



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 09:55 PM UTC

PDB ID : 8ZMU / pdb_00008zmu
Title : GLUTAMATE DEHYDROGENASE (W89F-MUTANT) FROM THERMO-
COCCUS PROFUNDUS IN THE UNLIGANDED STATE
Authors : Wakabayashi, T.; Matsui, Y.; Masayoshi, M.
Deposited on : 2024-05-23
Resolution : 2.03 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

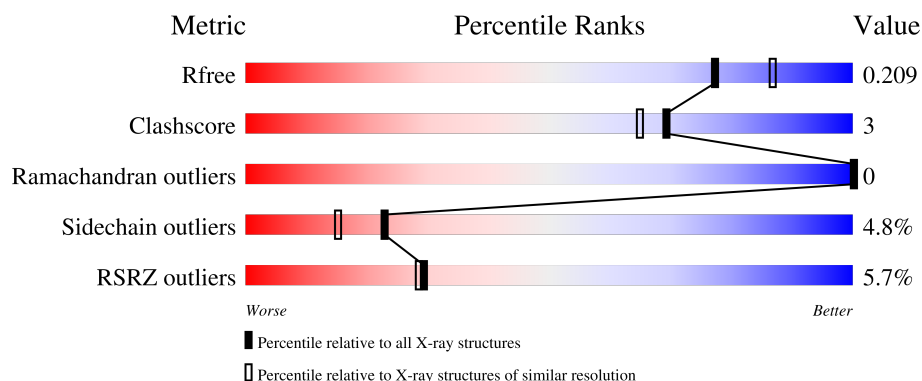
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.03 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	13299 (2.04-2.00)
Clashscore	190562	1022 (2.02-2.02)
Ramachandran outliers	187476	1014 (2.02-2.02)
Sidechain outliers	187428	1014 (2.02-2.02)
RSRZ outliers	180081	13314 (2.04-2.00)

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	419	2% 89% 9% ..
1	B	419	% 89% 10% ..
1	C	419	2% 88% 11% .
1	D	419	4% 87% 11% ..
1	E	419	2% 89% 9% ..

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Mol	Chain	Length	Quality of chain
1	F	419	

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	ACY	A	512	-	-	X	-
4	GOL	C	513	-	-	X	-
4	GOL	C	517	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 20653 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 Glutamate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	415	Total	C	N	O	S	0	1	0
			3262	2082	558	608	14			
1	B	416	Total	C	N	O	S	0	0	0
			3259	2080	556	609	14			
1	C	415	Total	C	N	O	S	0	0	0
			3254	2077	555	608	14			
1	D	415	Total	C	N	O	S	0	0	0
			3243	2068	554	607	14			
1	E	416	Total	C	N	O	S	0	1	0
			3273	2089	560	610	14			
1	F	398	Total	C	N	O	S	0	0	0
			3106	1982	527	583	14			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	89	PHE	TRP	engineered mutation	UNP O74024
B	89	PHE	TRP	engineered mutation	UNP O74024
C	89	PHE	TRP	engineered mutation	UNP O74024
D	89	PHE	TRP	engineered mutation	UNP O74024
E	89	PHE	TRP	engineered mutation	UNP O74024
F	89	PHE	TRP	engineered mutation	UNP O74024

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



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

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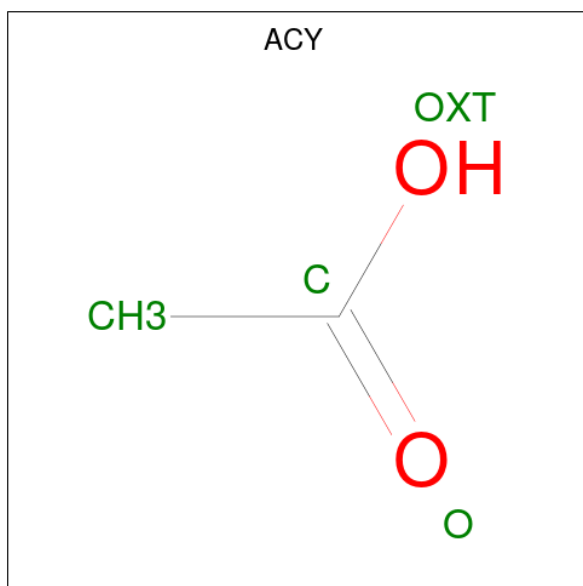
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total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

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	E	1	Total O S 5 4 1	0	0
2	E	1	Total O S 5 4 1	0	0
2	E	1	Total O S 5 4 1	0	0
2	E	1	Total O S 5 4 1	0	0
2	F	1	Total O S 5 4 1	0	0
2	F	1	Total O S 5 4 1	0	0
2	F	1	Total O S 5 4 1	0	0
2	F	1	Total O S 5 4 1	0	0

- Molecule 3 is ACETIC ACID (CCD ID: ACY) (formula: C₂H₄O₂).



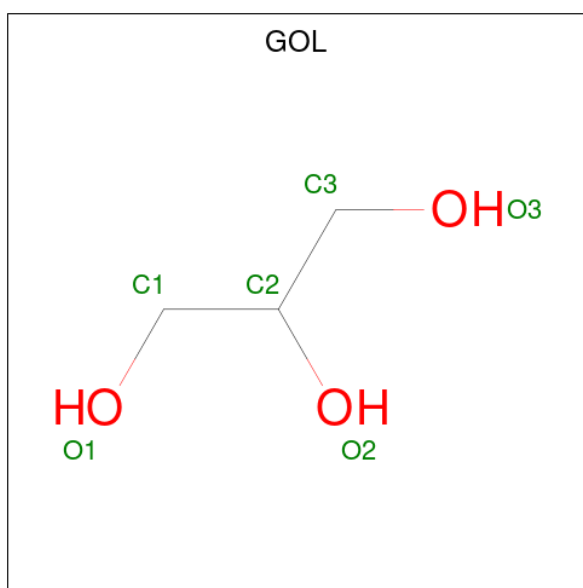
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0

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

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



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

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

- Molecule 5 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	Na	0	0
			1	1		

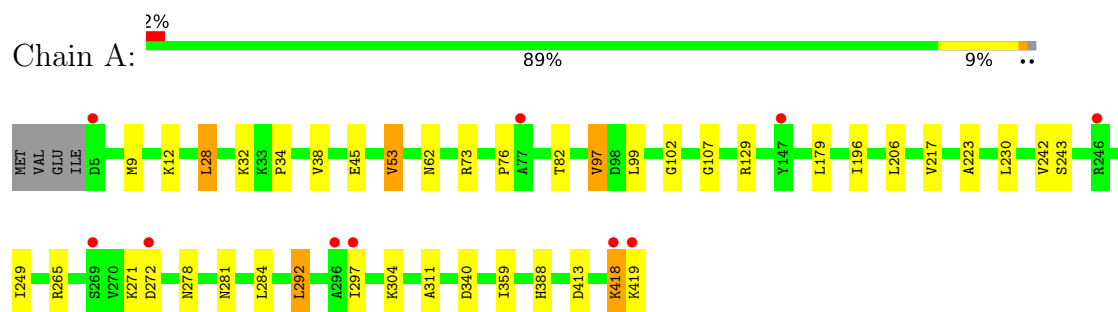
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	190	Total	O	0	0
			190	190		
6	B	184	Total	O	0	0
			184	184		
6	C	172	Total	O	0	0
			172	172		
6	D	93	Total	O	0	0
			93	93		
6	E	221	Total	O	0	0
			221	221		
6	F	88	Total	O	0	0
			88	88		

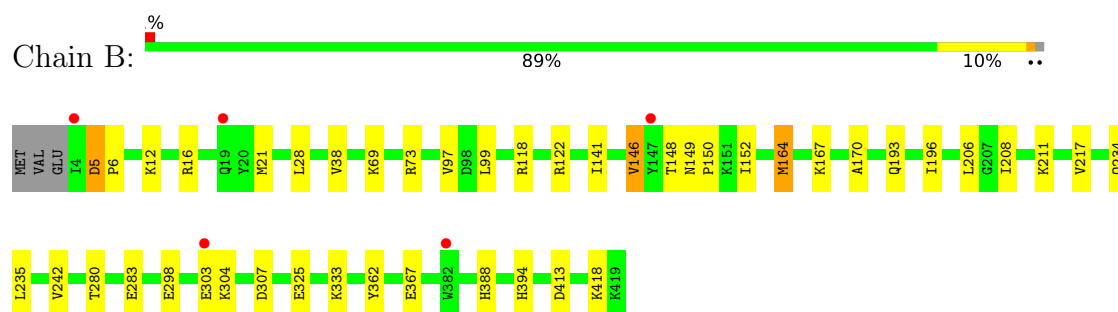
3 Residue-property plots [i](#)

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.

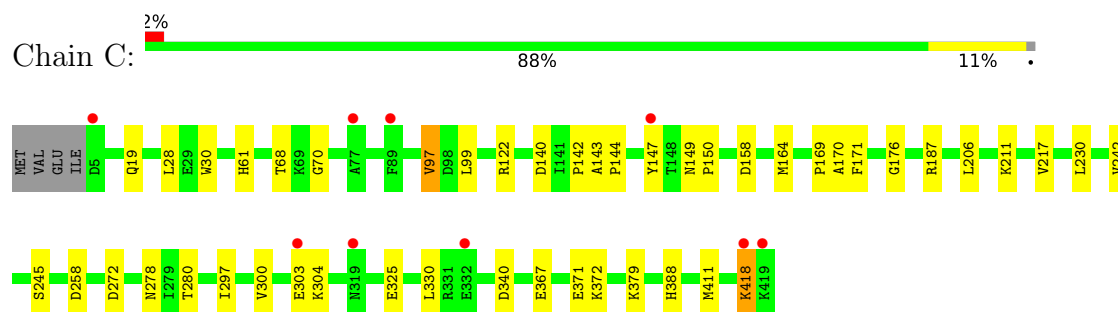
- Molecule 1: Glutamate dehydrogenase



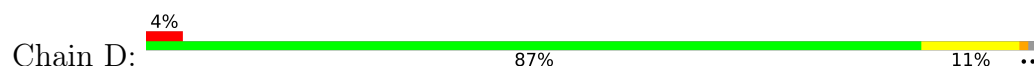
- Molecule 1: Glutamate dehydrogenase

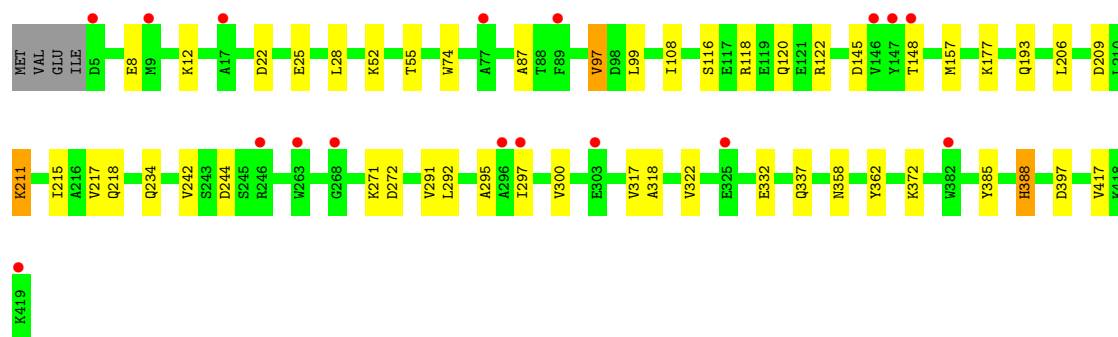


- Molecule 1: Glutamate dehydrogenase

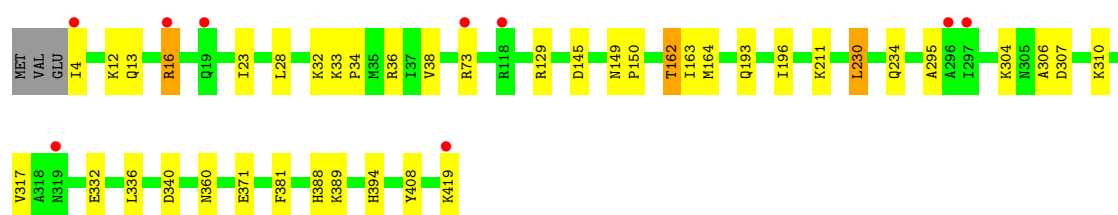
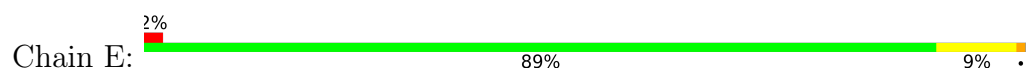


- Molecule 1: Glutamate dehydrogenase

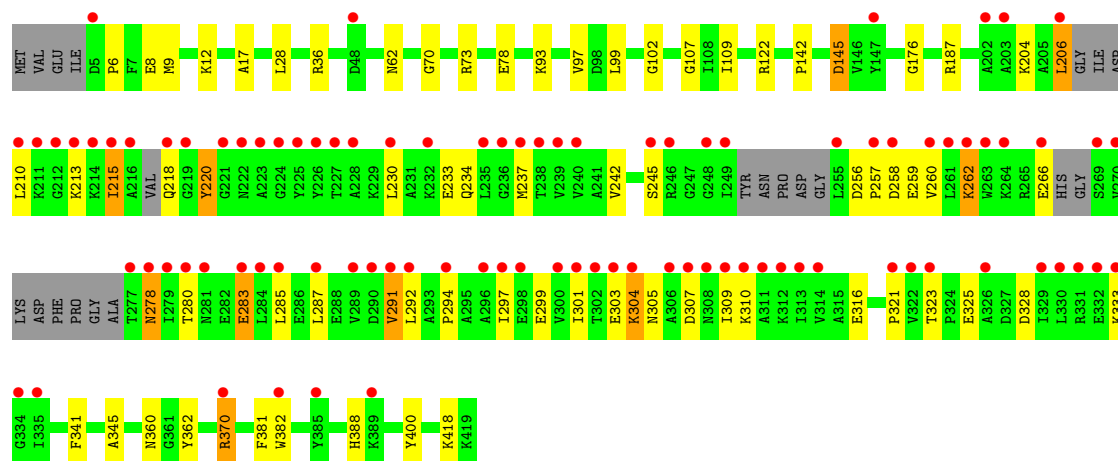
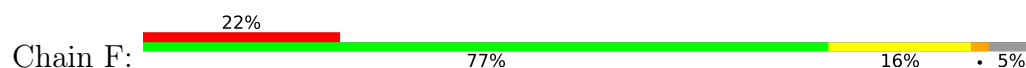




● Molecule 1: Glutamate dehydrogenase



● Molecule 1: Glutamate dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	113.36Å 163.25Å 132.60Å 90.00° 113.91° 90.00°	Depositor
Resolution (Å)	21.80 – 2.03 21.80 – 2.03	Depositor EDS
% Data completeness (in resolution range)	99.6 (21.80-2.03) 99.6 (21.80-2.03)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.10 (at 2.02Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.172 , 0.199 0.184 , 0.209	Depositor DCC
R_{free} test set	14014 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	26.9	Xtriage
Anisotropy	0.500	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 38.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.013 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	20653	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.24% 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: NA, GOL, ACY, SO4

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.15	3/3337 (0.1%)	1.30	6/4519 (0.1%)
1	B	1.16	5/3331 (0.2%)	1.33	3/4512 (0.1%)
1	C	1.20	4/3326 (0.1%)	1.37	6/4505 (0.1%)
1	D	1.17	3/3314 (0.1%)	1.42	5/4490 (0.1%)
1	E	1.13	4/3345 (0.1%)	1.31	3/4530 (0.1%)
1	F	1.13	1/3166 (0.0%)	1.37	1/4282 (0.0%)
All	All	1.16	20/19819 (0.1%)	1.35	24/26838 (0.1%)

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	306	ALA	C-O	6.63	1.31	1.24
1	E	394	HIS	CE1-NE2	6.49	1.39	1.32
1	A	249	ILE	C-O	6.48	1.30	1.24
1	C	330	LEU	C-O	6.40	1.31	1.24
1	C	158	ASP	C-O	6.16	1.31	1.24
1	C	297	ILE	C-O	5.89	1.30	1.23
1	D	417	VAL	C-O	5.83	1.29	1.24
1	A	34	PRO	C-O	-5.69	1.16	1.23
1	D	362	TYR	C-O	5.64	1.30	1.23
1	E	34	PRO	C-O	-5.54	1.17	1.23
1	B	146	VAL	C-O	5.35	1.29	1.24
1	B	152	ILE	C-O	-5.32	1.17	1.24
1	C	61	HIS	CE1-NE2	5.23	1.37	1.32
1	B	413	ASP	C-O	5.19	1.30	1.24
1	E	13	GLN	C-O	5.18	1.30	1.24
1	A	223	ALA	C-O	5.13	1.29	1.24
1	B	148	THR	C-O	5.12	1.30	1.23
1	D	87	ALA	C-O	5.10	1.29	1.24
1	F	145	ASP	C-O	-5.05	1.17	1.24
1	B	394	HIS	CE1-NE2	5.03	1.37	1.32

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	304	LYS	N-CA-C	8.30	120.33	111.28
1	C	258	ASP	CA-CB-CG	-6.70	105.90	112.60
1	F	418	LYS	N-CA-C	-6.28	105.44	113.23
1	A	53	VAL	CB-CA-C	6.04	119.58	111.19
1	E	162	THR	CA-CB-OG1	-5.90	100.75	109.60
1	E	381	PHE	CA-CB-CG	-5.87	107.93	113.80
1	B	418	LYS	N-CA-C	-5.85	106.14	113.28
1	D	55	THR	CA-CB-OG1	-5.76	100.96	109.60
1	C	280	THR	CA-CB-OG1	-5.76	100.97	109.60
1	C	418	LYS	N-CA-C	-5.67	105.86	112.89
1	A	82	THR	CA-CB-OG1	-5.61	101.19	109.60
1	B	5	ASP	CA-CB-CG	-5.60	107.00	112.60
1	D	397	ASP	N-CA-C	-5.57	105.21	112.23
1	C	340	ASP	CA-CB-CG	5.41	118.01	112.60
1	D	22	ASP	CA-CB-CG	5.34	117.94	112.60
1	A	340	ASP	CA-CB-CG	5.28	117.88	112.60
1	A	76	PRO	CB-CA-C	5.27	117.88	110.98
1	D	388	HIS	CA-C-O	-5.25	114.86	120.42
1	D	385	TYR	CA-C-O	-5.22	115.52	121.00
1	E	340	ASP	CA-CB-CG	5.18	117.78	112.60
1	C	300	VAL	N-CA-C	-5.15	105.82	113.39
1	A	413	ASP	CA-CB-CG	-5.14	107.46	112.60
1	A	281	ASN	CB-CA-C	5.04	118.87	110.81
1	C	147	TYR	CA-C-O	-5.00	114.41	120.32

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	3262	0	3271	12	0
1	B	3259	0	3260	22	0
1	C	3254	0	3258	27	0
1	D	3243	0	3240	19	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	3273	0	3281	14	0
1	F	3106	0	3100	47	0
2	A	50	0	0	0	0
2	B	35	0	0	0	0
2	C	45	0	0	0	0
2	D	30	0	0	1	0
2	E	35	0	0	0	0
2	F	20	0	0	0	0
3	A	8	0	6	2	0
3	B	8	0	6	1	0
3	C	12	0	9	0	0
3	E	4	0	3	0	0
4	A	6	0	8	1	0
4	B	12	0	16	2	0
4	C	30	0	40	14	0
4	E	12	0	16	0	0
5	B	1	0	0	0	0
6	A	190	0	0	1	0
6	B	184	0	0	0	0
6	C	172	0	0	1	0
6	D	93	0	0	0	0
6	E	221	0	0	2	0
6	F	88	0	0	1	0
All	All	20653	0	19514	137	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 3.

All (137) 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:F:262:LYS:HE3	1:F:262:LYS:HA	1.48	0.95
1:F:218:GLN:HB3	1:F:294:PRO:HA	1.52	0.92
3:A:512:ACY:H1	1:F:36:ARG:HD2	1.52	0.92
1:B:193:GLN:HE22	1:B:234:GLN:HE22	1.19	0.91
1:B:303:GLU:HG3	1:B:325:GLU:HB2	1.54	0.90
1:E:193:GLN:HE22	1:E:234:GLN:HE22	1.25	0.84
4:B:510:GOL:H32	1:D:122:ARG:HE	1.45	0.82
1:F:262:LYS:O	1:F:266:GLU:HG2	1.81	0.81
1:C:272:ASP:H	1:C:278:ASN:HD21	1.28	0.80
1:F:262:LYS:HE3	1:F:262:LYS:CA	2.15	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:C:513:GOL:H12	1:F:122:ARG:HE	1.51	0.75
1:C:169:PRO:HA	4:C:517:GOL:H11	1.69	0.74
1:F:303:GLU:HG3	1:F:325:GLU:HB2	1.68	0.73
1:B:303:GLU:H	1:B:303:GLU:CD	1.97	0.73
1:C:122:ARG:HH21	4:C:513:GOL:H11	1.56	0.71
1:B:303:GLU:HG3	1:B:325:GLU:CB	2.25	0.67
1:C:171:PHE:CD1	4:C:517:GOL:H32	2.29	0.67
1:C:171:PHE:HB2	4:C:517:GOL:H12	1.77	0.66
1:D:193:GLN:HE22	1:D:234:GLN:HE22	1.43	0.65
1:C:245:SER:HA	4:C:514:GOL:H32	1.78	0.65
1:F:215:ILE:HG21	1:F:237:MET:HE3	1.79	0.65
1:A:243:SER:HB3	1:A:284:LEU:HD13	1.78	0.65
1:F:256:ASP:O	1:F:260:VAL:HG23	1.98	0.63
1:F:258:ASP:O	1:F:262:LYS:HG2	1.98	0.63
1:C:171:PHE:H	4:C:517:GOL:H12	1.64	0.62
1:F:301:ILE:HA	1:F:305:ASN:HD21	1.64	0.62
1:F:230:LEU:O	1:F:234:GLN:HB2	2.00	0.61
1:F:262:LYS:HA	1:F:262:LYS:CE	2.27	0.60
1:F:6:PRO:HA	1:F:9:MET:HE3	1.83	0.60
1:F:301:ILE:HD13	1:F:309:ILE:HG12	1.83	0.60
1:F:292:LEU:HG	1:F:294:PRO:HD3	1.82	0.60
1:D:74:TRP:NE1	1:D:108:ILE:HD12	2.18	0.58
1:F:93:LYS:NZ	6:F:604:HOH:O	2.36	0.58
1:D:218:GLN:HE22	1:D:300:VAL:HG21	1.70	0.57
4:C:513:GOL:H12	1:F:122:ARG:NE	2.20	0.56
1:F:230:LEU:HD22	1:F:234:GLN:OE1	2.05	0.56
1:C:149:ASN:HB2	1:C:150:PRO:CD	2.36	0.56
1:E:196:ILE:HG13	1:E:230:LEU:HB3	1.87	0.56
1:F:70:GLY:O	1:F:142:PRO:HA	2.07	0.55
1:D:177:LYS:HZ1	1:D:358:ASN:HD21	1.55	0.55
1:F:220:TYR:OH	1:F:257:PRO:HB3	2.07	0.55
1:C:176:GLY:HA2	1:C:187:ARG:HD3	1.89	0.54
1:B:362:TYR:OH	1:C:372:LYS:HD3	2.08	0.54
1:B:164:MET:HE3	1:B:167:LYS:H	1.73	0.53
1:C:325:GLU:H	1:C:325:GLU:CD	2.16	0.53
1:E:360:ASN:HA	1:F:360:ASN:HD21	1.74	0.53
1:C:171:PHE:HB2	4:C:517:GOL:C1	2.38	0.53
1:F:278:ASN:HD22	1:F:278:ASN:N	2.06	0.52
1:D:177:LYS:NZ	1:D:358:ASN:HD21	2.07	0.52
1:D:244:ASP:HB2	2:D:501:SO4:O3	2.09	0.52
1:B:118:ARG:O	1:B:122:ARG:HG3	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:285:LEU:HB3	1:F:309:ILE:HD11	1.92	0.51
1:E:16:ARG:HG2	6:E:703:HOH:O	2.10	0.51
1:F:259:GLU:HA	1:F:262:LYS:HG3	1.93	0.51
1:A:217:VAL:O	1:A:242:VAL:HA	2.11	0.51
1:C:149:ASN:HB2	1:C:150:PRO:HD2	1.91	0.50
1:A:271:LYS:HG2	1:A:278:ASN:ND2	2.27	0.49
1:B:164:MET:HE1	1:B:170:ALA:HB2	1.95	0.49
1:D:209:ASP:OD1	1:D:211:LYS:HB2	2.12	0.49
1:F:215:ILE:HB	1:F:291:VAL:HG13	1.93	0.49
3:A:512:ACY:CH3	1:F:36:ARG:HD2	2.34	0.49
1:E:295:ALA:HB2	1:E:317:VAL:HB	1.95	0.49
1:C:171:PHE:H	4:C:517:GOL:H31	1.77	0.48
1:E:193:GLN:HE22	1:E:234:GLN:NE2	2.04	0.48
1:E:33:LYS:HE3	6:E:612:HOH:O	2.13	0.48
1:C:68:THR:O	1:C:140:ASP:HA	2.14	0.48
1:B:164:MET:CE	1:B:170:ALA:HB2	2.44	0.47
1:D:295:ALA:HB2	1:D:317:VAL:HB	1.96	0.47
1:F:17:ALA:HB1	1:F:400:TYR:HB3	1.97	0.47
1:B:73:ARG:HG3	1:B:146:VAL:HB	1.97	0.47
1:F:97:VAL:HG23	1:F:99:LEU:HG	1.96	0.47
1:A:97:VAL:HG23	1:A:99:LEU:HG	1.97	0.47
1:B:307:ASP:HA	1:B:333:LYS:CE	2.45	0.46
1:B:97:VAL:HG23	1:B:99:LEU:HG	1.97	0.46
1:C:122:ARG:HE	4:C:513:GOL:H11	1.80	0.46
1:F:256:ASP:HA	1:F:257:PRO:HD3	1.78	0.46
1:A:9:MET:HE3	1:A:9:MET:HB2	1.82	0.46
1:A:272:ASP:H	1:A:278:ASN:HD21	1.62	0.46
1:B:38:VAL:HG22	1:E:38:VAL:HG22	1.98	0.46
1:A:28:LEU:O	1:A:32:LYS:HG3	2.16	0.45
1:F:206:LEU:HD12	1:F:206:LEU:HA	1.77	0.45
1:F:316:GLU:HB3	1:F:321:PRO:HG2	1.98	0.45
1:B:280:THR:OG1	1:B:283:GLU:HG3	2.16	0.45
1:C:379:LYS:HE3	6:C:622:HOH:O	2.16	0.45
1:C:169:PRO:CA	4:C:517:GOL:H11	2.43	0.45
1:C:170:ALA:N	4:C:517:GOL:H31	2.31	0.45
1:A:129[A]:ARG:CZ	1:E:162:THR:HG21	2.46	0.45
1:B:217:VAL:O	1:B:242:VAL:HA	2.17	0.45
1:F:370:ARG:HE	1:F:370:ARG:HB3	1.48	0.45
1:E:149:ASN:HB2	1:E:150:PRO:CD	2.47	0.45
1:F:78:GLU:HB3	1:F:109:ILE:HG12	1.99	0.44
1:C:97:VAL:HG23	1:C:99:LEU:HG	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:157:MET:C	1:D:157:MET:SD	3.00	0.44
1:D:322:VAL:HG11	1:D:337:GLN:NE2	2.32	0.44
1:F:299:GLU:HG3	1:F:323:THR:HG22	2.00	0.44
1:C:217:VAL:O	1:C:242:VAL:HA	2.17	0.44
1:B:5:ASP:HB3	1:B:6:PRO:HD3	2.00	0.44
1:D:372:LYS:HD2	1:F:362:TYR:OH	2.17	0.44
1:F:341:PHE:O	1:F:345:ALA:HB2	2.18	0.43
1:D:217:VAL:O	1:D:242:VAL:HA	2.19	0.43
1:E:23:ILE:HA	1:E:408:TYR:CE2	2.53	0.43
1:F:176:GLY:HA2	1:F:187:ARG:HD3	2.01	0.43
1:C:143:ALA:HB1	1:C:144:PRO:HD2	2.01	0.43
1:D:297:ILE:HG13	1:D:300:VAL:CG1	2.49	0.43
1:F:303:GLU:CD	1:F:303:GLU:H	2.25	0.43
1:D:295:ALA:HA	1:D:318:ALA:HB2	2.00	0.43
4:A:513:GOL:H31	6:A:767:HOH:O	2.19	0.43
1:F:73:ARG:O	1:F:107:GLY:HA2	2.18	0.42
1:D:271:LYS:HG2	1:D:272:ASP:OD2	2.19	0.42
1:C:171:PHE:H	4:C:517:GOL:C3	2.32	0.42
1:F:304:LYS:HE3	1:F:304:LYS:HB3	1.44	0.42
1:F:381:PHE:CD1	1:F:381:PHE:C	2.98	0.42
3:B:508:ACY:H1	1:E:36:ARG:HD3	2.01	0.42
1:C:30:TRP:CD1	1:C:411:MET:HE3	2.54	0.42
1:B:69:LYS:HD2	1:B:141:ILE:HB	2.02	0.42
1:C:303:GLU:OE1	1:C:304:LYS:HE2	2.20	0.42
1:C:272:ASP:H	1:C:278:ASN:ND2	2.06	0.41
1:B:307:ASP:HA	1:B:333:LYS:HE3	2.02	0.41
1:E:129[A]:ARG:HG2	1:E:163:ILE:HD11	2.02	0.41
1:F:73:ARG:HG2	1:F:145:ASP:OD1	2.20	0.41
1:A:292:LEU:HB2	1:A:311:ALA:HB2	2.03	0.41
1:C:70:GLY:O	1:C:142:PRO:HA	2.21	0.41
1:A:62:ASN:O	1:A:102:GLY:HA3	2.21	0.41
1:B:208:ILE:HA	4:B:511:GOL:H12	2.03	0.41
1:F:62:ASN:O	1:F:102:GLY:HA3	2.21	0.41
1:F:218:GLN:NE2	1:F:285:LEU:HD21	2.36	0.41
1:F:280:THR:HG22	1:F:283:GLU:OE1	2.20	0.41
1:A:418:LYS:H	1:A:418:LYS:HG2	1.73	0.41
1:D:215:ILE:HA	1:D:291:VAL:O	2.20	0.41
1:D:116:SER:O	1:D:120:GLN:HG3	2.21	0.41
1:F:213:LYS:HD3	1:F:213:LYS:HA	1.93	0.41
1:D:97:VAL:HG23	1:D:99:LEU:HG	2.03	0.41
1:A:73:ARG:O	1:A:107:GLY:HA2	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:28:LEU:HD13	1:E:32:LYS:HD3	2.03	0.40
1:B:193:GLN:HE22	1:B:234:GLN:NE2	2.01	0.40
1:B:149:ASN:HB2	1:B:150:PRO:HD2	2.02	0.40
1:B:164:MET:HE3	1:B:167:LYS:N	2.35	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	414/419 (99%)	399 (96%)	15 (4%)	0	100	100
1	B	414/419 (99%)	404 (98%)	10 (2%)	0	100	100
1	C	413/419 (99%)	402 (97%)	11 (3%)	0	100	100
1	D	413/419 (99%)	399 (97%)	14 (3%)	0	100	100
1	E	415/419 (99%)	403 (97%)	12 (3%)	0	100	100
1	F	386/419 (92%)	358 (93%)	28 (7%)	0	100	100
All	All	2455/2514 (98%)	2365 (96%)	90 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	339/342 (99%)	321 (95%)	18 (5%)	20	13
1	B	338/342 (99%)	326 (96%)	12 (4%)	31	26
1	C	338/342 (99%)	327 (97%)	11 (3%)	33	29
1	D	336/342 (98%)	322 (96%)	14 (4%)	26	20
1	E	340/342 (99%)	323 (95%)	17 (5%)	22	15
1	F	321/342 (94%)	296 (92%)	25 (8%)	11	6
All	All	2012/2052 (98%)	1915 (95%)	97 (5%)	23	15

All (97) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	LYS
1	A	28	LEU
1	A	38	VAL
1	A	45	GLU
1	A	53	VAL
1	A	97	VAL
1	A	179	LEU
1	A	196	ILE
1	A	206	LEU
1	A	230	LEU
1	A	265	ARG
1	A	292	LEU
1	A	297	ILE
1	A	304	LYS
1	A	359	ILE
1	A	388	HIS
1	A	418	LYS
1	A	419	LYS
1	B	12	LYS
1	B	16	ARG
1	B	21	MET
1	B	28	LEU
1	B	164	MET
1	B	196	ILE
1	B	206	LEU
1	B	211	LYS
1	B	235	LEU
1	B	298	GLU
1	B	367	GLU
1	B	388	HIS

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Mol	Chain	Res	Type
1	C	19	GLN
1	C	28	LEU
1	C	97	VAL
1	C	164	MET
1	C	206	LEU
1	C	211	LYS
1	C	230	LEU
1	C	367	GLU
1	C	371	GLU
1	C	388	HIS
1	C	418	LYS
1	D	8	GLU
1	D	12	LYS
1	D	25	GLU
1	D	28	LEU
1	D	52	LYS
1	D	97	VAL
1	D	118	ARG
1	D	145	ASP
1	D	148	THR
1	D	206	LEU
1	D	211	LYS
1	D	292	LEU
1	D	332	GLU
1	D	388	HIS
1	E	4	ILE
1	E	12	LYS
1	E	16	ARG
1	E	73	ARG
1	E	145	ASP
1	E	164	MET
1	E	211	LYS
1	E	230	LEU
1	E	304	LYS
1	E	307	ASP
1	E	310	LYS
1	E	332	GLU
1	E	336	LEU
1	E	371	GLU
1	E	388	HIS
1	E	389	LYS
1	E	419	LYS

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Mol	Chain	Res	Type
1	F	8	GLU
1	F	12	LYS
1	F	28	LEU
1	F	204	LYS
1	F	206	LEU
1	F	210	LEU
1	F	215	ILE
1	F	220	TYR
1	F	233	GLU
1	F	242	VAL
1	F	245	SER
1	F	262	LYS
1	F	278	ASN
1	F	283	GLU
1	F	287	LEU
1	F	291	VAL
1	F	297	ILE
1	F	304	LYS
1	F	307	ASP
1	F	310	LYS
1	F	328	ASP
1	F	333	LYS
1	F	370	ARG
1	F	382	TRP
1	F	388	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	281	ASN
1	A	360	ASN
1	A	386	ASN
1	B	13	GLN
1	B	120	GLN
1	B	234	GLN
1	B	308	ASN
1	B	386	ASN
1	B	388	HIS
1	C	267	HIS
1	C	278	ASN
1	C	308	ASN
1	C	386	ASN

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Mol	Chain	Res	Type
1	D	218	GLN
1	D	234	GLN
1	D	281	ASN
1	D	308	ASN
1	D	358	ASN
1	D	386	ASN
1	E	120	GLN
1	E	193	GLN
1	E	308	ASN
1	E	386	ASN
1	F	13	GLN
1	F	120	GLN
1	F	278	ASN
1	F	308	ASN
1	F	360	ASN

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](#)

Of 62 ligands modelled in this entry, 1 is monoatomic - leaving 61 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	ACY	C	510	-	3,3,3	1.48	0	3,3,3	0.60	0
2	SO4	A	501	-	4,4,4	0.22	0	6,6,6	0.55	0
2	SO4	E	505	-	4,4,4	0.35	0	6,6,6	0.27	0
2	SO4	C	506	-	4,4,4	0.42	0	6,6,6	0.08	0
2	SO4	B	507	-	4,4,4	0.42	0	6,6,6	0.06	0
4	GOL	C	513	-	5,5,5	0.26	0	5,5,5	0.88	0
3	ACY	A	512	-	3,3,3	1.11	0	3,3,3	1.23	0
3	ACY	A	511	-	3,3,3	1.13	0	3,3,3	1.00	0
2	SO4	B	503	-	4,4,4	0.23	0	6,6,6	0.28	0
2	SO4	A	509	-	4,4,4	0.33	0	6,6,6	0.21	0
2	SO4	A	505	-	4,4,4	0.16	0	6,6,6	0.43	0
4	GOL	B	511	-	5,5,5	0.25	0	5,5,5	0.63	0
2	SO4	B	501	-	4,4,4	0.23	0	6,6,6	0.23	0
2	SO4	E	503	-	4,4,4	0.42	0	6,6,6	0.06	0
2	SO4	F	503	-	4,4,4	0.29	0	6,6,6	0.15	0
3	ACY	C	512	-	3,3,3	1.29	0	3,3,3	1.12	0
4	GOL	B	510	-	5,5,5	0.08	0	5,5,5	0.56	0
2	SO4	C	503	-	4,4,4	0.39	0	6,6,6	0.40	0
2	SO4	F	502	-	4,4,4	0.30	0	6,6,6	0.15	0
2	SO4	F	501	-	4,4,4	0.37	0	6,6,6	0.17	0
3	ACY	E	508	-	3,3,3	1.41	0	3,3,3	0.75	0
4	GOL	C	515	-	5,5,5	0.12	0	5,5,5	0.41	0
2	SO4	A	506	-	4,4,4	0.38	0	6,6,6	0.22	0
2	SO4	C	501	-	4,4,4	0.35	0	6,6,6	0.44	0
3	ACY	B	509	-	3,3,3	1.25	0	3,3,3	0.70	0
2	SO4	F	504	-	4,4,4	0.36	0	6,6,6	0.20	0
2	SO4	C	504	-	4,4,4	0.20	0	6,6,6	0.14	0
2	SO4	E	507	-	4,4,4	0.37	0	6,6,6	0.41	0
2	SO4	A	503	-	4,4,4	0.29	0	6,6,6	0.22	0
4	GOL	C	516	-	5,5,5	0.20	0	5,5,5	0.47	0
2	SO4	A	507	-	4,4,4	0.29	0	6,6,6	0.27	0
2	SO4	D	502	-	4,4,4	0.30	0	6,6,6	0.20	0
2	SO4	D	505	-	4,4,4	0.41	0	6,6,6	0.08	0
2	SO4	E	501	-	4,4,4	0.35	0	6,6,6	0.22	0
2	SO4	D	501	-	4,4,4	0.37	0	6,6,6	0.26	0
4	GOL	E	510	-	5,5,5	0.19	0	5,5,5	0.44	0
2	SO4	B	505	-	4,4,4	0.37	0	6,6,6	0.17	0
2	SO4	C	508	-	4,4,4	0.21	0	6,6,6	0.31	0
2	SO4	E	504	-	4,4,4	0.32	0	6,6,6	0.15	0
3	ACY	C	511	-	3,3,3	0.84	0	3,3,3	1.12	0
2	SO4	E	506	-	4,4,4	0.35	0	6,6,6	0.27	0
2	SO4	B	502	-	4,4,4	0.47	0	6,6,6	0.32	0
2	SO4	C	505	-	4,4,4	0.35	0	6,6,6	0.15	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	ACY	B	508	-	3,3,3	0.83	0	3,3,3	1.60	1 (33%)
4	GOL	C	514	-	5,5,5	0.16	0	5,5,5	0.57	0
4	GOL	E	509	-	5,5,5	0.18	0	5,5,5	0.41	0
2	SO4	D	504	-	4,4,4	0.31	0	6,6,6	0.21	0
4	GOL	A	513	-	5,5,5	0.16	0	5,5,5	0.42	0
2	SO4	B	504	5	4,4,4	0.25	0	6,6,6	0.14	0
2	SO4	C	509	-	4,4,4	0.26	0	6,6,6	0.25	0
2	SO4	D	506	-	4,4,4	0.33	0	6,6,6	0.21	0
2	SO4	A	502	-	4,4,4	0.34	0	6,6,6	0.06	0
2	SO4	C	507	-	4,4,4	0.30	0	6,6,6	0.20	0
4	GOL	C	517	-	5,5,5	0.32	0	5,5,5	1.06	0
2	SO4	D	503	-	4,4,4	0.27	0	6,6,6	0.36	0
2	SO4	A	510	-	4,4,4	0.26	0	6,6,6	0.19	0
2	SO4	A	504	-	4,4,4	0.13	0	6,6,6	0.20	0
2	SO4	B	506	-	4,4,4	0.26	0	6,6,6	0.34	0
2	SO4	A	508	-	4,4,4	0.32	0	6,6,6	0.22	0
2	SO4	E	502	-	4,4,4	0.52	0	6,6,6	0.17	0
2	SO4	C	502	-	4,4,4	0.51	0	6,6,6	0.31	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	A	513	-	-	2/4/4/4	-
4	GOL	B	510	-	-	2/4/4/4	-
4	GOL	E	510	-	-	2/4/4/4	-
4	GOL	C	513	-	-	2/4/4/4	-
4	GOL	C	515	-	-	3/4/4/4	-
4	GOL	C	517	-	-	2/4/4/4	-
4	GOL	E	509	-	-	0/4/4/4	-
4	GOL	C	514	-	-	2/4/4/4	-
4	GOL	B	511	-	-	2/4/4/4	-
4	GOL	C	516	-	-	0/4/4/4	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	508	ACY	O-C-CH3	-2.21	113.46	122.53

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	511	GOL	O1-C1-C2-C3
4	C	515	GOL	O1-C1-C2-C3
4	C	517	GOL	O1-C1-C2-C3
4	A	513	GOL	C1-C2-C3-O3
4	B	510	GOL	O1-C1-C2-C3
4	C	513	GOL	C1-C2-C3-O3
4	C	514	GOL	O1-C1-C2-C3
4	E	510	GOL	O1-C1-C2-C3
4	B	511	GOL	O1-C1-C2-O2
4	C	515	GOL	O1-C1-C2-O2
4	C	517	GOL	O1-C1-C2-O2
4	E	510	GOL	O1-C1-C2-O2
4	A	513	GOL	O2-C2-C3-O3
4	C	514	GOL	O1-C1-C2-O2
4	C	513	GOL	O2-C2-C3-O3
4	B	510	GOL	O1-C1-C2-O2
4	C	515	GOL	C1-C2-C3-O3

There are no ring outliers.

9 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	513	GOL	4	0
3	A	512	ACY	2	0
4	B	511	GOL	1	0
4	B	510	GOL	1	0
2	D	501	SO4	1	0
3	B	508	ACY	1	0
4	C	514	GOL	1	0
4	A	513	GOL	1	0
4	C	517	GOL	9	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	415/419 (99%)	-0.24	10 (2%) 59 60	17, 30, 54, 94	1 (0%)
1	B	416/419 (99%)	-0.22	5 (1%) 76 76	21, 32, 54, 72	0
1	C	415/419 (99%)	-0.05	9 (2%) 62 62	24, 33, 57, 80	0
1	D	415/419 (99%)	0.42	17 (4%) 41 41	28, 43, 70, 93	0
1	E	416/419 (99%)	-0.30	9 (2%) 62 62	14, 31, 53, 78	1 (0%)
1	F	398/419 (94%)	0.97	92 (23%) 2 2	24, 48, 117, 138	0
All	All	2475/2514 (98%)	0.09	142 (5%) 29 28	14, 35, 76, 138	2 (0%)

All (142) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	263	TRP	9.7
1	F	216	ALA	9.2
1	F	294	PRO	6.6
1	F	210	LEU	6.4
1	F	261	LEU	6.4
1	F	270	VAL	6.3
1	D	148	THR	5.9
1	E	4	ILE	5.7
1	F	255	LEU	5.5
1	F	203	ALA	5.4
1	D	325	GLU	5.4
1	F	211	LYS	5.3
1	F	382	TRP	5.3
1	F	249	ILE	5.2
1	F	260	VAL	5.1
1	E	297	ILE	5.0
1	F	239	VAL	5.0
1	F	285	LEU	5.0
1	F	297	ILE	4.9

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Mol	Chain	Res	Type	RSRZ
1	F	218	GLN	4.8
1	F	277	THR	4.8
1	F	225	TYR	4.8
1	E	319	ASN	4.7
1	F	279	ILE	4.7
1	F	309	ILE	4.7
1	F	333	LYS	4.6
1	F	292	LEU	4.6
1	F	330	LEU	4.5
1	F	335	ILE	4.4
1	B	303	GLU	4.3
1	F	212	GLY	4.2
1	F	248	GLY	4.2
1	F	312	LYS	4.2
1	C	147	TYR	4.2
1	A	419	LYS	4.1
1	A	77	ALA	4.0
1	F	287	LEU	4.0
1	F	284	LEU	4.0
1	F	236	GLY	3.9
1	F	308	ASN	3.9
1	A	418	LYS	3.9
1	E	19	GLN	3.9
1	F	280	THR	3.8
1	C	89	PHE	3.8
1	F	306	ALA	3.8
1	F	235	LEU	3.8
1	F	313	ILE	3.8
1	F	257	PRO	3.8
1	F	240	VAL	3.7
1	A	269	SER	3.7
1	A	147	TYR	3.7
1	D	147	TYR	3.7
1	F	48	ASP	3.6
1	F	329	ILE	3.6
1	F	334	GLY	3.6
1	F	245	SER	3.5
1	F	269	SER	3.5
1	F	301	ILE	3.4
1	F	278	ASN	3.4
1	F	215	ILE	3.4
1	F	300	VAL	3.3

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Mol	Chain	Res	Type	RSRZ
1	F	310	LYS	3.3
1	F	289	VAL	3.3
1	F	246	ARG	3.3
1	F	262	LYS	3.2
1	F	291	VAL	3.1
1	F	221	GLY	3.1
1	F	281	ASN	3.1
1	F	322	VAL	3.1
1	F	389	LYS	3.0
1	F	206	LEU	3.0
1	F	307	ASP	3.0
1	F	5	ASP	3.0
1	F	298	GLU	2.9
1	D	17	ALA	2.9
1	F	238	THR	2.9
1	F	266	GLU	2.9
1	F	326	ALA	2.9
1	B	147	TYR	2.9
1	C	5	ASP	2.8
1	F	290	ASP	2.8
1	F	323	THR	2.8
1	F	385	TYR	2.8
1	F	331	ARG	2.8
1	D	77	ALA	2.8
1	B	19	GLN	2.8
1	D	246	ARG	2.8
1	C	77	ALA	2.8
1	D	89	PHE	2.8
1	F	214	LYS	2.8
1	F	227	THR	2.8
1	F	304	LYS	2.8
1	F	321	PRO	2.8
1	F	226	TYR	2.7
1	B	4	ILE	2.7
1	F	224	GLY	2.7
1	F	237	MET	2.6
1	F	213	LYS	2.6
1	D	5	ASP	2.6
1	F	230	LEU	2.6
1	D	268	GLY	2.5
1	E	419	LYS	2.5
1	D	419	LYS	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	5	ASP	2.5
1	F	228	ALA	2.5
1	F	332	GLU	2.5
1	F	302	THR	2.5
1	D	303	GLU	2.4
1	F	219	GLY	2.4
1	D	297	ILE	2.4
1	F	296	ALA	2.4
1	F	311	ALA	2.4
1	D	146	VAL	2.4
1	F	303	GLU	2.3
1	F	314	VAL	2.3
1	D	263	TRP	2.3
1	A	296	ALA	2.3
1	E	296	ALA	2.3
1	D	382	TRP	2.3
1	E	73	ARG	2.2
1	E	118	ARG	2.2
1	C	419	LYS	2.2
1	F	370	ARG	2.2
1	C	319	ASN	2.2
1	D	296	ALA	2.2
1	F	264	LYS	2.2
1	A	272	ASP	2.2
1	E	16	ARG	2.2
1	B	382	TRP	2.2
1	F	258	ASP	2.2
1	C	303	GLU	2.1
1	F	147	TYR	2.1
1	F	202	ALA	2.1
1	F	232	LYS	2.1
1	D	9	MET	2.1
1	A	297	ILE	2.1
1	C	418	LYS	2.1
1	A	246	ARG	2.0
1	C	332	GLU	2.0
1	F	283	GLU	2.0
1	F	223	ALA	2.0
1	F	222	ASN	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
2	SO4	D	505	5/5	0.59	0.17	72,74,81,99	0
2	SO4	C	509	5/5	0.66	0.15	92,97,100,102	0
3	ACY	C	512	4/4	0.68	0.23	57,59,61,65	0
3	ACY	C	511	4/4	0.73	0.17	49,56,57,59	0
4	GOL	B	511	6/6	0.74	0.20	62,81,83,90	0
2	SO4	E	503	5/5	0.77	0.14	63,79,82,86	0
2	SO4	C	506	5/5	0.77	0.15	69,73,84,97	0
4	GOL	C	515	6/6	0.77	0.20	54,63,69,78	0
2	SO4	D	502	5/5	0.80	0.11	86,90,97,105	0
2	SO4	A	501	5/5	0.81	0.14	62,67,77,90	0
2	SO4	B	503	5/5	0.82	0.11	58,75,76,93	0
2	SO4	C	505	5/5	0.82	0.10	82,88,96,103	0
2	SO4	A	504	5/5	0.82	0.14	53,71,76,87	0
2	SO4	C	507	5/5	0.82	0.12	71,88,100,120	0
2	SO4	E	507	5/5	0.83	0.11	63,65,66,69	0
2	SO4	A	505	5/5	0.84	0.14	45,58,96,96	0
4	GOL	E	509	6/6	0.84	0.16	58,73,84,86	0
2	SO4	F	502	5/5	0.85	0.09	71,92,101,106	0
2	SO4	B	507	5/5	0.85	0.13	78,89,97,97	0
4	GOL	B	510	6/6	0.86	0.15	54,54,56,68	0
2	SO4	D	506	5/5	0.86	0.11	69,82,104,109	0
4	GOL	C	513	6/6	0.86	0.17	50,56,58,75	0
2	SO4	B	504	5/5	0.86	0.11	77,82,85,95	0
2	SO4	A	507	5/5	0.86	0.10	66,69,84,90	0
2	SO4	E	504	5/5	0.87	0.09	68,71,77,98	0
2	SO4	D	503	5/5	0.87	0.11	54,55,77,81	0
2	SO4	D	504	5/5	0.88	0.10	75,76,98,102	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	F	504	5/5	0.88	0.10	63,78,83,93	0
2	SO4	B	506	5/5	0.88	0.08	54,69,76,77	0
2	SO4	D	501	5/5	0.88	0.07	66,74,82,95	0
4	GOL	A	513	6/6	0.88	0.15	41,45,67,67	0
4	GOL	C	517	6/6	0.89	0.18	45,52,54,61	0
2	SO4	A	510	5/5	0.89	0.10	44,69,86,110	0
4	GOL	C	514	6/6	0.90	0.17	51,70,83,83	0
2	SO4	E	505	5/5	0.90	0.09	70,78,89,93	0
3	ACY	A	511	4/4	0.90	0.17	49,57,63,69	0
3	ACY	B	509	4/4	0.90	0.18	52,63,66,70	0
4	GOL	E	510	6/6	0.90	0.12	54,61,67,70	0
4	GOL	C	516	6/6	0.91	0.12	42,63,74,78	0
2	SO4	E	506	5/5	0.91	0.09	50,68,76,78	0
2	SO4	A	503	5/5	0.91	0.08	60,66,76,90	0
2	SO4	F	501	5/5	0.91	0.08	68,68,77,91	0
2	SO4	C	501	5/5	0.92	0.14	49,57,65,69	0
2	SO4	A	509	5/5	0.92	0.08	68,73,82,85	0
2	SO4	A	508	5/5	0.92	0.10	63,78,79,91	0
2	SO4	F	503	5/5	0.93	0.08	54,68,87,97	0
2	SO4	A	506	5/5	0.93	0.14	53,54,59,62	0
2	SO4	C	504	5/5	0.94	0.09	57,59,71,74	0
3	ACY	A	512	4/4	0.94	0.15	35,36,38,42	0
2	SO4	B	505	5/5	0.94	0.07	65,66,71,77	0
2	SO4	B	501	5/5	0.95	0.11	47,52,63,71	0
3	ACY	E	508	4/4	0.95	0.11	34,35,36,38	0
3	ACY	C	510	4/4	0.95	0.10	42,42,44,46	0
2	SO4	B	502	5/5	0.95	0.08	41,49,56,56	0
3	ACY	B	508	4/4	0.96	0.10	33,35,37,40	0
2	SO4	E	502	5/5	0.96	0.08	48,51,59,66	0
2	SO4	C	508	5/5	0.96	0.08	43,45,55,58	0
2	SO4	E	501	5/5	0.96	0.14	42,55,66,80	0
2	SO4	C	503	5/5	0.97	0.11	46,52,58,69	0
2	SO4	A	502	5/5	0.97	0.13	42,54,55,55	0
5	NA	B	512	1/1	0.97	0.06	39,39,39,39	0
2	SO4	C	502	5/5	0.98	0.08	42,44,52,52	0

6.5 Other polymers

There are no such residues in this entry.