



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:11 AM UTC

PDB ID : 2YMW / pdb_00002ymw
Title : Structure of the epsilon-lysine oxidase from *Marinomonas mediterranea*
Authors : Medrano, F.J.; Romero, A.
Deposited on : 2012-10-10
Resolution : 2.41 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

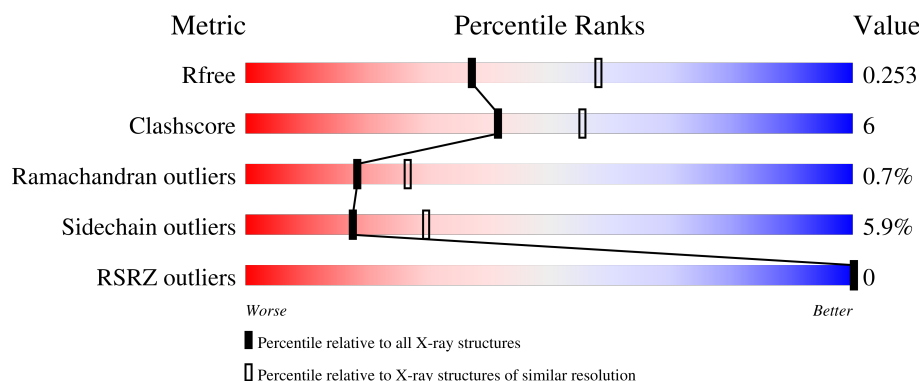
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.41 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	6062 (2.44-2.40)
Clashscore	190562	6562 (2.44-2.40)
Ramachandran outliers	187476	6481 (2.44-2.40)
Sidechain outliers	187428	6482 (2.44-2.40)
RSRZ outliers	180081	6066 (2.44-2.40)

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	746	
1	B	746	

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	EOH	A	1691	-	-	X	-
3	EOH	A	1696	-	-	X	-
3	EOH	A	1720	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 11836 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 L-LYSINE 6-OXIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	680	Total	C	N	O	S	0	0	0
			5348	3361	888	1076	23			
1	B	685	Total	C	N	O	S	0	0	0
			5378	3379	893	1083	23			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP F2JXJ3
A	-18	GLY	-	expression tag	UNP F2JXJ3
A	-17	SER	-	expression tag	UNP F2JXJ3
A	-16	SER	-	expression tag	UNP F2JXJ3
A	-15	HIS	-	expression tag	UNP F2JXJ3
A	-14	HIS	-	expression tag	UNP F2JXJ3
A	-13	HIS	-	expression tag	UNP F2JXJ3
A	-12	HIS	-	expression tag	UNP F2JXJ3
A	-11	HIS	-	expression tag	UNP F2JXJ3
A	-10	HIS	-	expression tag	UNP F2JXJ3
A	-9	SER	-	expression tag	UNP F2JXJ3
A	-8	SER	-	expression tag	UNP F2JXJ3
A	-7	GLY	-	expression tag	UNP F2JXJ3
A	-6	LEU	-	expression tag	UNP F2JXJ3
A	-5	VAL	-	expression tag	UNP F2JXJ3
A	-4	PRO	-	expression tag	UNP F2JXJ3
A	-3	ARG	-	expression tag	UNP F2JXJ3
A	-2	GLY	-	expression tag	UNP F2JXJ3
A	-1	SER	-	expression tag	UNP F2JXJ3
A	0	HIS	-	expression tag	UNP F2JXJ3
B	-19	MET	-	expression tag	UNP F2JXJ3
B	-18	GLY	-	expression tag	UNP F2JXJ3
B	-17	SER	-	expression tag	UNP F2JXJ3
B	-16	SER	-	expression tag	UNP F2JXJ3
B	-15	HIS	-	expression tag	UNP F2JXJ3

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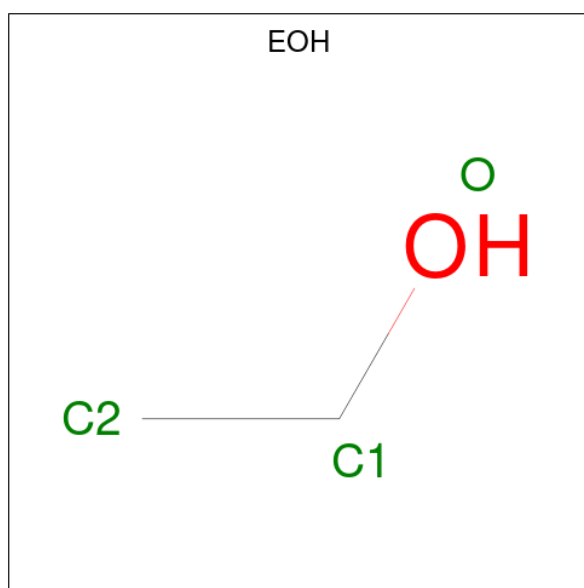
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Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	HIS	-	expression tag	UNP F2JXJ3
B	-13	HIS	-	expression tag	UNP F2JXJ3
B	-12	HIS	-	expression tag	UNP F2JXJ3
B	-11	HIS	-	expression tag	UNP F2JXJ3
B	-10	HIS	-	expression tag	UNP F2JXJ3
B	-9	SER	-	expression tag	UNP F2JXJ3
B	-8	SER	-	expression tag	UNP F2JXJ3
B	-7	GLY	-	expression tag	UNP F2JXJ3
B	-6	LEU	-	expression tag	UNP F2JXJ3
B	-5	VAL	-	expression tag	UNP F2JXJ3
B	-4	PRO	-	expression tag	UNP F2JXJ3
B	-3	ARG	-	expression tag	UNP F2JXJ3
B	-2	GLY	-	expression tag	UNP F2JXJ3
B	-1	SER	-	expression tag	UNP F2JXJ3
B	0	HIS	-	expression tag	UNP F2JXJ3

- Molecule 2 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	2	Total Na 2 2	0	0
2	B	2	Total Na 2 2	0	0

- Molecule 3 is ETHANOL (CCD ID: EOH) (formula: C₂H₆O).



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

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

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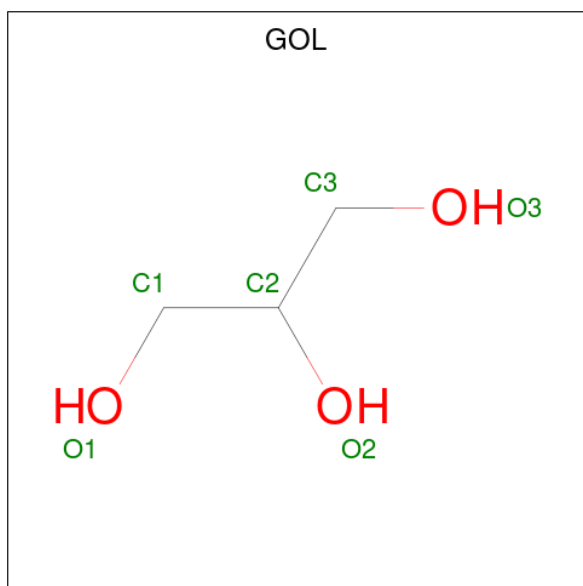
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			3	2	1		
3	B	1	Total	C	O	0	0
			3	2	1		
3	B	1	Total	C	O	0	0
			3	2	1		
3	B	1	Total	C	O	0	0
			3	2	1		
3	B	1	Total	C	O	0	0
			3	2	1		
3	B	1	Total	C	O	0	0
			3	2	1		
3	B	1	Total	C	O	0	0
			3	2	1		
3	B	1	Total	C	O	0	0
			3	2	1		
3	B	1	Total	C	O	0	0
			3	2	1		
3	B	1	Total	C	O	0	0
			3	2	1		
3	B	1	Total	C	O	0	0
			3	2	1		
3	B	1	Total	C	O	0	0
			3	2	1		
3	B	1	Total	C	O	0	0
			3	2	1		
3	B	1	Total	C	O	0	0
			3	2	1		
3	B	1	Total	C	O	0	0
			3	2	1		
3	B	1	Total	C	O	0	0
			3	2	1		
3	B	1	Total	C	O	0	0
			3	2	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			3	2	1		
3	B	1	Total	C	O	0	0
			3	2	1		
3	B	1	Total	C	O	0	0
			3	2	1		
3	B	1	Total	C	O	0	0
			3	2	1		
3	B	1	Total	C	O	0	0
			3	2	1		
3	B	1	Total	C	O	0	0
			3	2	1		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

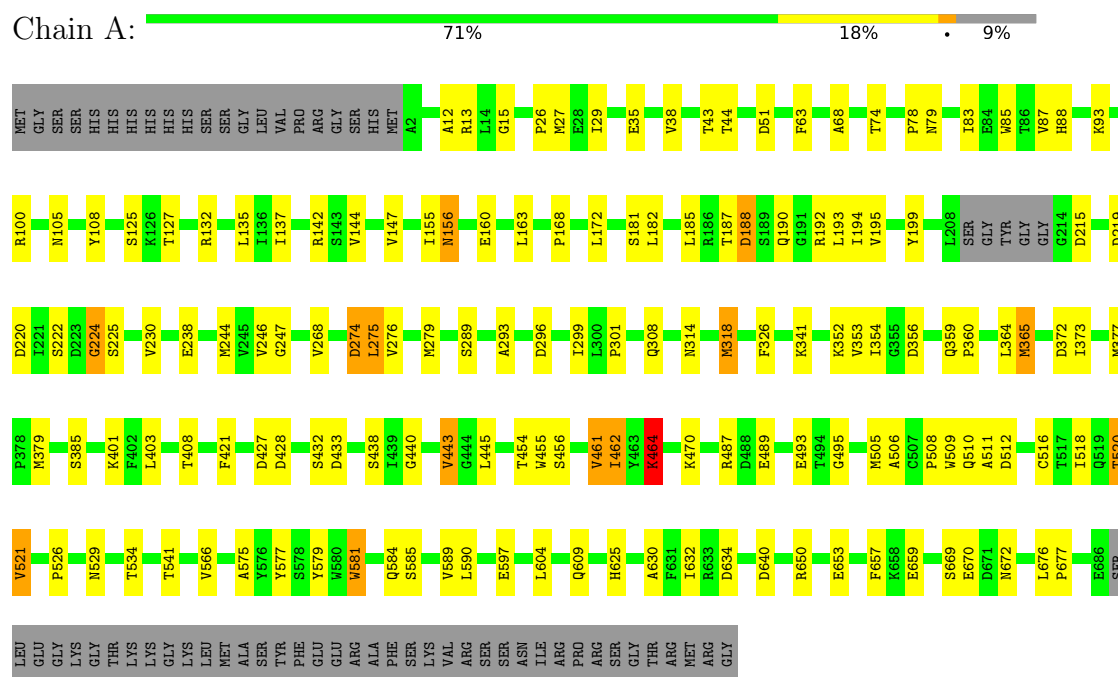
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	473	Total 473	O 473	0	0
5	B	408	Total 408	O 408	0	0

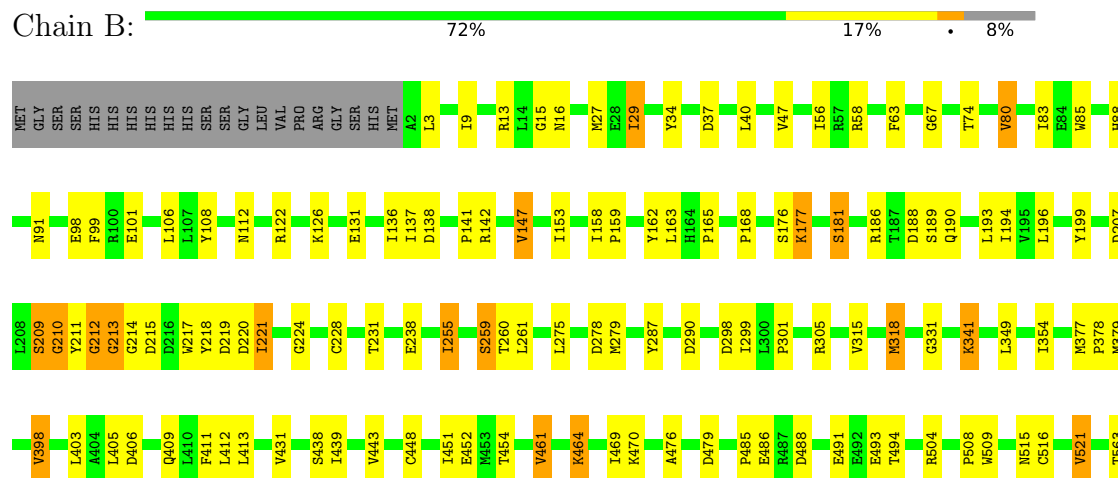
3 Residue-property plots

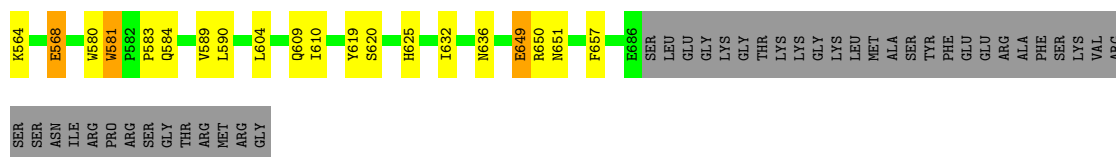
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: L-LYSINE 6-OXIDASE



• Molecule 1: L-LYSINE 6-OXIDASE





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	70.68Å 127.88Å 106.90Å 90.00° 107.18° 90.00°	Depositor
Resolution (Å)	35.28 – 2.41 35.28 – 2.41	Depositor EDS
% Data completeness (in resolution range)	98.3 (35.28-2.41) 89.7 (35.28-2.41)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.92 (at 2.42Å)	Xtriage
Refinement program	BUSTER 2.10.0	Depositor
R, R_{free}	0.192 , 0.279 (Not available) , 0.253	Depositor DCC
R_{free} test set	3476 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	34.7	Xtriage
Anisotropy	0.787	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 50.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	0.097 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11836	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.71% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: TRQ, NA, EOH, 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	0.90	2/5466 (0.0%)	1.36	31/7445 (0.4%)
1	B	0.87	1/5498 (0.0%)	1.38	40/7489 (0.5%)
All	All	0.88	3/10964 (0.0%)	1.37	71/14934 (0.5%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	318	MET	SD-CE	-7.13	1.61	1.79
1	A	43	THR	CA-C	5.54	1.55	1.52
1	B	318	MET	SD-CE	-5.35	1.66	1.79

All (71) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	211	TYR	N-CA-C	-10.15	101.15	112.72
1	B	219	ASP	CA-CB-CG	9.05	121.65	112.60
1	A	354	ILE	N-CA-C	9.00	119.80	110.62
1	A	155	ILE	N-CA-C	-8.10	99.14	109.58
1	A	625	HIS	N-CA-C	6.96	121.08	112.59
1	A	299	ILE	N-CA-C	6.89	118.22	111.67
1	A	188	ASP	CA-CB-CG	6.81	119.41	112.60
1	B	657	PHE	CA-CB-CG	6.67	120.47	113.80
1	B	625	HIS	N-CA-C	6.63	121.02	112.26
1	B	438	SER	CA-C-N	6.62	129.87	120.53
1	B	438	SER	C-N-CA	6.62	129.87	120.53
1	B	181	SER	N-CA-C	6.49	119.18	108.99
1	A	219	ASP	CA-CB-CG	6.43	119.03	112.60
1	B	163	LEU	N-CA-C	6.35	119.01	111.33
1	B	341	LYS	CA-C-N	6.31	128.64	120.44
1	B	341	LYS	C-N-CA	6.31	128.64	120.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	80	VAL	N-CA-C	6.20	117.51	108.58
1	A	326	PHE	CA-CB-CG	6.07	119.87	113.80
1	B	521	VAL	N-CA-C	6.02	116.83	108.89
1	A	597	GLU	CA-C-N	5.94	126.57	119.98
1	A	597	GLU	C-N-CA	5.94	126.57	119.98
1	A	105	ASN	CA-CB-CG	5.93	118.53	112.60
1	B	213	GLY	N-CA-C	5.90	119.71	110.91
1	A	521	VAL	N-CA-C	5.89	117.06	108.58
1	B	494	THR	N-CA-C	5.80	119.33	112.72
1	B	34	TYR	N-CA-C	5.78	118.89	109.59
1	B	210	GLY	N-CA-C	5.71	126.70	113.18
1	A	296	ASP	N-CA-C	5.69	118.21	111.33
1	A	35	GLU	N-CA-C	-5.66	103.06	110.53
1	B	207	ASP	CA-CB-CG	5.62	118.22	112.60
1	A	456	SER	N-CA-C	-5.61	105.16	112.23
1	B	209	SER	N-CA-C	-5.60	98.87	110.80
1	A	51	ASP	CA-CB-CG	5.58	118.18	112.60
1	A	529	ASN	N-CA-C	-5.56	104.52	112.13
1	A	301	PRO	CA-C-N	5.52	127.51	120.56
1	A	301	PRO	C-N-CA	5.52	127.51	120.56
1	B	188	ASP	CA-CB-CG	5.49	118.09	112.60
1	A	156	ASN	CA-CB-CG	5.46	118.06	112.60
1	B	486	GLU	CA-C-N	5.46	128.85	121.05
1	B	486	GLU	C-N-CA	5.46	128.85	121.05
1	A	224	GLY	CA-C-N	5.39	128.52	120.87
1	A	224	GLY	C-N-CA	5.39	128.52	120.87
1	B	636	ASN	CA-C-N	5.37	127.74	120.38
1	B	636	ASN	C-N-CA	5.37	127.74	120.38
1	A	181	SER	N-CA-C	5.37	117.34	109.24
1	A	401	LYS	N-CA-C	5.36	116.81	110.97
1	B	651	ASN	CA-CB-CG	5.35	117.95	112.60
1	B	315	VAL	CB-CA-C	5.35	120.06	111.29
1	A	172	LEU	CA-C-N	5.22	127.80	120.28
1	A	172	LEU	C-N-CA	5.22	127.80	120.28
1	A	356	ASP	CA-CB-CG	5.21	117.81	112.60
1	A	487	ARG	N-CA-C	5.19	117.83	110.10
1	B	176	SER	CA-C-N	5.18	127.18	120.44
1	B	176	SER	C-N-CA	5.18	127.18	120.44
1	A	268	VAL	CA-C-N	5.12	125.67	119.98
1	A	268	VAL	C-N-CA	5.12	125.67	119.98
1	A	464	LYS	N-CA-C	-5.12	107.20	113.50
1	B	315	VAL	CA-C-N	5.11	128.85	120.63

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	315	VAL	C-N-CA	5.11	128.85	120.63
1	B	509	TRP	N-CA-C	5.10	117.58	111.71
1	B	479	ASP	CA-CB-CG	5.10	117.70	112.60
1	B	398	VAL	N-CA-CB	5.09	116.47	110.82
1	A	634	ASP	CA-CB-CG	5.08	117.68	112.60
1	B	649	GLU	N-CA-C	5.07	118.73	112.24
1	B	413	LEU	CA-C-N	5.06	125.46	119.94
1	B	413	LEU	C-N-CA	5.06	125.46	119.94
1	B	508	PRO	CA-C-N	5.05	127.55	120.28
1	B	508	PRO	C-N-CA	5.05	127.55	120.28
1	B	37	ASP	CA-CB-CG	5.02	117.62	112.60
1	B	619	TYR	CA-C-N	5.02	126.96	120.44
1	B	619	TYR	C-N-CA	5.02	126.96	120.44

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	5348	0	5000	69	0
1	B	5378	0	5024	66	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	129	0	258	14	0
3	B	84	0	168	6	0
4	B	12	0	16	1	0
5	A	473	0	0	4	0
5	B	408	0	0	0	0
All	All	11836	0	10466	136	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (136) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:443:VAL:HG23	1:A:584:GLN:HG3	1.37	1.01
1:A:279:MET:HE1	1:A:408:THR:HG23	1.40	0.99
1:B:318:MET:HE3	1:B:379:MET:H	1.41	0.86
1:A:318:MET:HE3	1:A:379:MET:H	1.39	0.85
1:A:443:VAL:CG2	1:A:584:GLN:HG3	2.06	0.85
1:A:589:VAL:HG11	1:A:632:ILE:HD12	1.63	0.81
1:B:9:ILE:HB	1:B:439:ILE:HD11	1.62	0.79
1:B:589:VAL:HG11	1:B:632:ILE:HD12	1.65	0.79
1:B:331:GLY:H	3:B:1696:EOH:H23	1.55	0.72
1:B:461:VAL:HG22	1:B:493:GLU:HG3	1.71	0.72
1:A:26:PRO:HD2	3:A:1720:EOH:H11	1.72	0.71
1:B:379:MET:HG3	3:B:1703:EOH:H12	1.73	0.70
1:B:221:ILE:HD13	3:B:1704:EOH:H22	1.76	0.68
1:B:255:ILE:HG22	1:B:583:PRO:HB3	1.77	0.66
1:A:318:MET:HG2	1:A:377:MET:SD	2.35	0.66
1:B:515:ASN:HB3	3:B:1713:EOH:H12	1.78	0.65
1:B:318:MET:HE3	1:B:379:MET:N	2.12	0.65
1:A:314:ASN:HB2	3:A:1691:EOH:H22	1.79	0.64
1:A:464:LYS:HG3	1:A:470:LYS:HG3	1.78	0.63
1:A:137:ILE:HG12	1:A:168:PRO:HD3	1.80	0.63
1:A:443:VAL:HG23	1:A:584:GLN:CG	2.23	0.61
1:A:577:TYR:CD2	3:A:1691:EOH:H12	2.37	0.59
1:A:318:MET:HE3	1:A:379:MET:N	2.15	0.59
1:B:112:ASN:HB2	1:B:354:ILE:HD11	1.85	0.58
1:B:461:VAL:CG2	1:B:493:GLU:HG3	2.32	0.58
1:B:209:SER:O	1:B:213:GLY:HA3	2.03	0.58
1:B:406:ASP:H	1:B:409:GLN:HE21	1.51	0.58
1:B:3:LEU:HD22	1:B:238:GLU:HG3	1.84	0.57
1:B:464:LYS:HE3	1:B:470:LYS:HG3	1.87	0.57
1:A:142:ARG:HB2	1:A:185:LEU:HD12	1.88	0.56
1:A:657:PHE:HE1	1:A:677:PRO:HB3	1.71	0.56
1:B:177:LYS:HD2	1:B:177:LYS:H	1.72	0.55
1:A:518:ILE:HD13	1:A:577:TYR:CE1	2.42	0.54
1:B:378:PRO:HB2	1:B:403:LEU:HD23	1.89	0.54
1:A:534:THR:HG23	1:A:566:VAL:HG22	1.90	0.54
1:A:579:TYR:OH	3:A:1691:EOH:H11	2.07	0.54
1:A:508:PRO:HB2	1:A:510:GLN:OE1	2.08	0.53
1:B:451:ILE:HD11	1:B:516:CYS:SG	2.49	0.53
1:A:577:TYR:O	3:A:1689:EOH:H12	2.08	0.53
1:B:301:PRO:O	1:B:305:ARG:HD2	2.08	0.53
1:A:341:LYS:NZ	5:A:2232:HOH:O	2.29	0.52
1:B:15:GLY:HA3	1:B:58:ARG:HG2	1.89	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:195:VAL:HG11	1:A:244:MET:HG2	1.91	0.52
1:A:293:ALA:O	1:A:421:PHE:HA	2.10	0.52
1:B:137:ILE:HD11	1:B:218:TYR:HB3	1.90	0.52
1:A:100:ARG:NH1	5:A:2074:HOH:O	2.43	0.52
1:A:516:CYS:SG	1:A:581:TRQ:HB3	2.50	0.52
1:A:462:ILE:HD11	1:A:489:GLU:HG2	1.92	0.51
1:A:577:TYR:HD2	3:A:1691:EOH:H12	1.76	0.51
1:B:287:TYR:HB3	1:B:411:PHE:CE1	2.45	0.51
1:B:589:VAL:HG12	1:B:590:LEU:N	2.25	0.51
1:B:318:MET:HG2	1:B:377:MET:SD	2.50	0.51
1:A:589:VAL:O	1:A:609:GLN:HA	2.10	0.51
1:A:93:LYS:HE2	1:A:505:MET:HB2	1.93	0.50
1:A:669:SER:HA	1:A:672:ASN:OD1	2.11	0.50
1:A:182:LEU:HD22	1:A:222:SER:HB3	1.92	0.50
1:B:9:ILE:HB	1:B:439:ILE:CD1	2.36	0.50
1:B:80:VAL:HG21	1:B:83:ILE:HG13	1.92	0.50
1:B:147:VAL:HG23	1:B:189:SER:HA	1.92	0.50
1:B:649:GLU:HG2	3:B:1690:EOH:H21	1.94	0.50
1:B:443:VAL:CG2	1:B:584:GLN:HG3	2.42	0.49
1:A:589:VAL:HG12	1:A:590:LEU:N	2.26	0.49
1:A:83:ILE:HG12	1:A:230:VAL:HG22	1.95	0.49
1:B:261:LEU:HD23	1:B:439:ILE:HG12	1.95	0.49
1:A:630:ALA:HB2	1:A:650:ARG:HB2	1.94	0.49
1:A:359:GLN:HB3	1:A:360:PRO:HD2	1.94	0.48
1:A:185:LEU:HD22	1:A:193:LEU:HD11	1.96	0.48
1:B:91:ASN:HB3	1:B:137:ILE:HB	1.96	0.48
1:A:365:MET:CG	1:A:373:ILE:HG23	2.44	0.47
1:B:141:PRO:HD3	1:B:162:TYR:CD1	2.48	0.47
1:A:88:HIS:O	1:A:224:GLY:HA3	2.15	0.47
1:A:132:ARG:HA	1:A:135:LEU:HD12	1.96	0.47
1:B:210:GLY:C	1:B:212:GLY:H	2.19	0.47
1:B:29:ILE:HB	3:B:1700:EOH:H12	1.96	0.47
1:A:187:THR:HA	1:A:192:ARG:O	2.13	0.47
1:B:215:ASP:O	1:B:217:TRP:HD1	1.98	0.47
1:B:448:CYS:HB3	1:B:580:TRP:CZ3	2.50	0.47
1:A:12:ALA:HB3	1:A:246:VAL:HG22	1.96	0.46
1:B:3:LEU:HD12	1:B:67:GLY:HA2	1.97	0.46
1:B:136:ILE:HD11	1:B:485:PRO:HB3	1.96	0.46
1:B:488:ASP:HB3	1:B:491:GLU:HB3	1.96	0.46
1:A:520:THR:HG22	1:A:575:ALA:HB1	1.97	0.46
1:A:308:GLN:HB2	5:A:2178:HOH:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:470:LYS:HB3	1:A:495:GLY:HA2	1.97	0.46
1:B:15:GLY:O	1:B:199:TYR:HA	2.16	0.46
1:A:509:TRP:HB2	3:A:1696:EOH:H23	1.96	0.46
1:A:85:TRP:CD2	1:A:193:LEU:HB2	2.51	0.45
1:B:99:PHE:CZ	1:B:101:GLU:HA	2.51	0.45
1:B:620:SER:HB2	4:B:1715:GOL:H12	1.99	0.45
1:B:137:ILE:HG12	1:B:168:PRO:HD3	1.99	0.45
1:B:454:THR:HG21	1:B:504:ARG:HB2	1.99	0.45
1:A:274:ASP:O	1:A:275:LEU:C	2.60	0.45
1:A:506:ALA:HB1	1:A:511:ALA:HB3	1.98	0.45
1:B:126:LYS:HD2	1:B:131:GLU:HB3	1.99	0.45
1:A:15:GLY:O	1:A:199:TYR:HA	2.17	0.45
1:B:589:VAL:CG1	1:B:590:LEU:N	2.81	0.45
1:B:63:PHE:HB2	1:B:193:LEU:HB3	2.00	0.44
1:A:365:MET:HG3	1:A:373:ILE:HG23	1.99	0.44
1:B:85:TRP:CZ3	1:B:228:CYS:HB2	2.53	0.44
1:A:188:ASP:OD2	1:A:192:ARG:NH1	2.50	0.44
1:A:438:SER:HA	3:A:1720:EOH:C2	2.47	0.44
1:A:516:CYS:O	1:A:585:SER:OG	2.33	0.44
1:A:506:ALA:CB	1:A:511:ALA:HB3	2.47	0.44
1:B:158:ILE:HG21	1:B:165:PRO:HG3	1.99	0.44
1:B:16:ASN:N	1:B:56:ILE:O	2.51	0.43
1:A:512:ASP:OD2	3:A:1696:EOH:H22	2.18	0.43
3:A:1705:EOH:H21	3:A:1725:EOH:H22	2.00	0.43
1:A:461:VAL:HG13	1:A:493:GLU:HA	2.01	0.43
1:B:108:TYR:CE2	1:B:398:VAL:HG11	2.53	0.43
1:A:438:SER:HA	3:A:1720:EOH:H23	2.01	0.42
1:B:186:ARG:O	1:B:194:ILE:N	2.50	0.42
1:A:579:TYR:CE2	3:A:1691:EOH:H11	2.53	0.42
1:B:279:MET:HE2	1:B:412:LEU:HD21	2.01	0.42
1:B:580:TRP:HB3	1:B:581:TRQ:CE3	2.50	0.42
1:B:98:GLU:HG2	1:B:122:ARG:HG2	2.00	0.42
1:B:443:VAL:HG23	1:B:584:GLN:HG3	2.01	0.42
1:A:63:PHE:HB2	1:A:193:LEU:HB3	2.02	0.42
1:B:74:THR:HB	1:B:190:GLN:HB3	2.01	0.42
1:A:93:LYS:HZ1	3:A:1696:EOH:HO	1.63	0.41
1:A:108:TYR:CE2	1:A:352:LYS:HG3	2.55	0.41
1:B:9:ILE:HD11	1:B:469:ILE:HG12	2.02	0.41
1:B:349:LEU:HD11	1:B:405:LEU:O	2.19	0.41
1:A:144:VAL:HG11	1:A:185:LEU:HB3	2.02	0.41
1:B:181:SER:OG	1:B:199:TYR:HE2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:13:ARG:HH12	1:B:452:GLU:HG2	1.86	0.41
1:B:153:ILE:O	1:B:181:SER:HB2	2.21	0.41
1:A:289:SER:HB3	5:A:2195:HOH:O	2.21	0.41
1:B:88:HIS:O	1:B:224:GLY:HA3	2.21	0.41
1:A:13:ARG:HA	1:A:247:GLY:O	2.20	0.40
1:A:87:VAL:HB	1:A:185:LEU:HD11	2.02	0.40
1:B:142:ARG:HH12	1:B:159:PRO:HD3	1.86	0.40
1:A:432:SER:HB2	3:A:1701:EOH:H11	2.04	0.40
1:A:29:ILE:HG12	1:A:440:GLY:HA3	2.03	0.40
1:A:385:SER:HB2	1:A:521:VAL:HG22	2.04	0.40
1:A:403:LEU:HD11	1:A:445:LEU:HB3	2.04	0.40
1:B:259:SER:OG	1:B:409:GLN:NE2	2.54	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	675/746 (90%)	624 (92%)	46 (7%)	5 (1%)	18	27
1	B	682/746 (91%)	629 (92%)	48 (7%)	5 (1%)	18	27
All	All	1357/1492 (91%)	1253 (92%)	94 (7%)	10 (1%)	18	27

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	68	ALA
1	A	455	TRP
1	A	156	ASN
1	B	212	GLY
1	B	476	ALA
1	B	568	GLU

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Mol	Chain	Res	Type
1	A	526	PRO
1	B	40	LEU
1	A	78	PRO
1	B	214	GLY

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	591/644 (92%)	552 (93%)	39 (7%)	15	25
1	B	593/644 (92%)	562 (95%)	31 (5%)	21	34
All	All	1184/1288 (92%)	1114 (94%)	70 (6%)	18	30

All (70) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	MET
1	A	38	VAL
1	A	44	THR
1	A	74	THR
1	A	79	ASN
1	A	125	SER
1	A	127	THR
1	A	147	VAL
1	A	160	GLU
1	A	163	LEU
1	A	190	GLN
1	A	194	ILE
1	A	215	ASP
1	A	220	ASP
1	A	225	SER
1	A	238	GLU
1	A	274	ASP
1	A	275	LEU
1	A	276	VAL

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Mol	Chain	Res	Type
1	A	353	VAL
1	A	364	LEU
1	A	365	MET
1	A	372	ASP
1	A	427	ASP
1	A	428	ASP
1	A	433	ASP
1	A	443	VAL
1	A	454	THR
1	A	461	VAL
1	A	462	ILE
1	A	464	LYS
1	A	520	THR
1	A	541	THR
1	A	604	LEU
1	A	640	ASP
1	A	653	GLU
1	A	659	GLU
1	A	670	GLU
1	A	676	LEU
1	B	27	MET
1	B	29	ILE
1	B	47	VAL
1	B	106	LEU
1	B	138	ASP
1	B	147	VAL
1	B	177	LYS
1	B	196	LEU
1	B	220	ASP
1	B	221	ILE
1	B	231	THR
1	B	255	ILE
1	B	259	SER
1	B	260	THR
1	B	275	LEU
1	B	278	ASP
1	B	290	ASP
1	B	298	ASP
1	B	299	ILE
1	B	341	LYS
1	B	431	VAL
1	B	461	VAL

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Mol	Chain	Res	Type
1	B	464	LYS
1	B	521	VAL
1	B	563	THR
1	B	564	LYS
1	B	568	GLU
1	B	604	LEU
1	B	609	GLN
1	B	610	ILE
1	B	650	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	HIS
1	A	21	ASN
1	A	123	ASN
1	A	205	ASN
1	A	310	GLN
1	A	542	HIS
1	A	637	GLN
1	A	638	ASN
1	A	639	ASN
1	A	668	ASN
1	B	48	ASN
1	B	59	GLN
1	B	61	GLN
1	B	314	ASN
1	B	409	GLN
1	B	609	GLN
1	B	651	ASN
1	B	652	HIS
1	B	668	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
1	TRQ	A	581	1	16,17,18	3.22	8 (50%)	16,24,26	1.07	1 (6%)
1	TRQ	B	581	1	16,17,18	3.28	8 (50%)	16,24,26	1.03	1 (6%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	TRQ	A	581	1	-	0/5/19/21	0/2/2/2
1	TRQ	B	581	1	-	0/5/19/21	0/2/2/2

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	581	TRQ	O7-CZ2	8.85	1.43	1.23
1	B	581	TRQ	O7-CZ2	8.71	1.42	1.23
1	B	581	TRQ	O6-CH2	5.67	1.39	1.24
1	A	581	TRQ	O6-CH2	5.61	1.39	1.24
1	B	581	TRQ	CE2-NE1	-4.65	1.31	1.38
1	A	581	TRQ	CE2-NE1	-3.65	1.32	1.38
1	B	581	TRQ	CE2-CZ2	-3.34	1.38	1.50
1	A	581	TRQ	CE2-CZ2	-3.25	1.38	1.50
1	A	581	TRQ	CH2-CZ2	-3.15	1.38	1.53
1	B	581	TRQ	CH2-CZ2	-2.97	1.39	1.53
1	A	581	TRQ	CZ3-CH2	-2.56	1.39	1.45
1	B	581	TRQ	CE3-CZ3	2.52	1.41	1.35
1	A	581	TRQ	CE3-CZ3	2.42	1.41	1.35
1	B	581	TRQ	CB-CA	-2.17	1.49	1.53
1	A	581	TRQ	CB-CA	-2.14	1.49	1.53
1	B	581	TRQ	CZ3-CH2	-2.11	1.40	1.45

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	581	TRQ	O7-CZ2-CH2	2.33	121.36	118.39
1	A	581	TRQ	CB-CG-CD1	-2.02	123.03	126.97

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	581	TRQ	1	0
1	B	581	TRQ	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 77 ligands modelled in this entry, 4 are monoatomic - leaving 73 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	EOH	A	1705	-	2,2,2	0.55	0	1,1,1	0.73	0
3	EOH	B	1709	-	2,2,2	0.59	0	1,1,1	0.69	0
3	EOH	A	1719	-	2,2,2	0.49	0	1,1,1	0.47	0
3	EOH	B	1696	-	2,2,2	0.51	0	1,1,1	0.64	0
3	EOH	B	1702	-	2,2,2	0.57	0	1,1,1	0.63	0
3	EOH	A	1714	-	2,2,2	0.54	0	1,1,1	0.59	0
3	EOH	B	1801	-	2,2,2	0.53	0	1,1,1	0.69	0
3	EOH	A	1726	-	2,2,2	0.59	0	1,1,1	0.59	0
3	EOH	B	1700	-	2,2,2	0.50	0	1,1,1	0.78	0
3	EOH	A	1694	-	2,2,2	0.58	0	1,1,1	0.69	0
3	EOH	B	1708	-	2,2,2	0.59	0	1,1,1	0.57	0
3	EOH	B	1800	-	2,2,2	0.47	0	1,1,1	0.66	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	EOH	A	1707	-	2,2,2	0.45	0	1,1,1	0.67	0
4	GOL	B	1714	-	5,5,5	0.08	0	5,5,5	0.12	0
3	EOH	A	1689	-	2,2,2	0.49	0	1,1,1	0.70	0
3	EOH	A	1712	-	2,2,2	0.40	0	1,1,1	0.54	0
3	EOH	A	1698	-	2,2,2	0.54	0	1,1,1	0.62	0
3	EOH	B	1706	-	2,2,2	0.54	0	1,1,1	0.84	0
3	EOH	A	1692	-	2,2,2	0.59	0	1,1,1	0.61	0
3	EOH	A	1803	-	2,2,2	0.61	0	1,1,1	0.62	0
3	EOH	A	1801	-	2,2,2	0.54	0	1,1,1	0.73	0
3	EOH	A	1703	-	2,2,2	0.50	0	1,1,1	0.87	0
3	EOH	B	1695	-	2,2,2	0.53	0	1,1,1	0.67	0
3	EOH	B	1711	-	2,2,2	0.58	0	1,1,1	0.45	0
3	EOH	A	1723	-	2,2,2	0.50	0	1,1,1	0.70	0
3	EOH	A	1717	-	2,2,2	0.53	0	1,1,1	0.73	0
3	EOH	B	1689	-	2,2,2	0.52	0	1,1,1	0.66	0
3	EOH	A	1701	-	2,2,2	0.53	0	1,1,1	0.54	0
3	EOH	B	1710	-	2,2,2	0.64	0	1,1,1	0.69	0
3	EOH	A	1800	-	2,2,2	0.55	0	1,1,1	0.67	0
3	EOH	A	1697	-	2,2,2	0.57	0	1,1,1	0.66	0
3	EOH	A	1724	-	2,2,2	0.55	0	1,1,1	0.63	0
3	EOH	B	1705	-	2,2,2	0.54	0	1,1,1	0.74	0
3	EOH	B	1707	-	2,2,2	0.53	0	1,1,1	0.69	0
3	EOH	A	1725	-	2,2,2	0.55	0	1,1,1	0.64	0
3	EOH	A	1704	-	2,2,2	0.46	0	1,1,1	0.68	0
3	EOH	B	1712	-	2,2,2	0.58	0	1,1,1	0.78	0
3	EOH	B	1693	-	2,2,2	0.59	0	1,1,1	0.55	0
3	EOH	B	1697	-	2,2,2	0.59	0	1,1,1	0.61	0
3	EOH	A	1696	-	2,2,2	0.63	0	1,1,1	0.54	0
4	GOL	B	1715	-	5,5,5	0.06	0	5,5,5	0.22	0
3	EOH	B	1699	-	2,2,2	0.58	0	1,1,1	0.62	0
3	EOH	B	1694	-	2,2,2	0.59	0	1,1,1	0.81	0
3	EOH	A	1721	-	2,2,2	0.50	0	1,1,1	0.75	0
3	EOH	B	1691	-	2,2,2	0.51	0	1,1,1	0.54	0
3	EOH	B	1704	-	2,2,2	0.62	0	1,1,1	0.66	0
3	EOH	B	1802	-	2,2,2	0.56	0	1,1,1	0.64	0
3	EOH	A	1700	-	2,2,2	0.64	0	1,1,1	0.75	0
3	EOH	A	1722	-	2,2,2	0.52	0	1,1,1	0.58	0
3	EOH	A	1727	-	2,2,2	0.54	0	1,1,1	0.61	0
3	EOH	A	1695	-	2,2,2	0.45	0	1,1,1	0.76	0
3	EOH	A	1690	-	2,2,2	0.55	0	1,1,1	0.62	0
3	EOH	A	1709	-	2,2,2	0.66	0	1,1,1	0.59	0
3	EOH	A	1711	-	2,2,2	0.59	0	1,1,1	0.61	0
3	EOH	A	1718	-	2,2,2	0.60	0	1,1,1	0.65	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	EOH	B	1692	-	2,2,2	0.53	0	1,1,1	0.67	0
3	EOH	A	1710	-	2,2,2	0.50	0	1,1,1	0.73	0
3	EOH	B	1713	-	2,2,2	0.57	0	1,1,1	0.58	0
3	EOH	A	1720	-	2,2,2	0.60	0	1,1,1	0.50	0
3	EOH	B	1703	-	2,2,2	0.68	0	1,1,1	0.18	0
3	EOH	A	1713	-	2,2,2	0.51	0	1,1,1	0.62	0
3	EOH	A	1708	-	2,2,2	0.57	0	1,1,1	0.58	0
3	EOH	B	1698	-	2,2,2	0.63	0	1,1,1	0.54	0
3	EOH	B	1701	-	2,2,2	0.51	0	1,1,1	0.82	0
3	EOH	A	1699	-	2,2,2	0.55	0	1,1,1	0.75	0
3	EOH	A	1702	-	2,2,2	0.59	0	1,1,1	0.65	0
3	EOH	A	1693	-	2,2,2	0.55	0	1,1,1	0.70	0
3	EOH	B	1690	-	2,2,2	0.54	0	1,1,1	0.71	0
3	EOH	A	1715	-	2,2,2	0.60	0	1,1,1	0.65	0
3	EOH	A	1716	-	2,2,2	0.47	0	1,1,1	0.39	0
3	EOH	A	1706	-	2,2,2	0.52	0	1,1,1	0.64	0
3	EOH	A	1691	-	2,2,2	0.83	0	1,1,1	0.14	0
3	EOH	A	1802	-	2,2,2	0.54	0	1,1,1	0.70	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	B	1714	-	-	1/4/4/4	-
4	GOL	B	1715	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	1715	GOL	C1-C2-C3-O3
4	B	1715	GOL	O2-C2-C3-O3
4	B	1714	GOL	O1-C1-C2-C3

There are no ring outliers.

14 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1705	EOH	1	0
3	B	1696	EOH	1	0
3	B	1700	EOH	1	0
3	A	1689	EOH	1	0
3	A	1701	EOH	1	0
3	A	1725	EOH	1	0
3	A	1696	EOH	3	0
4	B	1715	GOL	1	0
3	B	1704	EOH	1	0
3	B	1713	EOH	1	0
3	A	1720	EOH	3	0
3	B	1703	EOH	1	0
3	B	1690	EOH	1	0
3	A	1691	EOH	5	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	679/746 (91%)	-1.36	0 100 100	15, 37, 61, 76	0
1	B	684/746 (91%)	-1.27	0 100 100	26, 44, 70, 86	0
All	All	1363/1492 (91%)	-1.31	0 100 100	15, 41, 65, 86	0

There are no RSRZ outliers to report.

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

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
1	TRQ	A	581	16/17	0.99	0.03	17,27,30,31	0
1	TRQ	B	581	16/17	0.99	0.02	22,33,39,43	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	EOH	B	1698	3/3	0.92	0.09	43,43,44,45	0
3	EOH	B	1695	3/3	0.93	0.18	76,76,77,78	0
3	EOH	A	1723	3/3	0.93	0.10	51,51,52,52	0
3	EOH	B	1697	3/3	0.94	0.08	37,37,40,42	0
3	EOH	A	1802	3/3	0.94	0.10	52,52,53,53	0
3	EOH	B	1699	3/3	0.94	0.09	58,58,59,60	0
3	EOH	A	1726	3/3	0.95	0.09	48,48,50,50	0
3	EOH	A	1699	3/3	0.95	0.08	51,51,52,52	0
3	EOH	A	1706	3/3	0.95	0.07	56,56,57,57	0
3	EOH	A	1722	3/3	0.95	0.08	48,48,48,49	0
3	EOH	A	1693	3/3	0.95	0.11	80,80,80,80	0
3	EOH	A	1724	3/3	0.95	0.08	57,57,57,58	0
3	EOH	B	1706	3/3	0.95	0.10	45,45,46,47	0
3	EOH	B	1712	3/3	0.95	0.06	35,35,36,37	0
3	EOH	A	1708	3/3	0.96	0.07	43,43,44,46	0
3	EOH	A	1710	3/3	0.96	0.08	45,45,46,47	0
3	EOH	A	1690	3/3	0.96	0.08	45,45,47,47	0
3	EOH	A	1698	3/3	0.96	0.07	34,34,38,42	0
3	EOH	A	1691	3/3	0.96	0.08	27,27,28,29	0
3	EOH	A	1692	3/3	0.96	0.08	53,53,56,57	0
3	EOH	B	1707	3/3	0.96	0.07	49,49,50,50	0
3	EOH	A	1801	3/3	0.96	0.11	68,68,69,69	0
3	EOH	B	1801	3/3	0.96	0.10	79,79,80,80	0
3	EOH	A	1709	3/3	0.97	0.05	31,31,32,33	0
3	EOH	A	1800	3/3	0.97	0.06	44,44,45,45	0
3	EOH	A	1705	3/3	0.97	0.10	65,65,66,66	0
3	EOH	A	1713	3/3	0.97	0.06	35,35,38,42	0
3	EOH	A	1803	3/3	0.97	0.06	28,28,32,35	0
3	EOH	B	1689	3/3	0.97	0.09	45,45,46,48	0
3	EOH	B	1690	3/3	0.97	0.08	48,48,48,49	0
3	EOH	A	1714	3/3	0.97	0.07	41,41,43,44	0
3	EOH	B	1696	3/3	0.97	0.07	32,32,34,35	0
3	EOH	A	1717	3/3	0.97	0.14	67,67,67,67	0
3	EOH	A	1719	3/3	0.97	0.10	39,39,42,43	0
3	EOH	A	1721	3/3	0.97	0.08	60,60,62,64	0
3	EOH	B	1701	3/3	0.97	0.13	28,28,32,36	0
3	EOH	B	1704	3/3	0.97	0.08	47,47,49,49	0
2	NA	A	1688	1/1	0.97	0.06	27,27,27,27	0
3	EOH	A	1707	3/3	0.97	0.09	41,41,46,48	0
3	EOH	B	1709	3/3	0.97	0.10	54,54,55,55	0
3	EOH	A	1702	3/3	0.97	0.08	29,29,31,36	0
3	EOH	B	1800	3/3	0.97	0.05	37,37,40,44	0
3	EOH	A	1725	3/3	0.97	0.07	45,45,45,46	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	EOH	A	1718	3/3	0.98	0.05	38,38,41,42	0
3	EOH	A	1727	3/3	0.98	0.06	50,50,51,52	0
3	EOH	A	1694	3/3	0.98	0.06	40,40,41,43	0
3	EOH	A	1720	3/3	0.98	0.12	34,34,37,38	0
3	EOH	A	1711	3/3	0.98	0.05	31,31,34,34	0
3	EOH	B	1702	3/3	0.98	0.06	31,31,31,32	0
3	EOH	B	1703	3/3	0.98	0.06	21,21,26,27	0
3	EOH	A	1696	3/3	0.98	0.08	33,33,35,36	0
3	EOH	B	1705	3/3	0.98	0.09	41,41,41,41	0
3	EOH	A	1703	3/3	0.98	0.06	47,47,48,50	0
3	EOH	A	1715	3/3	0.98	0.07	35,35,37,38	0
3	EOH	B	1691	3/3	0.98	0.05	26,26,28,29	0
3	EOH	B	1710	3/3	0.98	0.04	32,32,32,33	0
3	EOH	B	1711	3/3	0.98	0.05	37,37,40,40	0
3	EOH	B	1692	3/3	0.98	0.05	36,36,38,39	0
3	EOH	B	1694	3/3	0.98	0.10	32,32,34,34	0
3	EOH	A	1689	3/3	0.98	0.05	28,28,32,34	0
3	EOH	B	1802	3/3	0.98	0.10	51,51,51,52	0
4	GOL	B	1714	6/6	0.98	0.06	68,71,71,73	0
4	GOL	B	1715	6/6	0.98	0.04	38,40,46,48	0
3	EOH	A	1697	3/3	0.99	0.05	40,40,43,43	0
3	EOH	B	1708	3/3	0.99	0.07	43,43,43,44	0
3	EOH	A	1716	3/3	0.99	0.04	29,29,34,37	0
3	EOH	A	1695	3/3	0.99	0.04	14,14,15,16	0
3	EOH	B	1700	3/3	0.99	0.03	26,26,27,27	0
3	EOH	A	1704	3/3	0.99	0.04	38,38,39,41	0
3	EOH	B	1713	3/3	0.99	0.07	36,36,38,38	0
2	NA	B	1688	1/1	0.99	0.04	30,30,30,30	0
3	EOH	B	1693	3/3	0.99	0.04	29,29,30,31	0
3	EOH	A	1712	3/3	0.99	0.05	32,32,35,38	0
3	EOH	A	1700	3/3	0.99	0.04	24,24,25,26	0
3	EOH	A	1701	3/3	0.99	0.04	36,36,36,37	0
2	NA	A	1687	1/1	1.00	0.03	26,26,26,26	0
2	NA	B	1687	1/1	1.00	0.01	27,27,27,27	0

6.5 Other polymers ⓘ

There are no such residues in this entry.