	0
2	EDO	C	1515	4/4	0.89	0.12	19,25,26,32	0
2	EDO	C	1505	4/4	0.89	0.11	22,23,28,32	0
2	EDO	B	1509	4/4	0.89	0.12	33,34,35,36	0
3	GOL	B	1523	6/6	0.89	0.12	15,23,29,29	0
2	EDO	C	1507	4/4	0.89	0.13	32,33,37,40	0
2	EDO	A	1505	4/4	0.90	0.11	22,28,31,37	0
2	EDO	B	1502	4/4	0.90	0.12	18,27,28,33	0
2	EDO	B	1506	4/4	0.90	0.16	15,21,24,34	0
2	EDO	C	1504	4/4	0.91	0.09	22,27,28,33	0
2	EDO	C	1514	4/4	0.91	0.22	18,20,21,28	0
2	EDO	C	1520	4/4	0.91	0.21	25,26,28,32	0
2	EDO	D	1506	4/4	0.91	0.12	19,22,26,32	0
2	EDO	B	1518	4/4	0.91	0.11	35,37,41,43	0
2	EDO	B	1520	4/4	0.91	0.12	17,23,23,32	0
2	EDO	B	1504	4/4	0.92	0.09	20,21,21,23	0
2	EDO	B	1505	4/4	0.92	0.12	21,27,30,35	0
2	EDO	D	1505	4/4	0.92	0.10	25,28,30,38	0
2	EDO	B	1519	4/4	0.92	0.11	25,35,36,38	0
2	EDO	B	1513	4/4	0.92	0.09	21,31,35,39	0
2	EDO	C	1511	4/4	0.92	0.12	23,31,33,35	0
2	EDO	D	1507	4/4	0.93	0.11	17,18,20,20	0
2	EDO	B	1511	4/4	0.93	0.10	23,32,35,38	0
2	EDO	B	1510	4/4	0.93	0.09	29,31,31,38	0
2	EDO	A	1503	4/4	0.94	0.07	18,18,20,21	0
2	EDO	B	1508	4/4	0.94	0.10	14,19,19,21	0
2	EDO	D	1504	4/4	0.94	0.07	17,18,20,21	0
2	EDO	A	1510	4/4	0.94	0.17	15,22,23,26	0
2	EDO	A	1506	4/4	0.94	0.08	15,16,16,18	0
2	EDO	A	1502	4/4	0.95	0.07	16,16,17,21	0
2	EDO	D	1515	4/4	0.95	0.15	10,16,18,18	0
2	EDO	B	1503	4/4	0.96	0.07	14,15,16,16	0
2	EDO	C	1508	4/4	0.96	0.07	12,15,16,18	0
3	GOL	A	1511	6/6	0.96	0.07	8,9,15,21	0
2	EDO	D	1503	4/4	0.96	0.09	14,16,21,22	0
3	GOL	B	1522	6/6	0.96	0.08	7,8,14,20	0
2	EDO	C	1518	4/4	0.96	0.07	25,29,30,32	0
2	EDO	C	1512	4/4	0.97	0.17	12,19,23,28	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	EDO	D	1502	4/4	0.97	0.05	15,15,15,18	0
3	GOL	C	1521	6/6	0.97	0.06	7,7,15,18	0
2	EDO	D	1509	4/4	0.98	0.13	13,18,19,25	0
2	EDO	C	1502	4/4	0.98	0.05	12,14,15,15	0
3	GOL	D	1516	6/6	0.98	0.07	6,8,16,22	0

6.5 Other polymers [i](#)

There are no such residues in this entry.