Full wwPDB X-ray Structure Validation Report (i)

Sep 2, 2023 – 10:55 PM EDT

PDB ID	:	3RHM
Title	:	Crystal structure of the E673Q mutant oF C-Terminal domain of 10'FORMY LTETRAHYDROFOLATE DEHYDROGENASE
Authors	:	Tsybovsky, Y.
Deposited on	:	2011-04-11
Resolution	:	2.38 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 (i) 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 (i)) were used in the production of this report:

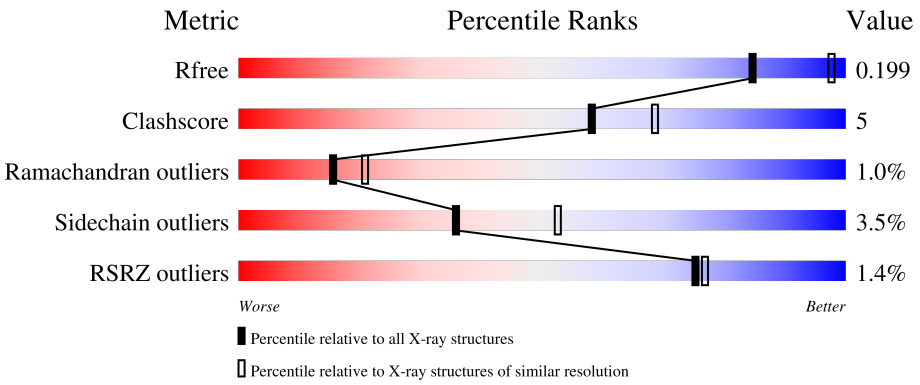
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.38 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}	130704	5509 (2.40-2.36)
Clashscore	141614	6082 (2.40-2.36)
Ramachandran outliers	138981	5973 (2.40-2.36)
Sidechain outliers	138945	5975 (2.40-2.36)
RSRZ outliers	127900	5397 (2.40-2.36)

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 <=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	517	.% 79% 16% • •
1	B	517	2% 79% 16% • •
1	C	517	.% 78% 17% • •
1	D	517	.% 80% 14% •• •

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 16121 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 Aldehyde dehydrogenase 1 family, member L1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	498	3826	2436	657	715	18	3	6	0
1	B	498	3817	2429	657	713	18	3	5	0
1	C	498	3806	2424	652	712	18	0	3	0
1	D	498	3825	2435	656	716	18	0	6	0

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	386	MET	-	expression tag	UNP Q5HZB2
A	387	ARG	-	expression tag	UNP Q5HZB2
A	388	GLY	-	expression tag	UNP Q5HZB2
A	389	SER	-	expression tag	UNP Q5HZB2
A	390	HIS	-	expression tag	UNP Q5HZB2
A	391	HIS	-	expression tag	UNP Q5HZB2
A	392	HIS	-	expression tag	UNP Q5HZB2
A	393	HIS	-	expression tag	UNP Q5HZB2
A	394	HIS	-	expression tag	UNP Q5HZB2
A	395	THR	-	expression tag	UNP Q5HZB2
A	396	THR	-	expression tag	UNP Q5HZB2
A	673	GLN	GLU	engineered mutation	UNP Q5HZB2
B	386	MET	-	expression tag	UNP Q5HZB2
B	387	ARG	-	expression tag	UNP Q5HZB2
B	388	GLY	-	expression tag	UNP Q5HZB2
B	389	SER	-	expression tag	UNP Q5HZB2
B	390	HIS	-	expression tag	UNP Q5HZB2
B	391	HIS	-	expression tag	UNP Q5HZB2
B	392	HIS	-	expression tag	UNP Q5HZB2
B	393	HIS	-	expression tag	UNP Q5HZB2
B	394	HIS	-	expression tag	UNP Q5HZB2

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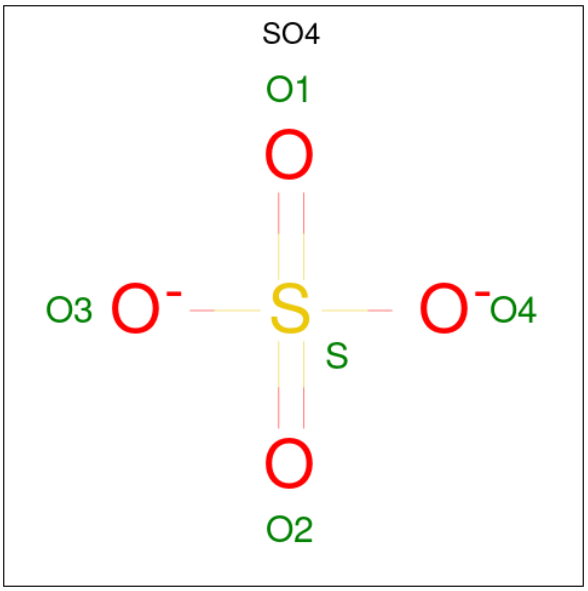
Chain	Residue	Modelled	Actual	Comment	Reference
B	395	THR	-	expression tag	UNP Q5HZB2
B	396	THR	-	expression tag	UNP Q5HZB2
B	673	GLN	GLU	engineered mutation	UNP Q5HZB2
C	386	MET	-	expression tag	UNP Q5HZB2
C	387	ARG	-	expression tag	UNP Q5HZB2
C	388	GLY	-	expression tag	UNP Q5HZB2
C	389	SER	-	expression tag	UNP Q5HZB2
C	390	HIS	-	expression tag	UNP Q5HZB2
C	391	HIS	-	expression tag	UNP Q5HZB2
C	392	HIS	-	expression tag	UNP Q5HZB2
C	393	HIS	-	expression tag	UNP Q5HZB2
C	394	HIS	-	expression tag	UNP Q5HZB2
C	395	THR	-	expression tag	UNP Q5HZB2
C	396	THR	-	expression tag	UNP Q5HZB2
C	673	GLN	GLU	engineered mutation	UNP Q5HZB2
D	386	MET	-	expression tag	UNP Q5HZB2
D	387	ARG	-	expression tag	UNP Q5HZB2
D	388	GLY	-	expression tag	UNP Q5HZB2
D	389	SER	-	expression tag	UNP Q5HZB2
D	390	HIS	-	expression tag	UNP Q5HZB2
D	391	HIS	-	expression tag	UNP Q5HZB2
D	392	HIS	-	expression tag	UNP Q5HZB2
D	393	HIS	-	expression tag	UNP Q5HZB2
D	394	HIS	-	expression tag	UNP Q5HZB2
D	395	THR	-	expression tag	UNP Q5HZB2
D	396	THR	-	expression tag	UNP Q5HZB2
D	673	GLN	GLU	engineered mutation	UNP Q5HZB2

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



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

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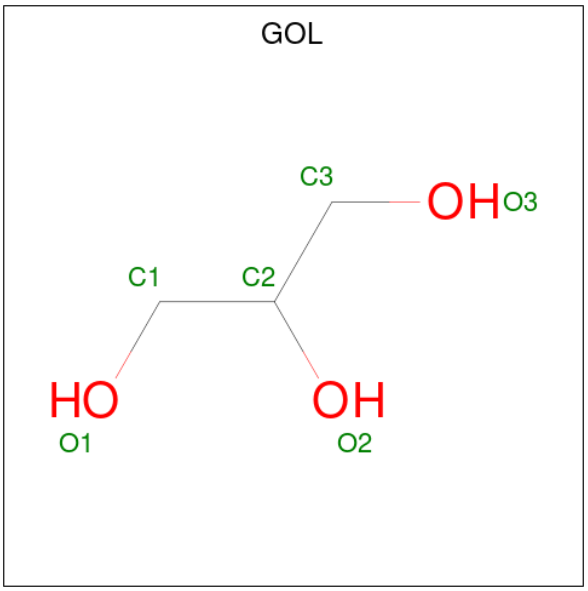
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
2	B	1	5	4	1	0	0
			Total	O	S		
2	C	1	5	4	1	0	0
			Total	O	S		
2	C	1	5	4	1	0	0
			Total	O	S		
2	C	1	5	4	1	0	0
			Total	O	S		
2	C	1	5	4	1	0	0
			Total	O	S		
2	C	1	5	4	1	0	0
			Total	O	S		
2	C	1	5	4	1	0	0
			Total	O	S		
2	C	1	5	4	1	0	0
			Total	O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

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

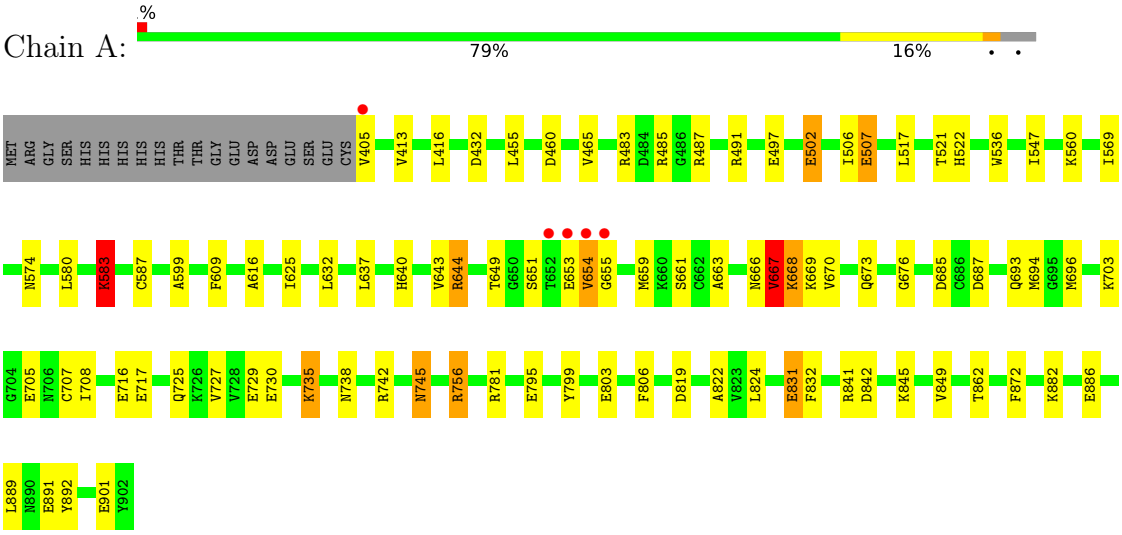
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	179	179	179	0	0
			Total	O		
4	B	156	156	156	0	0
			Total	O		
4	C	179	179	179	0	0
			Total	O		
4	D	169	169	169	0	0

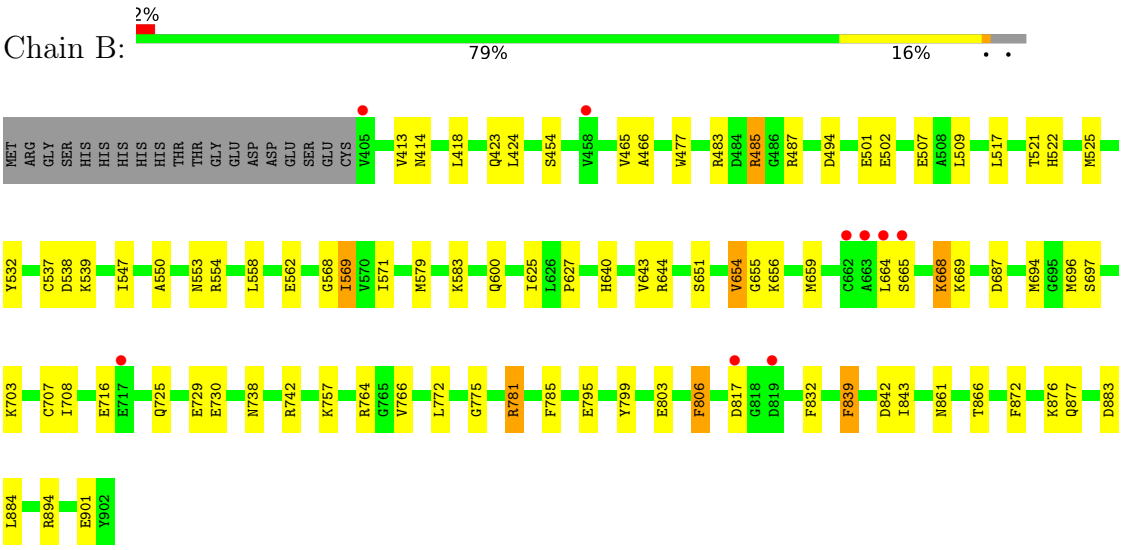
3 Residue-property plots (i)

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.

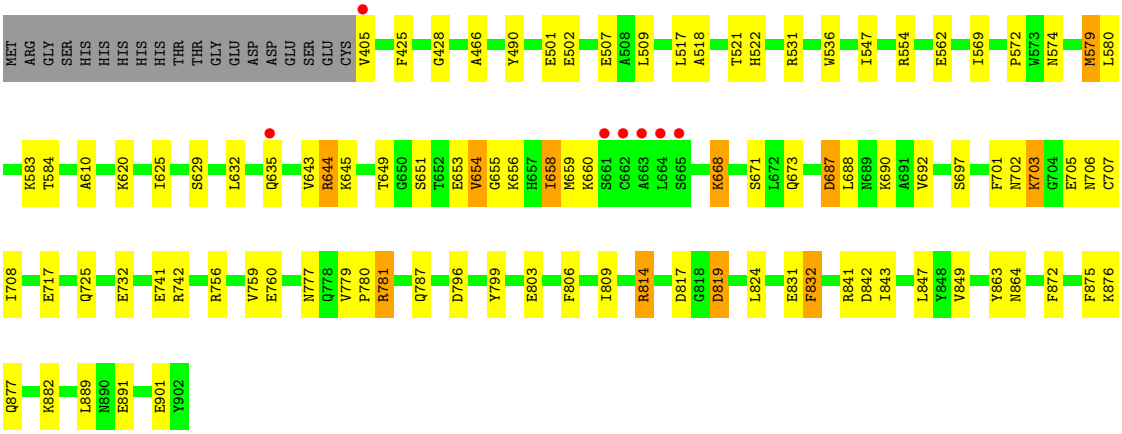
• Molecule 1: Aldehyde dehydrogenase 1 family, member L1



• Molecule 1: Aldehyde dehydrogenase 1 family, member L1

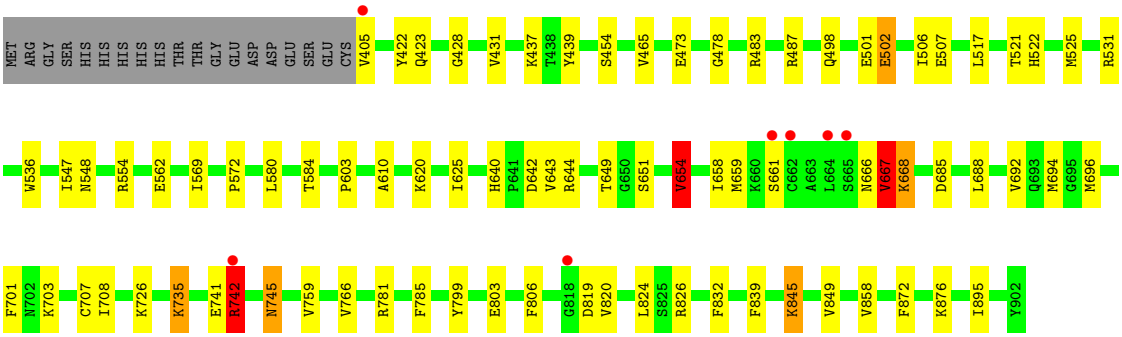


• Molecule 1: Aldehyde dehydrogenase 1 family, member L1



- Molecule 1: Aldehyde dehydrogenase 1 family, member L1

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	261.61Å 193.72Å 97.31Å 90.00° 109.05° 90.00°	Depositor
Resolution (Å)	49.81 – 2.38 49.81 – 2.38	Depositor EDS
% Data completeness (in resolution range)	96.7 (49.81-2.38) 96.7 (49.81-2.38)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.14 (at 2.37Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.171 , 0.202 0.169 , 0.199	Depositor DCC
R_{free} test set	8869 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å^2)	38.3	Xtriage
Anisotropy	0.167	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 38.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.012 for -h-2*l,-k,l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	16121	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 3.67% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: SO4, 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	1.40	19/3921 (0.5%)	1.04	13/5309 (0.2%)
1	B	1.33	16/3914 (0.4%)	1.05	19/5300 (0.4%)
1	C	1.46	17/3892 (0.4%)	1.14	17/5271 (0.3%)
1	D	1.36	14/3921 (0.4%)	1.05	10/5309 (0.2%)
All	All	1.39	66/15648 (0.4%)	1.07	59/21189 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	D	0	1
All	All	0	2

All (66) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	760	GLU	CG-CD	10.94	1.68	1.51
1	A	831	GLU	CG-CD	10.29	1.67	1.51
1	D	707	CYS	CB-SG	9.22	1.98	1.82
1	B	502	GLU	CD-OE2	9.06	1.35	1.25
1	A	831	GLU	CB-CG	8.54	1.68	1.52
1	A	560	LYS	CG-CD	-8.37	1.24	1.52
1	B	707	CYS	CB-SG	8.14	1.96	1.82
1	A	831	GLU	CD-OE1	7.96	1.34	1.25
1	C	562	GLU	CG-CD	7.93	1.63	1.51
1	C	707	CYS	CB-SG	7.80	1.95	1.82
1	C	732	GLU	CG-CD	7.70	1.63	1.51
1	D	501	GLU	CD-OE2	7.34	1.33	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	466	ALA	CA-CB	7.30	1.67	1.52
1	A	507	GLU	CG-CD	6.93	1.62	1.51
1	D	742	ARG	CG-CD	6.84	1.69	1.51
1	B	562	GLU	CG-CD	6.76	1.62	1.51
1	B	795	GLU	CB-CG	6.72	1.65	1.52
1	C	831	GLU	CG-CD	6.68	1.61	1.51
1	C	701	PHE	CB-CG	-6.63	1.40	1.51
1	A	707	CYS	CB-SG	6.46	1.93	1.82
1	C	732	GLU	CB-CG	6.35	1.64	1.52
1	C	741	GLU	CG-CD	6.34	1.61	1.51
1	B	501	GLU	CD-OE2	6.32	1.32	1.25
1	B	562	GLU	CD-OE1	6.31	1.32	1.25
1	A	609	PHE	CE1-CZ	6.27	1.49	1.37
1	C	501	GLU	CD-OE2	6.18	1.32	1.25
1	B	532	TYR	CD2-CE2	6.13	1.48	1.39
1	A	669	LYS	CE-NZ	6.05	1.64	1.49
1	D	667	VAL	CA-CB	6.01	1.67	1.54
1	D	562	GLU	CD-OE1	5.86	1.32	1.25
1	B	806	PHE	CD2-CE2	5.77	1.50	1.39
1	A	886	GLU	CG-CD	5.76	1.60	1.51
1	B	507	GLU	CG-CD	5.75	1.60	1.51
1	C	405	VAL	CA-CB	5.68	1.66	1.54
1	D	858	VAL	CB-CG1	5.66	1.64	1.52
1	A	502	GLU	CD-OE2	5.64	1.31	1.25
1	C	891	GLU	CD-OE2	5.59	1.31	1.25
1	D	507	GLU	CG-CD	5.58	1.60	1.51
1	C	875	PHE	CE1-CZ	5.50	1.47	1.37
1	C	502	GLU	CD-OE2	5.46	1.31	1.25
1	A	727	VAL	CB-CG1	-5.43	1.41	1.52
1	D	498	GLN	CG-CD	5.42	1.63	1.51
1	A	599	ALA	CA-CB	5.40	1.63	1.52
1	A	795	GLU	CB-CG	5.36	1.62	1.52
1	B	502	GLU	CG-CD	5.35	1.59	1.51
1	D	759	VAL	CA-CB	5.33	1.66	1.54
1	A	405	VAL	CB-CG2	5.28	1.64	1.52
1	C	759	VAL	CB-CG1	5.27	1.64	1.52
1	D	502	GLU	CG-CD	5.26	1.59	1.51
1	D	422	TYR	CD2-CE2	5.25	1.47	1.39
1	A	497	GLU	CD-OE2	5.22	1.31	1.25
1	A	716	GLU	CG-CD	5.22	1.59	1.51
1	B	502	GLU	CD-OE1	5.20	1.31	1.25
1	C	832	PHE	CE2-CZ	5.20	1.47	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	766	VAL	CB-CG1	5.19	1.63	1.52
1	D	502	GLU	CD-OE2	5.11	1.31	1.25
1	B	537	CYS	CB-SG	5.11	1.91	1.82
1	D	839	PHE	CE2-CZ	5.10	1.47	1.37
1	A	413	VAL	CB-CG1	-5.10	1.42	1.52
1	C	507	GLU	CG-CD	5.07	1.59	1.51
1	A	670	VAL	CB-CG2	-5.04	1.42	1.52
1	B	766	VAL	CB-CG1	5.04	1.63	1.52
1	B	839	PHE	CD2-CE2	5.02	1.49	1.39
1	B	716	GLU	CG-CD	5.01	1.59	1.51
1	A	616	ALA	CA-CB	5.01	1.62	1.52
1	B	466	ALA	CA-CB	5.00	1.62	1.52

All (59) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	814	ARG	NE-CZ-NH2	-13.02	113.79	120.30
1	C	644	ARG	NE-CZ-NH2	-12.38	114.11	120.30
1	C	531	ARG	NE-CZ-NH2	9.07	124.83	120.30
1	B	894	ARG	NE-CZ-NH1	-8.78	115.91	120.30
1	C	742	ARG	NE-CZ-NH2	-8.76	115.92	120.30
1	D	483	ARG	NE-CZ-NH2	-8.43	116.08	120.30
1	C	644	ARG	NE-CZ-NH1	8.35	124.47	120.30
1	C	842	ASP	CB-CG-OD2	7.66	125.19	118.30
1	B	687	ASP	CB-CG-OD2	7.46	125.02	118.30
1	A	491	ARG	NE-CZ-NH2	-7.37	116.61	120.30
1	C	796	ASP	CB-CG-OD2	7.12	124.70	118.30
1	D	642	ASP	CB-CG-OD2	7.03	124.63	118.30
1	B	764	ARG	NE-CZ-NH2	-6.87	116.87	120.30
1	A	485	ARG	NE-CZ-NH2	-6.80	116.90	120.30
1	C	687	ASP	CB-CG-OD2	6.54	124.18	118.30
1	B	487	ARG	NE-CZ-NH2	-6.53	117.04	120.30
1	C	531	ARG	NE-CZ-NH1	-6.38	117.11	120.30
1	B	554	ARG	NE-CZ-NH2	-6.34	117.13	120.30
1	B	483	ARG	NE-CZ-NH1	6.32	123.46	120.30
1	B	494	ASP	CB-CG-OD2	6.15	123.84	118.30
1	C	579	MET	CG-SD-CE	-6.09	90.45	100.20
1	C	817	ASP	CB-CG-OD1	5.96	123.67	118.30
1	A	694	MET	CG-SD-CE	5.87	109.59	100.20
1	B	483	ARG	NE-CZ-NH2	-5.84	117.38	120.30
1	B	694	MET	CG-SD-CE	5.83	109.54	100.20
1	D	694	MET	CG-SD-CE	5.82	109.51	100.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	416	LEU	CB-CG-CD1	-5.75	101.22	111.00
1	B	842	ASP	CB-CG-OD2	5.72	123.45	118.30
1	B	509	LEU	CB-CG-CD1	-5.70	101.31	111.00
1	D	531	ARG	NE-CZ-NH2	-5.67	117.47	120.30
1	A	432	ASP	CB-CG-OD1	5.65	123.39	118.30
1	D	473	GLU	CA-CB-CG	-5.62	101.03	113.40
1	B	554	ARG	NE-CZ-NH1	5.57	123.08	120.30
1	C	814	ARG	CG-CD-NE	-5.56	100.13	111.80
1	A	460	ASP	CB-CG-OD2	-5.55	113.31	118.30
1	D	667	VAL	N-CA-CB	5.52	123.65	111.50
1	B	485	ARG	NE-CZ-NH2	-5.52	117.54	120.30
1	B	817	ASP	CB-CG-OD1	5.50	123.25	118.30
1	D	487	ARG	NE-CZ-NH1	5.47	123.04	120.30
1	C	756	ARG	NE-CZ-NH1	5.47	123.04	120.30
1	D	644	ARG	NE-CZ-NH2	-5.43	117.59	120.30
1	A	455	LEU	CA-CB-CG	5.34	127.59	115.30
1	B	569	ILE	CG1-CB-CG2	-5.33	99.67	111.40
1	B	538	ASP	CB-CG-OD1	5.32	123.08	118.30
1	B	644	ARG	NE-CZ-NH1	-5.32	117.64	120.30
1	A	756	ARG	NE-CZ-NH1	5.26	122.93	120.30
1	D	826	ARG	NE-CZ-NH2	-5.24	117.68	120.30
1	A	583	LYS	CB-CG-CD	-5.21	98.06	111.60
1	A	841	ARG	NE-CZ-NH1	5.18	122.89	120.30
1	B	568	GLY	N-CA-C	-5.18	100.16	113.10
1	B	487	ARG	NE-CZ-NH1	5.10	122.85	120.30
1	C	644	ARG	CG-CD-NE	-5.10	101.10	111.80
1	A	842	ASP	CB-CG-OD2	5.05	122.84	118.30
1	C	554	ARG	NE-CZ-NH2	-5.05	117.78	120.30
1	D	642	ASP	CB-CG-OD1	-5.04	113.76	118.30
1	A	637	LEU	CB-CG-CD1	-5.02	102.47	111.00
1	A	687	ASP	CB-CG-OD2	5.02	122.82	118.30
1	C	809	ILE	CG1-CB-CG2	-5.01	100.38	111.40
1	C	756	ARG	NE-CZ-NH2	-5.01	117.80	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	882	LYS	Peptide
1	D	654	VAL	Peptide

5.2 Too-close contacts (i)

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	3826	0	3807	41	0
1	B	3817	0	3796	32	0
1	C	3806	0	3784	44	0
1	D	3825	0	3802	37	0
2	A	40	0	0	1	0
2	B	35	0	0	2	0
2	C	40	0	0	0	0
2	D	25	0	0	1	0
3	A	6	0	8	0	0
3	B	6	0	8	1	0
3	C	6	0	8	0	0
3	D	6	0	8	0	0
4	A	179	0	0	3	0
4	B	156	0	0	1	0
4	C	179	0	0	1	0
4	D	169	0	0	1	0
All	All	16121	0	15221	146	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 5.

All (146) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:799:TYR:CE2	1:B:803[B]:GLU:HG3	2.05	0.91
1:C:799:TYR:CE2	1:C:803[A]:GLU:HG3	2.11	0.85
1:D:610:ALA:HB2	1:D:625:ILE:HD12	1.58	0.85
1:A:654:VAL:CB	1:A:655:GLY:HA3	2.07	0.83
1:D:799:TYR:CE2	1:D:803[A]:GLU:HG3	2.15	0.81
1:B:799:TYR:CZ	1:B:803[B]:GLU:HG3	2.16	0.80
1:D:651:SER:OG	1:D:654:VAL:CB	2.30	0.79
1:A:651:SER:O	1:A:654:VAL:CB	2.31	0.79
1:C:687:ASP:OD1	1:C:690:LYS:HD2	1.86	0.76
1:A:725:GLN:NE2	1:A:729:GLU:OE2	2.19	0.75
1:B:843:ILE:HD12	1:C:843:ILE:HD12	1.73	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:799:TYR:CZ	1:A:803[A]:GLU:HG3	2.28	0.68
1:A:824:LEU:HD21	1:A:849:VAL:HG13	1.77	0.67
1:B:725:GLN:OE1	1:B:729:GLU:OE2	2.13	0.66
1:C:799:TYR:CZ	1:C:803[A]:GLU:HG3	2.30	0.66
1:B:569:ILE:HD12	1:B:571:ILE:HD11	1.76	0.66
1:B:643:VAL:O	1:B:668:LYS:HE3	1.96	0.66
1:C:799:TYR:CE2	1:C:803[A]:GLU:CG	2.79	0.65
1:A:799:TYR:CE2	1:A:803[A]:GLU:HG3	2.33	0.63
1:B:414:ASN:OD1	1:B:742:ARG:NH2	2.31	0.63
1:B:669:LYS:NZ	2:B:3024:SO4:O3	2.31	0.63
1:C:651:SER:OG	1:C:654:VAL:CB	2.47	0.61
1:C:702:ASN:O	1:C:703:LYS:HG2	2.01	0.61
1:C:777:ASN:O	1:C:787:GLN:HG3	2.01	0.61
1:C:643:VAL:O	1:C:668:LYS:HE3	2.00	0.60
1:C:717:GLU:OE1	1:C:814:ARG:HD2	2.02	0.60
1:A:654:VAL:CB	1:A:655:GLY:CA	2.77	0.60
1:D:423:GLN:HA	1:D:454:SER:OG	2.02	0.60
1:A:756:ARG:NH2	4:A:1134:HOH:O	2.35	0.58
1:B:558:LEU:C	1:B:558:LEU:HD12	2.24	0.58
1:C:572:PRO:HG2	1:C:579:MET:HG3	1.84	0.58
1:C:425:PHE:CD2	1:C:610:ALA:HB1	2.38	0.58
1:B:708:ILE:HG21	1:B:872:PHE:HE1	1.69	0.58
1:B:651:SER:OG	1:B:654:VAL:CB	2.53	0.56
1:C:654:VAL:N	1:C:655:GLY:HA3	2.20	0.56
1:A:487:ARG:NH1	1:C:490[B]:TYR:OH	2.33	0.56
1:C:824:LEU:HD21	1:C:849:VAL:HG13	1.86	0.56
1:A:521:THR:O	1:A:522:HIS:C	2.43	0.56
1:A:799:TYR:CE2	1:A:803[A]:GLU:CG	2.89	0.55
1:D:781[B]:ARG:HD2	1:D:785:PHE:CD2	2.42	0.55
1:A:502:GLU:O	1:A:506:ILE:HG13	2.07	0.54
1:B:521:THR:O	1:B:522:HIS:C	2.46	0.54
1:A:738:ASN:HB2	2:A:3020:SO4:O2	2.09	0.53
1:C:654:VAL:N	1:C:655:GLY:CA	2.72	0.53
1:D:439:TYR:OH	1:D:603:PRO:HG3	2.09	0.53
1:A:651:SER:OG	1:A:654:VAL:CB	2.56	0.52
1:C:779:VAL:HG22	1:C:787:GLN:HG2	1.90	0.52
1:C:877:GLN:NE2	1:D:666:ASN:O	2.41	0.52
1:D:708:ILE:CG2	1:D:872:PHE:HE1	2.23	0.51
1:A:651:SER:C	1:A:654:VAL:CB	2.79	0.51
1:D:502:GLU:O	1:D:506:ILE:HG13	2.09	0.51
1:A:587:CYS:HB2	1:A:892:TYR:CE1	2.46	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:901:GLU:OE2	4:A:1039:HOH:O	2.19	0.51
1:C:708:ILE:HG21	1:C:872:PHE:HE1	1.75	0.51
1:B:781[A]:ARG:HD2	1:B:785:PHE:CD2	2.45	0.51
1:C:901:GLU:OE2	4:C:1048:HOH:O	2.19	0.50
1:D:832:PHE:O	1:D:876:LYS:HE3	2.11	0.50
1:D:569:ILE:HD11	1:D:584:THR:OG1	2.12	0.50
1:D:685:ASP:OD2	1:D:845:LYS:NZ	2.45	0.50
1:C:521:THR:O	1:C:522:HIS:C	2.49	0.49
1:C:863:TYR:O	1:C:864:ASN:HB2	2.13	0.49
1:C:580:LEU:C	1:C:580:LEU:HD23	2.33	0.49
1:C:632:LEU:C	1:C:632:LEU:HD23	2.33	0.48
1:B:832:PHE:O	1:B:876:LYS:HE2	2.13	0.48
1:B:708:ILE:CG2	1:B:872:PHE:HE1	2.26	0.48
1:C:688:LEU:O	1:C:692:VAL:HG23	2.14	0.48
1:C:882:LYS:HE3	1:D:895:ILE:HB	1.96	0.48
1:D:572:PRO:HD3	1:D:649:THR:HB	1.95	0.48
1:C:819:ASP:C	1:C:819:ASP:OD2	2.52	0.48
1:A:666:ASN:O	1:B:877:GLN:NE2	2.43	0.48
1:D:824:LEU:HD21	1:D:849:VAL:HG13	1.95	0.48
1:C:569:ILE:HD11	1:C:584:THR:OG1	2.14	0.47
1:D:643:VAL:O	1:D:668:LYS:HE3	2.14	0.47
1:C:653:GLU:CA	1:C:654:VAL:CB	2.92	0.47
1:D:554:ARG:NE	2:D:3016:SO4:O4	2.34	0.47
1:B:424:LEU:HD23	1:B:627:PRO:HD2	1.95	0.47
1:C:673:GLN:NE2	1:C:872:PHE:HE2	2.12	0.47
1:D:428:GLY:HA3	1:D:620:LYS:HG2	1.97	0.47
1:A:483:ARG:HB3	1:D:548:ASN:HD21	1.79	0.47
1:A:708:ILE:HG21	1:A:872:PHE:HE1	1.79	0.47
1:C:583:LYS:HE2	1:C:649:THR:OG1	2.15	0.47
1:A:653:GLU:CB	1:A:654:VAL:HA	2.45	0.46
1:D:708:ILE:CG2	1:D:872:PHE:CE1	2.98	0.46
1:A:666:ASN:O	1:A:667:VAL:HG22	2.16	0.46
1:B:423:GLN:HA	1:B:454:SER:OG	2.16	0.46
1:D:692:VAL:O	1:D:696:MET:HG2	2.16	0.46
1:C:658:ILE:O	1:C:660:LYS:N	2.49	0.46
1:D:423:GLN:HB3	1:D:431:VAL:O	2.15	0.45
1:D:610:ALA:HB2	1:D:625:ILE:CD1	2.40	0.45
1:B:654:VAL:N	1:B:655:GLY:HA3	2.30	0.45
1:D:666:ASN:O	1:D:667:VAL:HG13	2.16	0.45
1:D:688:LEU:HD23	1:D:726:LYS:HE3	1.98	0.45
1:C:428:GLY:HA3	1:C:620:LYS:HG2	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:645:LYS:NZ	1:C:671:SER:HB3	2.32	0.45
1:B:883:ASP:O	1:B:884:LEU:HB2	2.16	0.45
1:D:517:LEU:HD11	1:D:701:PHE:CZ	2.51	0.45
1:A:644:ARG:O	1:A:644:ARG:HD2	2.17	0.45
1:B:738:ASN:HB2	2:B:3021:SO4:O1	2.17	0.45
1:B:696:MET:SD	1:B:730:GLU:HB2	2.57	0.45
1:D:465:VAL:HG11	1:D:640:HIS:CE1	2.52	0.44
1:A:632:LEU:C	1:A:632:LEU:HD23	2.38	0.44
1:C:574:ASN:HB2	1:C:705:GLU:O	2.17	0.44
1:A:661:SER:HA	4:A:1119:HOH:O	2.17	0.44
1:D:799:TYR:CZ	1:D:803[A]:GLU:HG3	2.52	0.44
1:B:465:VAL:HG11	1:B:640:HIS:CE1	2.52	0.44
1:D:735:LYS:H	1:D:745:ASN:HD21	1.66	0.44
1:A:891:GLU:H	1:A:891:GLU:HG2	1.67	0.43
1:B:477:TRP:CH2	1:B:485:ARG:HG3	2.52	0.43
1:B:600:GLN:HA	4:B:1104:HOH:O	2.17	0.43
1:D:580:LEU:C	1:D:580:LEU:HD23	2.38	0.43
1:D:742:ARG:HD3	1:D:742:ARG:HA	1.90	0.43
1:A:574:ASN:HB2	1:A:705:GLU:O	2.18	0.43
1:C:518:ALA:HA	1:C:522:HIS:HB2	2.00	0.43
1:A:643:VAL:O	1:A:668:LYS:HE3	2.19	0.43
1:D:521:THR:O	1:D:522:HIS:C	2.56	0.43
1:C:536:TRP:CG	1:C:889:LEU:HD11	2.54	0.43
1:D:478:GLY:HA3	4:D:1112:HOH:O	2.19	0.43
1:B:866:THR:OG1	3:B:2002:GOL:H11	2.19	0.42
1:C:847:LEU:HD23	1:C:847:LEU:HA	1.78	0.42
1:C:687:ASP:HB2	1:C:841:ARG:NH2	2.34	0.42
1:B:839:PHE:HA	1:B:861:ASN:OD1	2.20	0.42
1:C:572:PRO:CG	1:C:579:MET:HG3	2.49	0.42
1:A:517:LEU:CD2	1:A:522:HIS:CE1	3.03	0.42
1:A:685:ASP:OD2	1:A:845:LYS:NZ	2.52	0.41
1:A:649:THR:HG23	1:A:673:GLN:HB3	2.02	0.41
1:C:832:PHE:O	1:C:876:LYS:HE2	2.21	0.41
1:A:536:TRP:CG	1:A:889:LEU:HD11	2.56	0.41
1:B:539:LYS:HD3	1:D:536:TRP:CE2	2.55	0.41
1:C:654:VAL:H	1:C:655:GLY:CA	2.34	0.41
1:B:424:LEU:CD2	1:B:627:PRO:HD2	2.51	0.41
1:B:550:ALA:O	1:B:553:ASN:HB2	2.21	0.41
1:C:610:ALA:HB2	1:C:625:ILE:HD12	2.01	0.41
1:C:780:PRO:O	1:C:781:ARG:HB3	2.21	0.40
1:A:465:VAL:HG11	1:A:640:HIS:CE1	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:569:ILE:HD12	1:A:583:LYS:HB2	2.02	0.40
1:A:735:LYS:H	1:A:745:ASN:HD21	1.69	0.40
1:A:819:ASP:HB3	1:A:822:ALA:HB3	2.03	0.40
1:B:772:LEU:HD21	1:B:775:GLY:O	2.21	0.40
1:D:569:ILE:CD1	1:D:584:THR:OG1	2.69	0.40
1:A:580:LEU:HD23	1:A:580:LEU:C	2.42	0.40
1:A:676:GLY:HA2	1:A:832:PHE:HB3	2.04	0.40
1:A:745:ASN:C	1:A:745:ASN:HD22	2.24	0.40
1:D:708:ILE:HG21	1:D:872:PHE:HE1	1.86	0.40
1:A:696:MET:SD	1:A:730:GLU:HB2	2.61	0.40
1:A:862:THR:HB	1:B:901:GLU:HB2	2.03	0.40
1:D:569:ILE:HG21	1:D:569:ILE:HD13	1.84	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	502/517 (97%)	482 (96%)	16 (3%)	4 (1%)	19	27
1	B	501/517 (97%)	472 (94%)	24 (5%)	5 (1%)	15	21
1	C	499/517 (96%)	475 (95%)	19 (4%)	5 (1%)	15	21
1	D	502/517 (97%)	473 (94%)	23 (5%)	6 (1%)	13	17
All	All	2004/2068 (97%)	1902 (95%)	82 (4%)	20 (1%)	15	21

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	663	ALA
1	B	654	VAL
1	B	659	MET

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Mol	Chain	Res	Type
1	B	665	SER
1	C	654	VAL
1	C	656	LYS
1	C	659	MET
1	D	654	VAL
1	D	658	ILE
1	A	659	MET
1	B	656	LYS
1	B	664	LEU
1	A	667	VAL
1	D	659	MET
1	D	667	VAL
1	A	654	VAL
1	D	661	SER
1	C	781	ARG
1	C	658	ILE
1	D	820	VAL

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	402/426 (94%)	385 (96%)	17 (4%)	30	44
1	B	401/426 (94%)	386 (96%)	15 (4%)	34	50
1	C	399/426 (94%)	386 (97%)	13 (3%)	38	55
1	D	402/426 (94%)	389 (97%)	13 (3%)	39	56
All	All	1604/1704 (94%)	1546 (96%)	58 (4%)	36	51

All (58) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	507	GLU
1	A	547	ILE
1	A	583	LYS
1	A	625	ILE

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Mol	Chain	Res	Type
1	A	644	ARG
1	A	667	VAL
1	A	668	LYS
1	A	693	GLN
1	A	703	LYS
1	A	717	GLU
1	A	735	LYS
1	A	742	ARG
1	A	745	ASN
1	A	781[A]	ARG
1	A	781[B]	ARG
1	A	806	PHE
1	A	831	GLU
1	B	413	VAL
1	B	418	LEU
1	B	517	LEU
1	B	525	MET
1	B	547	ILE
1	B	579	MET
1	B	583	LYS
1	B	625	ILE
1	B	668	LYS
1	B	697	SER
1	B	703	LYS
1	B	757	LYS
1	B	781[A]	ARG
1	B	781[B]	ARG
1	B	806	PHE
1	C	509	LEU
1	C	517	LEU
1	C	547	ILE
1	C	629	SER
1	C	635	GLN
1	C	644	ARG
1	C	668	LYS
1	C	697	SER
1	C	703	LYS
1	C	706	ASN
1	C	725	GLN
1	C	806	PHE
1	C	819	ASP
1	D	405	VAL

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Mol	Chain	Res	Type
1	D	437	LYS
1	D	525	MET
1	D	547	ILE
1	D	668	LYS
1	D	703	LYS
1	D	735	LYS
1	D	741	GLU
1	D	742	ARG
1	D	745	ASN
1	D	806	PHE
1	D	819	ASP
1	D	845	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	600	GLN
1	A	693	GLN
1	A	706	ASN
1	A	745	ASN
1	A	750	ASN
1	A	877	GLN
1	B	407	ASN
1	B	673	GLN
1	B	706	ASN
1	B	725	GLN
1	B	750	ASN
1	C	407	ASN
1	C	600	GLN
1	C	673	GLN
1	C	706	ASN
1	C	725	GLN
1	C	750	ASN
1	D	407	ASN
1	D	457	GLN
1	D	548	ASN
1	D	673	GLN
1	D	706	ASN
1	D	745	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

32 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	SO4	B	3021	-	4,4,4	0.28	0	6,6,6	0.39	0
3	GOL	B	2002	-	5,5,5	0.36	0	5,5,5	0.92	0
2	SO4	D	3016	-	4,4,4	0.34	0	6,6,6	0.40	0
2	SO4	A	3011	-	4,4,4	0.23	0	6,6,6	0.53	0
2	SO4	A	3014	-	4,4,4	0.12	0	6,6,6	0.59	0
2	SO4	C	3010	-	4,4,4	0.29	0	6,6,6	0.51	0
2	SO4	A	3001	-	4,4,4	0.29	0	6,6,6	0.80	0
2	SO4	B	3004	-	4,4,4	0.16	0	6,6,6	0.76	0
2	SO4	B	3003	-	4,4,4	0.31	0	6,6,6	0.74	0
3	GOL	C	2003	-	5,5,5	0.70	0	5,5,5	0.91	0
2	SO4	C	3002	-	4,4,4	0.36	0	6,6,6	0.18	0
2	SO4	B	3024	-	4,4,4	0.26	0	6,6,6	1.36	1 (16%)
2	SO4	C	3019	-	4,4,4	0.33	0	6,6,6	0.58	0
2	SO4	A	3017	-	4,4,4	0.28	0	6,6,6	0.66	0
2	SO4	C	3013	-	4,4,4	0.28	0	6,6,6	0.32	0
2	SO4	B	3009	-	4,4,4	0.32	0	6,6,6	0.32	0
2	SO4	C	3022	-	4,4,4	0.28	0	6,6,6	1.26	1 (16%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	SO4	D	3008	-	4,4,4	0.52	0	6,6,6	0.37	0
2	SO4	A	3020	-	4,4,4	0.41	0	6,6,6	0.61	0
2	SO4	C	3007	-	4,4,4	0.23	0	6,6,6	0.85	0
2	SO4	B	3005	-	4,4,4	0.31	0	6,6,6	0.35	0
2	SO4	A	3025	-	4,4,4	0.31	0	6,6,6	1.64	2 (33%)
2	SO4	A	3006	-	4,4,4	0.37	0	6,6,6	0.98	0
3	GOL	A	2001	-	5,5,5	0.57	0	5,5,5	1.19	0
2	SO4	B	3015	-	4,4,4	0.33	0	6,6,6	0.82	0
2	SO4	D	3023	-	4,4,4	0.33	0	6,6,6	0.82	0
2	SO4	C	3027	-	4,4,4	0.37	0	6,6,6	1.20	0
2	SO4	C	3028	-	4,4,4	0.38	0	6,6,6	0.92	0
2	SO4	D	3012	-	4,4,4	0.26	0	6,6,6	0.35	0
3	GOL	D	2004	-	5,5,5	0.76	0	5,5,5	1.30	1 (20%)
2	SO4	D	3026	-	4,4,4	0.29	0	6,6,6	1.30	1 (16%)
2	SO4	A	3018	-	4,4,4	0.55	0	6,6,6	0.52	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
3	GOL	D	2004	-	-	2/4/4/4	-
3	GOL	B	2002	-	-	2/4/4/4	-
3	GOL	C	2003	-	-	2/4/4/4	-
3	GOL	A	2001	-	-	4/4/4/4	-

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	3025	SO4	O3-S-O1	-2.63	95.61	109.31
2	B	3024	SO4	O3-S-O1	-2.56	95.96	109.31
3	D	2004	GOL	O1-C1-C2	2.37	121.57	110.20
2	A	3025	SO4	O4-S-O3	2.32	118.97	109.06
2	C	3022	SO4	O3-S-O2	2.24	120.99	109.31
2	D	3026	SO4	O2-S-O1	-2.03	94.47	109.43

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	2001	GOL	C1-C2-C3-O3
3	B	2002	GOL	C1-C2-C3-O3
3	A	2001	GOL	O2-C2-C3-O3
3	A	2001	GOL	O1-C1-C2-C3
3	A	2001	GOL	O1-C1-C2-O2
3	D	2004	GOL	O1-C1-C2-O2
3	B	2002	GOL	O1-C1-C2-O2
3	C	2003	GOL	O1-C1-C2-O2
3	C	2003	GOL	O1-C1-C2-C3
3	D	2004	GOL	O1-C1-C2-C3

There are no ring outliers.

5 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	3021	SO4	1	0
3	B	2002	GOL	1	0
2	D	3016	SO4	1	0
2	B	3024	SO4	1	0
2	A	3020	SO4	1	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	498/517 (96%)	-0.20	5 (1%)	82	83	26, 38, 57, 110	22 (4%)
1	B	498/517 (96%)	-0.11	9 (1%)	68	70	25, 41, 63, 117	25 (5%)
1	C	498/517 (96%)	-0.26	7 (1%)	75	77	24, 36, 54, 112	21 (4%)
1	D	498/517 (96%)	-0.19	7 (1%)	75	77	25, 40, 59, 109	23 (4%)
All	All	1992/2068 (96%)	-0.19	28 (1%)	75	77	24, 39, 60, 117	91 (4%)

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	654	VAL	4.9
1	B	663	ALA	4.3
1	B	665	SER	3.8
1	B	662	CYS	3.6
1	D	661	SER	3.6
1	C	662	CYS	3.4
1	D	662	CYS	3.4
1	B	458	VAL	3.3
1	B	405	VAL	2.8
1	C	665	SER	2.7
1	D	818	GLY	2.7
1	C	635	GLN	2.7
1	D	405	VAL	2.7
1	A	653	GLU	2.6
1	B	664	LEU	2.5
1	C	663	ALA	2.5
1	D	742	ARG	2.5
1	B	817	ASP	2.4
1	C	664	LEU	2.3
1	A	652	THR	2.3
1	D	665	SER	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	819	ASP	2.2
1	C	405	VAL	2.2
1	C	661	SER	2.2
1	A	405	VAL	2.2
1	B	717	GLU	2.2
1	A	655	GLY	2.1
1	D	664	LEU	2.1

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 monosaccharides 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(Å^2)	Q<0.9
3	GOL	D	2004	6/6	0.83	0.23	75,77,80,81	0
2	SO4	A	3017	5/5	0.86	0.30	52,58,60,63	5
2	SO4	D	3023	5/5	0.86	0.21	54,56,58,60	5
2	SO4	A	3011	5/5	0.86	0.33	92,92,94,94	5
2	SO4	C	3010	5/5	0.87	0.23	80,80,84,87	5
2	SO4	B	3009	5/5	0.89	0.20	101,102,104,105	0
2	SO4	A	3018	5/5	0.90	0.20	40,46,48,52	5
2	SO4	C	3019	5/5	0.90	0.21	51,52,55,55	5
3	GOL	B	2002	6/6	0.91	0.17	61,69,72,73	0
3	GOL	A	2001	6/6	0.91	0.12	67,71,71,73	0
2	SO4	A	3020	5/5	0.92	0.19	34,37,39,42	5
2	SO4	A	3006	5/5	0.92	0.15	63,65,67,69	5
2	SO4	B	3005	5/5	0.93	0.23	75,76,78,80	5
3	GOL	C	2003	6/6	0.93	0.19	69,73,73,74	0
2	SO4	C	3007	5/5	0.93	0.15	69,70,73,73	5
2	SO4	D	3012	5/5	0.94	0.17	89,90,90,92	5

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	SO4	D	3008	5/5	0.94	0.17	61,62,64,67	5
2	SO4	C	3022	5/5	0.95	0.15	46,47,50,51	5
2	SO4	C	3027	5/5	0.95	0.16	58,63,66,66	5
2	SO4	B	3015	5/5	0.95	0.10	48,49,52,55	5
2	SO4	B	3024	5/5	0.95	0.25	63,66,70,72	5
2	SO4	D	3016	5/5	0.95	0.10	56,56,57,59	5
2	SO4	A	3014	5/5	0.96	0.19	41,42,47,48	5
2	SO4	A	3025	5/5	0.96	0.27	58,61,64,68	5
2	SO4	B	3021	5/5	0.96	0.11	49,49,52,54	5
2	SO4	D	3026	5/5	0.96	0.22	55,62,65,67	5
2	SO4	A	3001	5/5	0.97	0.17	44,46,49,53	5
2	SO4	B	3003	5/5	0.97	0.16	33,34,37,40	5
2	SO4	C	3028	5/5	0.97	0.11	42,51,55,57	5
2	SO4	C	3002	5/5	0.97	0.15	42,42,47,48	5
2	SO4	C	3013	5/5	0.98	0.11	51,53,55,58	5
2	SO4	B	3004	5/5	0.99	0.11	42,44,48,49	5

6.5 Other polymers (i)

There are no such residues in this entry.