



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:14 AM UTC

PDB ID : 9GGI / pdb_00009ggi
Title : Crystal structure of argininosuccinate lyase from Arabidopsis thaliana (AtASL)
Authors : Nielipinski, M.; Pietrzyk-Brzezinska, A.J.; Krzeszewska, D.; Sekula, B.
Deposited on : 2024-08-13
Resolution : 1.55 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

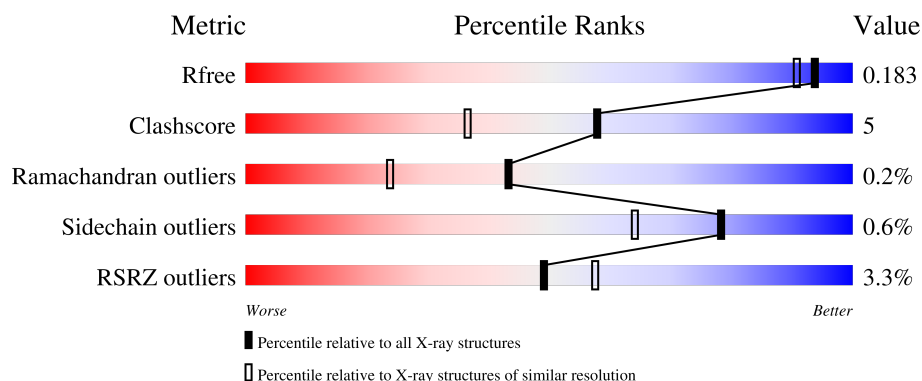
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.55 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.



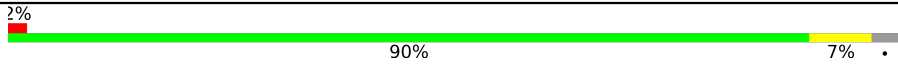
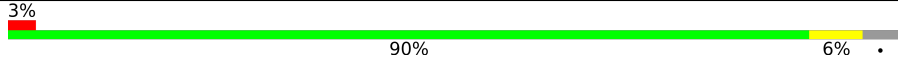
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2145 (1.56-1.56)
Clashscore	190562	2189 (1.56-1.56)
Ramachandran outliers	187476	2153 (1.56-1.56)
Sidechain outliers	187428	2150 (1.56-1.56)
RSRZ outliers	180081	2146 (1.56-1.56)

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	465	4% 83% 12% . .
1	B	465	3% 90% 7% .
1	C	465	3% 87% 8% . .
1	D	465	4% 88% 8% . .
1	E	465	3% 89% 7% .

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Mol	Chain	Length	Quality of chain
1	F	465	
1	G	465	
1	H	465	

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	SO4	A	604	-	-	X	-
3	GOL	E	607	-	-	X	-
3	GOL	E	608	-	-	X	-
3	GOL	H	607	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 33706 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 Argininosuccinate lyase, chloroplastic.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	450	Total	C	N	O	S	0	23	0
			3642	2308	613	699	22			
1	B	453	Total	C	N	O	S	0	15	0
			3632	2298	621	692	21			
1	C	448	Total	C	N	O	S	0	15	0
			3584	2268	611	685	20			
1	D	448	Total	C	N	O	S	0	15	0
			3584	2269	609	686	20			
1	E	451	Total	C	N	O	S	0	17	0
			3616	2289	614	693	20			
1	F	453	Total	C	N	O	S	0	19	0
			3651	2315	619	696	21			
1	G	448	Total	C	N	O	S	0	15	0
			3583	2268	610	685	20			
1	H	448	Total	C	N	O	S	0	16	0
			3586	2272	608	686	20			

There are 24 discrepancies between the modelled and reference sequences:

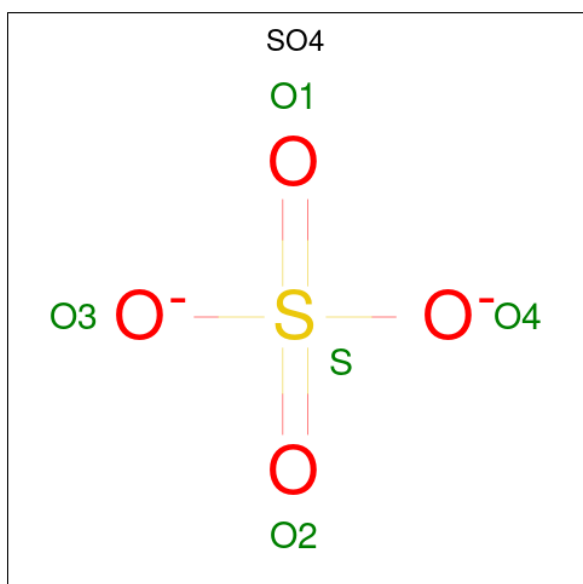
Chain	Residue	Modelled	Actual	Comment	Reference
A	53	SER	-	expression tag	UNP Q9LEU8
A	54	ASN	-	expression tag	UNP Q9LEU8
A	55	ALA	-	expression tag	UNP Q9LEU8
B	53	SER	-	expression tag	UNP Q9LEU8
B	54	ASN	-	expression tag	UNP Q9LEU8
B	55	ALA	-	expression tag	UNP Q9LEU8
C	53	SER	-	expression tag	UNP Q9LEU8
C	54	ASN	-	expression tag	UNP Q9LEU8
C	55	ALA	-	expression tag	UNP Q9LEU8
D	53	SER	-	expression tag	UNP Q9LEU8
D	54	ASN	-	expression tag	UNP Q9LEU8
D	55	ALA	-	expression tag	UNP Q9LEU8
E	53	SER	-	expression tag	UNP Q9LEU8

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Chain	Residue	Modelled	Actual	Comment	Reference
E	54	ASN	-	expression tag	UNP Q9LEU8
E	55	ALA	-	expression tag	UNP Q9LEU8
F	53	SER	-	expression tag	UNP Q9LEU8
F	54	ASN	-	expression tag	UNP Q9LEU8
F	55	ALA	-	expression tag	UNP Q9LEU8
G	53	SER	-	expression tag	UNP Q9LEU8
G	54	ASN	-	expression tag	UNP Q9LEU8
G	55	ALA	-	expression tag	UNP Q9LEU8
H	53	SER	-	expression tag	UNP Q9LEU8
H	54	ASN	-	expression tag	UNP Q9LEU8
H	55	ALA	-	expression tag	UNP Q9LEU8

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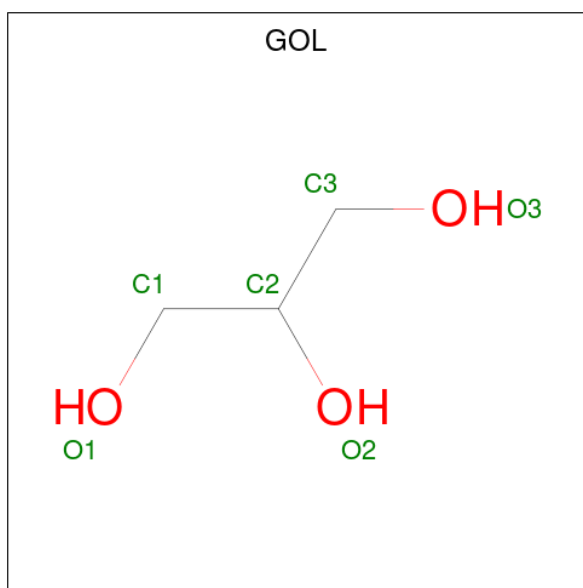
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total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

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

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



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

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	C	2	Total Cl 2 2	0	0
4	D	1	Total Cl 1 1	0	0
4	E	2	Total Cl 2 2	0	0
4	G	1	Total Cl 1 1	0	0

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

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

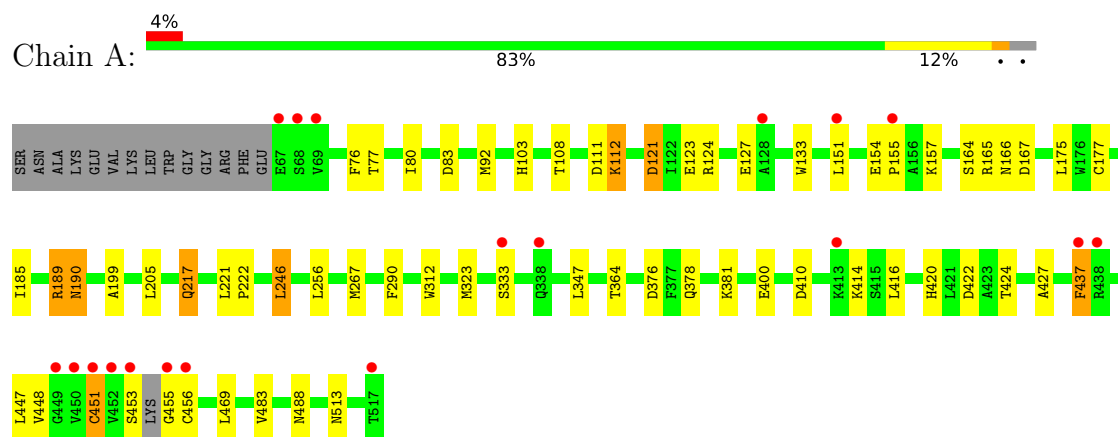
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	542	Total O 542 542	0	0
6	B	554	Total O 554 554	0	0
6	C	589	Total O 590 590	0	1
6	D	561	Total O 562 562	0	1
6	E	553	Total O 554 554	0	1
6	F	564	Total O 565 565	0	1
6	G	571	Total O 571 571	0	0
6	H	556	Total O 556 556	0	0

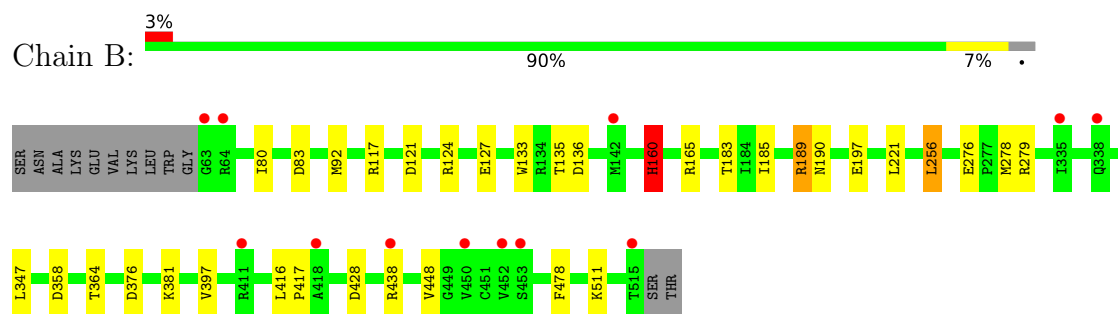
3 Residue-property plots [i](#)

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.

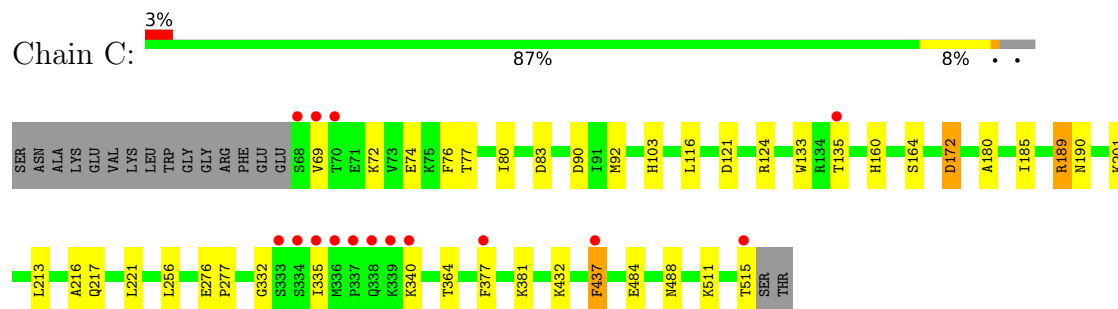
- Molecule 1: Argininosuccinate lyase, chloroplactic



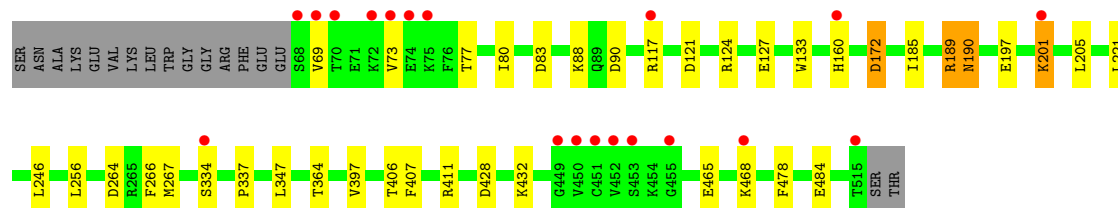
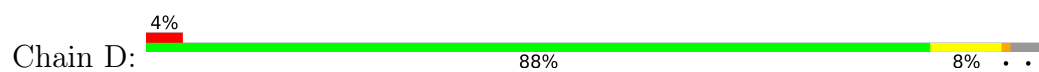
- Molecule 1: Argininosuccinate lyase, chloroplactic



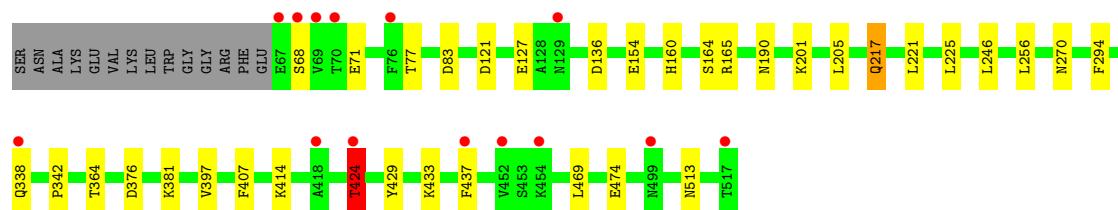
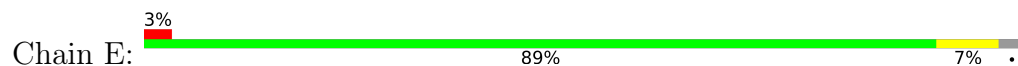
- Molecule 1: Argininosuccinate lyase, chloroplactic



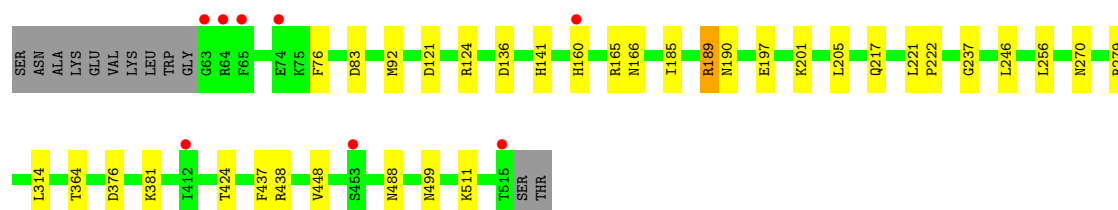
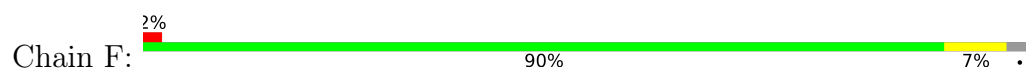
- Molecule 1: Argininosuccinate lyase, chloroplactic



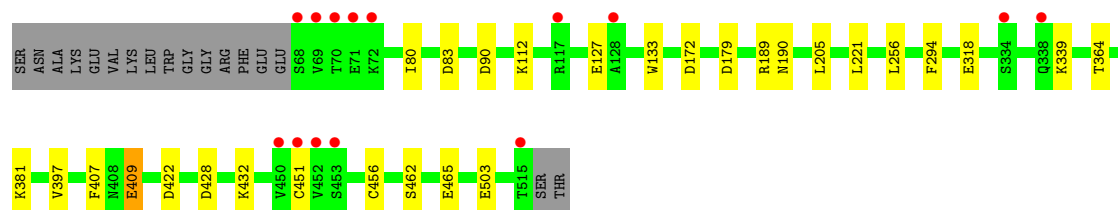
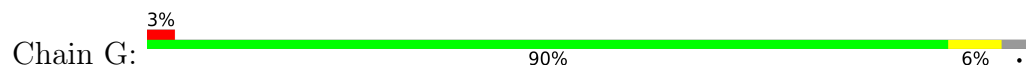
- Molecule 1: Argininosuccinate lyase, chloroplastic



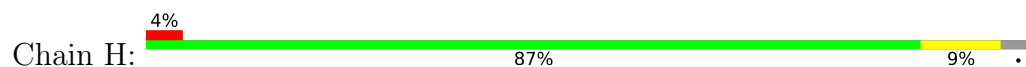
- Molecule 1: Argininosuccinate lyase, chloroplastic

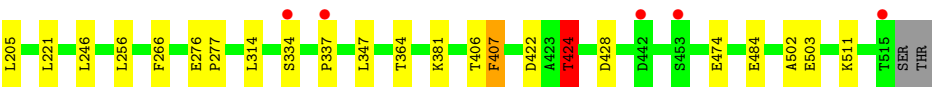


- Molecule 1: Argininosuccinate lyase, chloroplastic



- Molecule 1: Argininosuccinate lyase, chloroplastic





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	106.04Å 229.54Å 111.59Å 90.00° 90.56° 90.00°	Depositor
Resolution (Å)	48.45 – 1.55 48.45 – 1.55	Depositor EDS
% Data completeness (in resolution range)	94.1 (48.45-1.55) 94.1 (48.45-1.55)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.87 (at 1.55Å)	Xtriage
Refinement program	REFMAC 5.8.0425	Depositor
R, R_{free}	0.147 , 0.171 0.161 , 0.183	Depositor DCC
R_{free} test set	3603 reflections (0.50%)	wwPDB-VP
Wilson B-factor (Å ²)	16.8	Xtriage
Anisotropy	0.612	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 52.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.002 for l,k,-h 0.018 for h,-k,-l 0.013 for l,-k,h	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	33706	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.48% 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: SO4, NA, CL, 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	0.87	10/3767 (0.3%)	1.19	19/5093 (0.4%)
1	B	0.81	7/3741 (0.2%)	1.11	18/5057 (0.4%)
1	C	0.81	2/3692 (0.1%)	1.11	9/4990 (0.2%)
1	D	0.81	2/3692 (0.1%)	1.12	10/4994 (0.2%)
1	E	0.79	0/3731	1.15	16/5045 (0.3%)
1	F	0.80	5/3773 (0.1%)	1.13	10/5101 (0.2%)
1	G	0.75	0/3691	1.09	5/4992 (0.1%)
1	H	0.83	8/3697 (0.2%)	1.16	16/5000 (0.3%)
All	All	0.81	34/29784 (0.1%)	1.13	103/40272 (0.3%)

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	217[A]	GLN	C-O	7.63	1.30	1.24
1	A	217[B]	GLN	C-O	7.63	1.30	1.24
1	F	141	HIS	ND1-CE1	5.91	1.38	1.32
1	H	135[A]	THR	C-O	5.76	1.31	1.24
1	H	135[B]	THR	C-O	5.76	1.31	1.24
1	A	189[A]	ARG	C-O	5.76	1.30	1.24
1	A	189[B]	ARG	C-O	5.76	1.30	1.24
1	A	513[A]	ASN	C-O	5.68	1.31	1.23
1	A	513[B]	ASN	C-O	5.68	1.31	1.23
1	F	189[A]	ARG	C-O	5.63	1.30	1.24
1	F	189[B]	ARG	C-O	5.63	1.30	1.24
1	H	190[A]	ASN	C-O	5.57	1.30	1.24
1	H	190[B]	ASN	C-O	5.57	1.30	1.24
1	D	189[A]	ARG	C-O	5.57	1.30	1.24
1	D	189[B]	ARG	C-O	5.57	1.30	1.24
1	B	189[A]	ARG	C-O	5.52	1.30	1.24
1	B	189[B]	ARG	C-O	5.52	1.30	1.24
1	B	279[A]	ARG	C-O	5.51	1.31	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	279[B]	ARG	C-O	5.51	1.31	1.24
1	H	424[A]	THR	C-O	5.50	1.30	1.24
1	H	424[B]	THR	C-O	5.50	1.30	1.24
1	B	190[A]	ASN	C-O	5.48	1.30	1.24
1	B	190[B]	ASN	C-O	5.48	1.30	1.24
1	B	160	HIS	CE1-NE2	5.46	1.38	1.32
1	A	217[A]	GLN	CA-C	5.29	1.57	1.52
1	A	217[B]	GLN	CA-C	5.29	1.57	1.52
1	F	279[A]	ARG	C-O	5.25	1.30	1.24
1	F	279[B]	ARG	C-O	5.25	1.30	1.24
1	C	189[A]	ARG	C-O	5.15	1.30	1.24
1	C	189[B]	ARG	C-O	5.15	1.30	1.24
1	H	189[A]	ARG	C-O	5.08	1.30	1.24
1	H	189[B]	ARG	C-O	5.08	1.30	1.24
1	A	190[A]	ASN	C-O	5.07	1.30	1.24
1	A	190[B]	ASN	C-O	5.07	1.30	1.24

All (103) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	217[A]	GLN	O-C-N	-9.08	115.90	121.71
1	A	217[B]	GLN	O-C-N	-9.08	115.90	121.71
1	H	135[A]	THR	CA-C-O	8.41	129.60	120.10
1	H	135[B]	THR	CA-C-O	8.41	129.60	120.10
1	B	189[A]	ARG	CA-C-O	7.97	128.99	120.55
1	B	189[B]	ARG	CA-C-O	7.97	128.99	120.55
1	E	83	ASP	CA-CB-CG	7.76	120.36	112.60
1	A	189[A]	ARG	CA-C-O	7.74	128.76	120.55
1	A	189[B]	ARG	CA-C-O	7.74	128.76	120.55
1	F	83	ASP	CA-CB-CG	7.55	120.15	112.60
1	H	190[A]	ASN	CA-C-O	7.41	128.41	120.55
1	H	190[B]	ASN	CA-C-O	7.41	128.41	120.55
1	C	190[A]	ASN	CA-C-O	7.38	128.37	120.55
1	C	190[B]	ASN	CA-C-O	7.38	128.37	120.55
1	H	424[A]	THR	CA-C-O	7.27	128.31	120.10
1	H	424[B]	THR	CA-C-O	7.27	128.31	120.10
1	A	217[A]	GLN	CA-C-O	7.16	127.08	120.50
1	A	217[B]	GLN	CA-C-O	7.16	127.08	120.50
1	D	190[A]	ASN	CA-C-O	7.14	128.12	120.55
1	D	190[B]	ASN	CA-C-O	7.14	128.12	120.55
1	F	190[A]	ASN	CA-C-O	7.10	128.08	120.55
1	F	190[B]	ASN	CA-C-O	7.10	128.08	120.55

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	189[A]	ARG	CA-C-O	6.88	127.84	120.55
1	F	189[B]	ARG	CA-C-O	6.88	127.84	120.55
1	B	190[A]	ASN	CA-C-O	6.75	127.70	120.55
1	B	190[B]	ASN	CA-C-O	6.75	127.70	120.55
1	A	121[A]	ASP	CA-CB-CG	6.62	119.22	112.60
1	A	121[B]	ASP	CA-CB-CG	6.62	119.22	112.60
1	C	189[A]	ARG	CA-C-O	6.41	127.34	120.55
1	C	189[B]	ARG	CA-C-O	6.41	127.34	120.55
1	D	172	ASP	CA-CB-CG	6.41	119.00	112.60
1	A	111	ASP	CA-CB-CG	6.30	118.90	112.60
1	H	189[A]	ARG	CA-C-O	6.22	127.14	120.55
1	H	189[B]	ARG	CA-C-O	6.22	127.14	120.55
1	H	503	GLU	CB-CG-CD	6.22	123.17	112.60
1	G	422	ASP	CA-CB-CG	6.20	118.80	112.60
1	D	478	PHE	CA-CB-CG	-6.19	107.61	113.80
1	D	83	ASP	CA-CB-CG	6.17	118.77	112.60
1	A	513[A]	ASN	CA-C-O	6.16	128.99	121.84
1	A	513[B]	ASN	CA-C-O	6.16	128.99	121.84
1	C	515	THR	CA-CB-OG1	-6.13	100.40	109.60
1	D	189[A]	ARG	CA-C-O	6.11	127.02	120.55
1	D	189[B]	ARG	CA-C-O	6.11	127.02	120.55
1	E	424[A]	THR	CA-C-O	6.10	126.99	120.10
1	E	424[B]	THR	CA-C-O	6.10	126.99	120.10
1	B	276	GLU	CB-CA-C	-6.10	103.32	110.76
1	C	83	ASP	CA-CB-CG	6.08	118.68	112.60
1	H	407	PHE	CA-CB-CG	-6.01	107.79	113.80
1	E	217	GLN	CB-CA-C	-6.00	100.67	109.97
1	A	167	ASP	CA-CB-CG	5.99	118.59	112.60
1	H	428	ASP	CA-CB-CG	5.98	118.58	112.60
1	B	127	GLU	CB-CG-CD	5.88	122.60	112.60
1	G	294	PHE	CA-CB-CG	5.88	119.68	113.80
1	E	121	ASP	CA-CB-CG	5.85	118.45	112.60
1	B	83	ASP	CA-CB-CG	5.83	118.43	112.60
1	A	83	ASP	CA-CB-CG	5.83	118.43	112.60
1	A	190[A]	ASN	CA-C-O	5.83	126.73	120.55
1	A	190[B]	ASN	CA-C-O	5.83	126.73	120.55
1	E	338	GLN	CB-CG-CD	5.78	122.43	112.60
1	E	424[A]	THR	N-CA-C	5.78	118.36	111.71
1	E	424[B]	THR	N-CA-C	5.78	118.36	111.71
1	H	103	HIS	CA-CB-CG	5.71	119.51	113.80
1	G	428	ASP	CA-CB-CG	5.71	118.31	112.60
1	A	378	GLN	OE1-CD-NE2	5.70	128.30	122.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	83	ASP	CA-CB-CG	5.68	118.28	112.60
1	A	422	ASP	CA-CB-CG	5.66	118.25	112.60
1	E	136	ASP	CA-CB-CG	5.66	118.25	112.60
1	E	127	GLU	N-CA-CB	5.65	119.04	110.28
1	A	151	LEU	N-CA-CB	-5.64	101.80	110.20
1	H	71	GLU	CB-CA-C	-5.59	102.07	110.90
1	F	376	ASP	CA-CB-CG	5.57	118.17	112.60
1	E	71	GLU	N-CA-CB	5.56	118.07	110.01
1	A	376	ASP	CA-CB-CG	5.53	118.12	112.60
1	F	217	GLN	CB-CA-C	-5.51	101.42	109.97
1	F	279[A]	ARG	CA-C-O	5.45	126.08	119.11
1	F	279[B]	ARG	CA-C-O	5.45	126.08	119.11
1	H	422	ASP	CA-CB-CG	5.42	118.02	112.60
1	E	127	GLU	CB-CG-CD	5.41	121.80	112.60
1	B	279[A]	ARG	CA-C-O	5.40	126.02	119.11
1	B	279[B]	ARG	CA-C-O	5.40	126.02	119.11
1	D	428	ASP	CA-CB-CG	5.39	117.99	112.60
1	E	342	PRO	N-CA-CB	-5.36	98.26	102.28
1	B	376	ASP	CA-CB-CG	5.33	117.93	112.60
1	D	465	GLU	N-CA-CB	5.29	118.49	110.28
1	B	279[A]	ARG	N-CA-C	5.29	119.58	113.18
1	B	279[B]	ARG	N-CA-C	5.29	119.58	113.18
1	B	478	PHE	CA-CB-CG	-5.25	108.55	113.80
1	H	83	ASP	CA-CB-CG	5.25	117.85	112.60
1	E	71	GLU	CB-CG-CD	5.23	121.48	112.60
1	H	71	GLU	N-CA-CB	5.21	117.62	110.07
1	C	103	HIS	CA-CB-CG	5.12	118.92	113.80
1	D	201	LYS	CB-CG-CD	5.12	123.06	111.30
1	E	294	PHE	CA-CB-CG	5.11	118.91	113.80
1	B	428	ASP	CA-CB-CG	5.11	117.71	112.60
1	B	189[A]	ARG	O-C-N	-5.09	116.73	122.12
1	B	189[B]	ARG	O-C-N	-5.09	116.73	122.12
1	B	358	ASP	CA-CB-CG	5.09	117.69	112.60
1	B	136	ASP	CA-CB-CG	5.08	117.68	112.60
1	C	172	ASP	CA-CB-CG	5.08	117.68	112.60
1	F	136	ASP	CA-CB-CG	5.08	117.68	112.60
1	C	340	LYS	CB-CG-CD	-5.06	99.67	111.30
1	E	376	ASP	CA-CB-CG	5.06	117.66	112.60
1	G	179	ASP	CA-CB-CG	5.00	117.60	112.60

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	3642	0	3687	62	0
1	B	3632	0	3669	29	0
1	C	3584	0	3624	39	0
1	D	3584	0	3619	40	0
1	E	3616	0	3657	32	0
1	F	3651	0	3693	39	0
1	G	3583	0	3619	29	0
1	H	3586	0	3630	39	0
2	A	20	0	0	2	0
2	B	25	0	0	1	0
2	C	15	0	0	1	0
2	D	35	0	0	1	0
2	E	15	0	0	0	0
2	F	35	0	0	0	0
2	G	30	0	0	0	0
2	H	20	0	0	1	0
3	A	12	0	15	5	0
3	B	18	0	24	4	0
3	C	24	0	32	4	0
3	D	12	0	16	0	0
3	E	18	0	24	9	0
3	F	18	0	24	4	0
3	G	12	0	16	1	0
3	H	18	0	24	5	0
4	C	2	0	0	0	0
4	D	1	0	0	0	0
4	E	2	0	0	0	0
4	G	1	0	0	0	0
5	C	1	0	0	0	0
6	A	542	0	0	9	0
6	B	554	0	0	12	0
6	C	590	0	0	15	0
6	D	562	0	0	11	0
6	E	554	0	0	12	0
6	F	565	0	0	11	0
6	G	571	0	0	10	0
6	H	556	0	0	11	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	33706	0	29373	264	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 (264) 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:92:MET:HE1	6:B:822:HOH:O	1.50	1.11
1:F:92:MET:HE1	6:F:735:HOH:O	1.58	1.02
1:H:92:MET:HE1	6:H:728:HOH:O	1.59	1.01
1:E:270[B]:ASN:OD1	6:E:701:HOH:O	1.80	0.99
1:A:205[B]:LEU:HD23	1:A:221:LEU:HD22	1.44	0.96
1:C:484:GLU:HG2	6:C:1187:HOH:O	1.69	0.90
1:B:135[B]:THR:HG22	6:B:731:HOH:O	1.71	0.90
1:A:483[A]:VAL:HG12	2:A:604:SO4:O2	1.70	0.89
1:F:165:ARG:H	3:F:610:GOL:H12	1.36	0.89
1:F:205[A]:LEU:HD23	1:F:221[A]:LEU:HD22	1.57	0.87
1:E:205[A]:LEU:HD23	1:E:221[A]:LEU:HD22	1.54	0.87
1:B:364[B]:THR:HG22	1:D:364[B]:THR:HG22	1.57	0.87
1:A:267[B]:MET:HE2	6:A:1160:HOH:O	1.74	0.86
1:B:364[B]:THR:HG22	1:D:364[B]:THR:CG2	2.06	0.84
1:E:364[B]:THR:HG22	1:G:364[B]:THR:HG22	1.57	0.84
1:A:364[B]:THR:HG22	1:C:364[B]:THR:CG2	2.07	0.84
1:A:92:MET:HE3	1:A:175:LEU:HD13	1.59	0.84
1:A:164:SER:OG	3:A:606:GOL:H11	1.79	0.82
1:A:488[A]:ASN:ND2	6:A:701:HOH:O	2.13	0.81
1:E:364[B]:THR:CG2	1:G:364[B]:THR:HG22	2.12	0.80
1:E:364[B]:THR:HG22	1:G:364[B]:THR:CG2	2.11	0.80
1:A:364[B]:THR:HG22	1:C:364[B]:THR:HG22	1.63	0.79
1:B:364[B]:THR:CG2	1:D:364[B]:THR:HG22	2.12	0.79
1:A:364[B]:THR:CG2	1:C:364[B]:THR:HG22	2.14	0.78
1:D:411:ARG:HG3	1:D:411:ARG:HH11	1.49	0.77
1:A:451:CYS:HG	1:A:456:CYS:HG	0.80	0.77
1:F:270[B]:ASN:OD1	6:F:702:HOH:O	2.03	0.76
1:H:205[A]:LEU:HD23	1:H:221[A]:LEU:HD22	1.68	0.74
1:A:424[A]:THR:HG22	6:A:812:HOH:O	1.86	0.74
1:C:432[A]:LYS:HE3	6:D:1097:HOH:O	1.87	0.74
1:E:190[B]:ASN:ND2	6:E:702:HOH:O	1.96	0.73
1:B:197[A]:GLU:OE2	6:B:701:HOH:O	2.06	0.73
1:F:76:PHE:CG	1:G:397[B]:VAL:HG21	2.24	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:453:SER:HG	1:A:455:GLY:N	1.88	0.71
1:H:406[B]:THR:HG22	6:H:863:HOH:O	1.89	0.71
3:E:607:GOL:H11	1:G:339:LYS:HZ1	1.55	0.71
1:D:205:LEU:HD23	1:D:221[A]:LEU:HD22	1.73	0.70
1:F:364[B]:THR:HG22	1:H:364[B]:THR:CG2	2.22	0.69
1:E:190[B]:ASN:ND2	6:E:704:HOH:O	2.24	0.69
1:A:103:HIS:ND1	1:A:267[B]:MET:HE1	2.08	0.68
1:F:364[B]:THR:CG2	1:H:364[B]:THR:HG22	2.24	0.68
1:F:381:LYS:NZ	6:F:705:HOH:O	2.24	0.68
1:A:400[B]:GLU:CD	1:D:69:VAL:HG13	2.19	0.67
1:A:103:HIS:HE1	6:A:1095:HOH:O	1.78	0.67
1:H:88:LYS:HE2	1:H:127:GLU:OE2	1.94	0.67
1:D:117:ARG:NH1	6:D:702:HOH:O	2.22	0.67
1:F:364[B]:THR:HG22	1:H:364[B]:THR:HG22	1.77	0.67
1:H:121:ASP:OD1	1:H:124:ARG:NH2	2.24	0.67
1:A:400[B]:GLU:OE1	1:D:69:VAL:HA	1.95	0.66
1:B:189[B]:ARG:HD2	6:B:987:HOH:O	1.95	0.66
1:D:197[A]:GLU:OE2	6:D:701:HOH:O	2.14	0.66
1:G:205:LEU:HD23	1:G:221[A]:LEU:HD22	1.77	0.66
1:H:474:GLU:HG3	6:H:1063:HOH:O	1.95	0.65
1:B:381:LYS:NZ	6:B:707:HOH:O	2.30	0.65
1:E:164:SER:OG	3:E:608:GOL:H11	1.95	0.65
1:F:437[A]:PHE:CD2	1:F:438:ARG:HD3	2.32	0.65
1:D:201:LYS:HB2	6:D:1222:HOH:O	1.96	0.65
1:F:437[A]:PHE:HD2	1:F:438:ARG:HD3	1.62	0.65
1:C:92:MET:HE1	6:C:734:HOH:O	1.98	0.64
1:F:270[A]:ASN:ND2	6:F:708:HOH:O	2.29	0.64
1:E:165:ARG:H	3:E:608:GOL:H32	1.63	0.63
1:B:124:ARG:NH2	6:B:702:HOH:O	2.24	0.63
1:A:364[B]:THR:HG22	1:C:364[B]:THR:HG21	1.80	0.62
1:H:424[A]:THR:HG22	6:H:840:HOH:O	1.99	0.62
1:C:180:ALA:HA	3:C:608:GOL:H2	1.81	0.62
3:E:607:GOL:H11	1:G:339:LYS:NZ	2.13	0.62
1:G:397[B]:VAL:HG23	6:G:1037:HOH:O	2.00	0.62
1:A:164:SER:HA	3:A:606:GOL:H32	1.81	0.61
1:E:154:GLU:HG3	6:E:766:HOH:O	2.00	0.61
1:A:420:HIS:CD2	1:A:448:VAL:HG13	2.36	0.60
1:H:511:LYS:HE2	6:H:702:HOH:O	1.99	0.60
1:G:381:LYS:NZ	6:G:703:HOH:O	2.25	0.60
1:H:381:LYS:NZ	6:H:708:HOH:O	2.34	0.60
6:B:1075:HOH:O	1:C:160:HIS:HE1	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:503[B]:GLU:HG3	6:G:863:HOH:O	2.02	0.60
3:E:607:GOL:H32	1:F:166:ASN:HD21	1.67	0.59
1:A:483[A]:VAL:CG1	2:A:604:SO4:O2	2.49	0.59
1:B:117:ARG:HG3	6:B:969:HOH:O	2.02	0.59
1:G:381:LYS:NZ	6:G:711:HOH:O	2.36	0.59
1:A:364[B]:THR:CG2	1:C:364[B]:THR:CG2	2.75	0.59
1:D:121:ASP:OD1	1:D:124:ARG:NH2	2.33	0.58
1:D:406:THR:HG23	6:D:878:HOH:O	2.02	0.58
1:A:166:ASN:HD21	3:B:607:GOL:C3	2.17	0.58
1:A:456:CYS:SG	6:A:1048:HOH:O	2.57	0.58
1:B:278:MET:HA	3:B:606:GOL:H11	1.86	0.58
1:D:88:LYS:HE2	1:D:127[A]:GLU:OE2	2.04	0.58
1:E:160[B]:HIS:HE2	3:E:608:GOL:H31	1.67	0.58
1:C:121:ASP:OD1	1:C:124:ARG:NH2	2.35	0.57
1:B:364[B]:THR:HG22	1:D:364[B]:THR:HG21	1.84	0.57
1:E:474:GLU:HG3	6:E:941:HOH:O	2.03	0.57
1:E:364[B]:THR:CG2	1:G:364[B]:THR:CG2	2.79	0.57
1:E:424[A]:THR:HG22	6:E:791:HOH:O	2.04	0.57
1:B:364[B]:THR:CG2	1:D:364[B]:THR:CG2	2.78	0.56
1:F:270[A]:ASN:ND2	6:F:710:HOH:O	2.37	0.56
1:A:400[B]:GLU:OE2	1:D:73:VAL:HG23	2.05	0.56
1:C:488[A]:ASN:ND2	6:C:717:HOH:O	2.37	0.56
1:G:409:GLU:HG3	6:G:856:HOH:O	2.06	0.55
1:A:447:LEU:HD13	1:A:469:LEU:HD12	1.87	0.55
1:B:121:ASP:HB2	6:B:702:HOH:O	2.06	0.55
1:H:502:ALA:HB1	3:H:607:GOL:C1	2.37	0.55
1:A:381:LYS:NZ	6:A:709:HOH:O	2.41	0.54
1:E:424[A]:THR:HG23	6:E:848:HOH:O	2.07	0.54
1:A:166:ASN:HD21	3:B:607:GOL:H31	1.72	0.54
1:E:160[B]:HIS:HB3	1:F:437[B]:PHE:CZ	2.42	0.54
1:C:180:ALA:CA	3:C:608:GOL:H2	2.37	0.53
1:F:205[A]:LEU:HD23	1:F:221[A]:LEU:CD2	2.34	0.53
1:E:221[B]:LEU:HD11	1:E:407:PHE:CE1	2.43	0.53
1:C:164:SER:OG	3:C:609:GOL:H32	2.09	0.53
1:D:201:LYS:O	1:D:201:LYS:HG2	2.08	0.53
1:G:462:SER:OG	1:G:465:GLU:HG3	2.08	0.53
1:A:451:CYS:HA	1:A:456:CYS:SG	2.49	0.52
1:F:237:GLY:HA2	3:F:609:GOL:H31	1.90	0.52
1:F:448:VAL:HG11	1:H:337:PRO:HD2	1.90	0.52
1:A:451:CYS:CB	1:A:456:CYS:HG	2.23	0.52
1:C:437:PHE:CZ	1:D:160:HIS:HB3	2.45	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:312:TRP:CE2	1:A:323[B]:MET:HE1	2.44	0.52
1:A:424[A]:THR:HG23	6:A:776:HOH:O	2.10	0.52
1:C:377:PHE:CE2	6:C:883:HOH:O	2.54	0.52
1:E:217:GLN:HE22	1:E:424[A]:THR:HG21	1.74	0.52
3:C:607:GOL:H32	6:C:788:HOH:O	2.10	0.52
1:A:103:HIS:CE1	1:A:267[B]:MET:HE1	2.45	0.51
1:A:437:PHE:CE2	1:B:160:HIS:HB3	2.45	0.51
1:E:221[B]:LEU:HD11	1:E:407:PHE:CZ	2.45	0.51
1:F:189[A]:ARG:HH12	3:F:609:GOL:C1	2.23	0.51
1:A:121[A]:ASP:OD1	1:A:124:ARG:NH2	2.40	0.50
1:E:165:ARG:H	3:E:608:GOL:C3	2.24	0.50
1:A:347[B]:LEU:HD12	1:D:77:THR:HG22	1.94	0.50
1:H:190[B]:ASN:ND2	6:H:723:HOH:O	2.44	0.50
1:H:502:ALA:HB1	3:H:607:GOL:H12	1.93	0.50
1:A:185:ILE:O	1:A:189[A]:ARG:HG2	2.12	0.50
1:D:185:ILE:O	1:D:189[A]:ARG:HG2	2.13	0.49
1:A:246[A]:LEU:HD12	1:A:290:PHE:HB2	1.94	0.49
1:E:513:ASN:HB3	6:E:1042:HOH:O	2.10	0.49
1:B:438[A]:ARG:NE	6:B:717:HOH:O	2.46	0.49
1:A:108:THR:HG21	3:A:605:GOL:C3	2.42	0.49
1:F:165:ARG:N	3:F:610:GOL:H12	2.17	0.49
1:B:185:ILE:O	1:B:189[A]:ARG:HG2	2.12	0.49
1:D:334:SER:OG	2:D:602:SO4:O4	2.30	0.49
1:H:165:ARG:NH2	6:H:713:HOH:O	2.37	0.49
1:F:364[B]:THR:HG22	1:H:364[B]:THR:HG21	1.94	0.49
1:A:217[A]:GLN:OE1	1:A:424[A]:THR:HG21	2.13	0.48
1:D:266:PHE:HB2	1:H:266:PHE:HB3	1.95	0.48
1:F:424[B]:THR:HG21	6:F:1103:HOH:O	2.13	0.48
1:B:448:VAL:HG11	1:D:337:PRO:HD2	1.95	0.48
1:H:197[A]:GLU:CD	6:H:704:HOH:O	2.57	0.48
1:C:74:GLU:HG2	6:C:1126:HOH:O	2.13	0.48
1:G:318[B]:GLU:HG2	6:G:769:HOH:O	2.14	0.48
1:G:221[A]:LEU:HD21	1:G:407:PHE:CE2	2.49	0.47
3:E:606:GOL:H11	1:F:438:ARG:HB2	1.95	0.47
1:A:410[A]:ASP:OD2	1:A:414:LYS:HE2	2.14	0.47
1:E:364[B]:THR:HG21	1:G:364[B]:THR:HG22	1.96	0.47
1:H:484:GLU:HG3	6:H:1022:HOH:O	2.13	0.47
1:C:511:LYS:HE2	6:C:703:HOH:O	2.13	0.47
1:F:499:ASN:HB2	6:F:944:HOH:O	2.14	0.47
1:D:189[A]:ARG:NH2	6:D:718:HOH:O	2.45	0.47
1:G:318[B]:GLU:OE1	6:G:701:HOH:O	2.21	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:221[A]:LEU:HD21	1:D:407:PHE:CE2	2.50	0.47
1:H:165:ARG:H	3:H:605:GOL:HO3	1.61	0.46
1:C:185:ILE:O	1:C:189[A]:ARG:HG2	2.14	0.46
1:F:488:ASN:HA	6:F:701:HOH:O	2.15	0.46
1:B:347[B]:LEU:HD12	1:C:77:THR:HG22	1.96	0.46
1:H:95:LYS:NZ	6:H:718:HOH:O	2.42	0.46
1:A:112:LYS:HB3	1:A:112:LYS:HE3	1.40	0.46
1:C:201:LYS:NZ	6:C:731:HOH:O	2.49	0.46
1:A:154:GLU:N	1:A:155:PRO:CD	2.79	0.46
1:B:416:LEU:N	1:B:417:PRO:CD	2.79	0.45
1:C:437:PHE:CE2	1:D:160:HIS:HB3	2.50	0.45
1:H:334:SER:OG	2:H:601:SO4:O2	2.30	0.45
1:A:205[B]:LEU:CD2	1:A:221:LEU:HD22	2.31	0.45
1:A:157:LYS:HD3	1:B:438[B]:ARG:HG3	1.98	0.45
1:D:484:GLU:HG2	6:D:1134:HOH:O	2.17	0.45
1:B:165:ARG:HG2	2:B:603:SO4:O2	2.16	0.45
1:D:432:LYS:NZ	6:D:724:HOH:O	2.50	0.45
1:F:364[B]:THR:CG2	1:H:364[B]:THR:CG2	2.88	0.45
1:F:364[B]:THR:HG21	1:H:364[B]:THR:HG22	1.99	0.45
1:C:135:THR:HG22	6:C:1126:HOH:O	2.16	0.45
1:D:80:ILE:HD11	1:D:133:TRP:CE3	2.52	0.45
1:F:197[A]:GLU:OE1	6:F:703:HOH:O	2.21	0.45
1:F:121[A]:ASP:OD1	1:F:124:ARG:NH2	2.43	0.44
1:F:438:ARG:HG2	1:F:438:ARG:HH21	1.81	0.44
1:H:502:ALA:CB	3:H:607:GOL:H11	2.47	0.44
1:A:77:THR:HG22	1:D:347[B]:LEU:HD12	1.99	0.44
1:E:381:LYS:NZ	6:E:724:HOH:O	2.50	0.44
1:B:80:ILE:HD11	1:B:133:TRP:CE3	2.53	0.44
1:A:190[B]:ASN:ND2	6:A:716:HOH:O	2.48	0.44
1:E:469:LEU:HD23	6:E:960:HOH:O	2.17	0.44
1:A:217[A]:GLN:HE21	1:B:256:LEU:HD23	1.82	0.44
1:A:364[B]:THR:HG21	1:C:364[B]:THR:HG22	1.95	0.44
1:B:397:VAL:HG21	1:C:76:PHE:CG	2.53	0.44
1:E:221[B]:LEU:HD21	1:E:225:LEU:HD11	1.99	0.44
1:G:451:CYS:HB3	1:G:456:CYS:O	2.18	0.44
1:D:246:LEU:HD12	1:D:246:LEU:C	2.43	0.43
1:A:199:ALA:HB1	1:A:222:PRO:HB3	2.00	0.43
1:C:276:GLU:HB2	1:C:277:PRO:HD2	1.99	0.43
1:E:364[B]:THR:HG22	1:G:364[B]:THR:HG21	1.93	0.43
1:C:80:ILE:HD11	1:C:133:TRP:CE3	2.53	0.43
1:G:80:ILE:HD11	1:G:133:TRP:CE3	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:76:PHE:CG	1:D:397:VAL:HG21	2.54	0.43
1:H:205[A]:LEU:HD23	1:H:221[A]:LEU:CD2	2.42	0.43
1:A:427:ALA:HB2	1:C:335:ILE:HD11	1.99	0.43
1:G:432:LYS:HE3	6:G:1161:HOH:O	2.18	0.43
1:B:511:LYS:HE2	6:B:709:HOH:O	2.18	0.43
1:C:484:GLU:CD	1:C:484:GLU:H	2.27	0.43
1:G:190[A]:ASN:ND2	6:G:708:HOH:O	2.48	0.43
1:A:400[B]:GLU:OE2	1:D:69:VAL:HG13	2.19	0.43
1:A:416:LEU:HD13	1:A:483[B]:VAL:HG21	2.00	0.43
1:B:183:THR:HG23	6:B:1038:HOH:O	2.17	0.43
1:C:276:GLU:HB2	1:C:277:PRO:CD	2.49	0.43
1:E:246:LEU:C	1:E:246:LEU:HD12	2.43	0.43
1:A:80:ILE:HD11	1:A:133:TRP:CE3	2.54	0.43
1:A:177:CYS:HB2	1:A:246[A]:LEU:HD11	2.01	0.43
1:D:190[A]:ASN:ND2	6:D:729:HOH:O	2.52	0.43
1:F:189[B]:ARG:HH22	1:F:511:LYS:HG2	1.84	0.43
1:D:411:ARG:HG3	1:D:411:ARG:NH1	2.25	0.42
1:C:72:LYS:NZ	6:C:732:HOH:O	2.49	0.42
1:E:77:THR:HG22	1:H:347[B]:LEU:HD12	2.01	0.42
1:F:189[B]:ARG:NH2	6:F:728:HOH:O	2.53	0.42
1:H:276:GLU:HB2	1:H:277:PRO:HD2	2.00	0.42
1:E:397:VAL:HG21	1:H:76:PHE:CG	2.54	0.42
1:G:221[A]:LEU:HD21	1:G:407:PHE:CZ	2.54	0.42
1:C:377:PHE:CZ	6:C:883:HOH:O	2.73	0.42
1:D:88:LYS:HE3	6:D:1114:HOH:O	2.19	0.42
1:A:108:THR:CG2	3:A:605:GOL:H32	2.50	0.42
6:F:1020:HOH:O	1:G:397[B]:VAL:HG22	2.19	0.42
1:F:246:LEU:C	1:F:246:LEU:HD12	2.44	0.41
1:A:217[A]:GLN:NE2	1:B:256:LEU:HD23	2.34	0.41
1:E:414:LYS:HD2	6:E:710:HOH:O	2.20	0.41
1:E:201:LYS:NZ	6:E:726:HOH:O	2.52	0.41
1:F:314:LEU:HD21	1:H:314:LEU:HD21	2.03	0.41
1:C:332:GLY:O	6:C:702:HOH:O	2.22	0.41
1:H:112:LYS:HB3	1:H:112:LYS:HE3	1.83	0.41
1:A:189[B]:ARG:HD2	6:A:977:HOH:O	2.20	0.41
1:C:90:ASP:OD1	1:C:172:ASP:OD2	2.39	0.41
1:C:189[A]:ARG:NH2	6:C:735:HOH:O	2.53	0.41
1:D:90:ASP:OD1	1:D:172:ASP:OD2	2.38	0.41
3:E:607:GOL:H32	1:F:166:ASN:ND2	2.35	0.41
1:G:90:ASP:OD1	1:G:172:ASP:OD2	2.39	0.41
1:H:154:GLU:N	1:H:155:PRO:CD	2.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:603:SO4:O2	6:C:701:HOH:O	2.21	0.41
1:A:165:ARG:HG2	3:A:606:GOL:O3	2.20	0.41
1:A:166:ASN:ND2	3:B:607:GOL:H31	2.36	0.41
1:F:437[B]:PHE:CE1	1:F:438:ARG:HD3	2.56	0.41
1:G:318[B]:GLU:CD	6:G:809:HOH:O	2.62	0.41
1:A:123:GLU:O	1:A:127:GLU:HG2	2.20	0.41
1:C:381:LYS:NZ	6:C:712:HOH:O	2.35	0.41
1:E:429:TYR:CZ	1:E:433:LYS:HE3	2.55	0.41
1:C:216:ALA:O	1:C:217[A]:GLN:HG3	2.21	0.41
1:C:213:LEU:HD12	1:C:213:LEU:HA	1.97	0.40
1:C:221:LEU:HD13	1:C:221:LEU:C	2.46	0.40
1:D:264:ASP:HB3	1:D:267:MET:HE3	2.03	0.40
1:G:189[B]:ARG:NE	3:G:609:GOL:O2	2.55	0.40
1:D:468:LYS:HD2	6:D:754:HOH:O	2.21	0.40
1:H:221[A]:LEU:HD21	1:H:407:PHE:CZ	2.57	0.40
1:G:112:LYS:HE3	1:G:112:LYS:HB3	1.92	0.40
1:H:142:MET:HE3	1:H:142:MET:HB2	1.89	0.40
1:H:221[A]:LEU:HD21	1:H:407:PHE:CE2	2.57	0.40
1:F:185:ILE:O	1:F:189[A]:ARG:HG2	2.21	0.40
1:H:502:ALA:HB1	3:H:607:GOL:H11	2.04	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	469/465 (101%)	460 (98%)	7 (2%)	2 (0%)	30	13
1	B	466/465 (100%)	459 (98%)	6 (1%)	1 (0%)	43	24
1	C	461/465 (99%)	453 (98%)	7 (2%)	1 (0%)	43	24
1	D	461/465 (99%)	452 (98%)	7 (2%)	2 (0%)	30	13

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	466/465 (100%)	458 (98%)	6 (1%)	2 (0%)	30	13
1	F	470/465 (101%)	463 (98%)	6 (1%)	1 (0%)	43	24
1	G	461/465 (99%)	454 (98%)	6 (1%)	1 (0%)	43	24
1	H	462/465 (99%)	456 (99%)	5 (1%)	1 (0%)	43	24
All	All	3716/3720 (100%)	3655 (98%)	50 (1%)	11 (0%)	43	18

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	256[A]	LEU
1	A	256[B]	LEU
1	B	256	LEU
1	C	256	LEU
1	D	256[A]	LEU
1	D	256[B]	LEU
1	E	256[A]	LEU
1	E	256[B]	LEU
1	F	256	LEU
1	G	256	LEU
1	H	256	LEU

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	413/402 (103%)	407 (98%)	6 (2%)	57	31
1	B	407/402 (101%)	406 (100%)	1 (0%)	87	79
1	C	403/402 (100%)	400 (99%)	3 (1%)	76	59
1	D	403/402 (100%)	403 (100%)	0	100	100
1	E	408/402 (102%)	404 (99%)	4 (1%)	68	47
1	F	411/402 (102%)	409 (100%)	2 (0%)	81	68
1	G	403/402 (100%)	401 (100%)	2 (0%)	81	68

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	H	404/402 (100%)	400 (99%)	4 (1%)	68 47
All	All	3252/3216 (101%)	3230 (99%)	22 (1%)	78 59

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	112	LYS
1	A	246[A]	LEU
1	A	246[B]	LEU
1	A	333	SER
1	A	437	PHE
1	A	451	CYS
1	B	160	HIS
1	C	69	VAL
1	C	116	LEU
1	C	437	PHE
1	E	68	SER
1	E	424[A]	THR
1	E	424[B]	THR
1	E	437	PHE
1	F	160	HIS
1	F	201	LYS
1	G	127	GLU
1	G	409	GLU
1	H	69	VAL
1	H	246	LEU
1	H	424[A]	THR
1	H	424[B]	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	166	ASN
1	A	270	ASN
1	A	403	GLN
1	A	460	ASN
1	B	104	GLN
1	B	202	ASN
1	B	338	GLN
1	B	403	GLN
1	B	459	GLN

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Mol	Chain	Res	Type
1	C	298	ASN
1	D	129	ASN
1	D	202	ASN
1	D	403	GLN
1	D	460	ASN
1	E	214	GLN
1	E	298	ASN
1	E	403	GLN
1	F	166	ASN
1	F	202	ASN
1	F	298	ASN
1	G	298	ASN
1	G	459	GLN
1	G	460	ASN
1	G	513	ASN
1	H	231	GLN
1	H	298	ASN
1	H	338	GLN
1	H	403	GLN
1	H	459	GLN

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 68 ligands modelled in this entry, 7 are 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	$\# Z > 2$
2	SO4	F	606	-	4,4,4	0.26	0	6,6,6	0.48	0
2	SO4	H	602	-	4,4,4	0.30	0	6,6,6	0.42	0
2	SO4	H	601	-	4,4,4	0.39	0	6,6,6	0.10	0
2	SO4	F	607	-	4,4,4	0.36	0	6,6,6	0.29	0
2	SO4	C	605	-	4,4,4	0.35	0	6,6,6	0.10	0
3	GOL	C	609	-	5,5,5	0.09	0	5,5,5	0.33	0
2	SO4	E	603	-	4,4,4	0.51	0	6,6,6	0.25	0
2	SO4	B	605	-	4,4,4	0.39	0	6,6,6	0.16	0
2	SO4	F	603	-	4,4,4	0.32	0	6,6,6	0.40	0
2	SO4	E	604	-	4,4,4	0.20	0	6,6,6	0.14	0
3	GOL	E	606	-	5,5,5	0.50	0	5,5,5	0.90	0
3	GOL	E	608	-	5,5,5	0.83	0	5,5,5	0.53	0
3	GOL	F	609	-	5,5,5	0.08	0	5,5,5	0.31	0
2	SO4	D	602	-	4,4,4	0.24	0	6,6,6	0.11	0
3	GOL	H	605	-	5,5,5	0.52	0	5,5,5	0.69	0
3	GOL	C	607	-	5,5,5	0.38	0	5,5,5	0.84	0
3	GOL	G	609	-	5,5,5	0.16	0	5,5,5	0.51	0
2	SO4	D	606	-	4,4,4	0.21	0	6,6,6	0.37	0
2	SO4	B	601	-	4,4,4	0.50	0	6,6,6	0.22	0
3	GOL	A	606	-	5,5,5	1.11	0	5,5,5	0.43	0
3	GOL	C	608	-	5,5,5	0.47	0	5,5,5	1.00	0
2	SO4	H	603	-	4,4,4	0.38	0	6,6,6	0.18	0
3	GOL	D	609	-	5,5,5	0.11	0	5,5,5	0.35	0
2	SO4	E	605	-	4,4,4	0.33	0	6,6,6	0.15	0
2	SO4	G	605	-	4,4,4	0.50	0	6,6,6	0.38	0
2	SO4	D	607	-	4,4,4	0.37	0	6,6,6	0.09	0
2	SO4	B	604	-	4,4,4	0.40	0	6,6,6	0.33	0
2	SO4	G	607	-	4,4,4	0.32	0	6,6,6	0.19	0
2	SO4	A	603	-	4,4,4	0.36	0	6,6,6	0.20	0
3	GOL	B	607	-	5,5,5	0.18	0	5,5,5	0.59	0
3	GOL	F	610	-	5,5,5	0.82	0	5,5,5	0.60	0
2	SO4	D	603	-	4,4,4	0.45	0	6,6,6	0.12	0
2	SO4	C	604	-	4,4,4	0.68	0	6,6,6	0.27	0
2	SO4	D	604	-	4,4,4	0.38	0	6,6,6	0.11	0
2	SO4	A	602	-	4,4,4	0.33	0	6,6,6	0.68	0
2	SO4	F	605	-	4,4,4	0.39	0	6,6,6	0.10	0
2	SO4	H	604	-	4,4,4	0.39	0	6,6,6	0.08	0
3	GOL	D	610	-	5,5,5	0.20	0	5,5,5	0.54	0
2	SO4	B	602	-	4,4,4	0.27	0	6,6,6	0.41	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GOL	E	607	-	5,5,5	0.49	0	5,5,5	0.66	0
2	SO4	G	604	-	4,4,4	0.43	0	6,6,6	0.14	0
3	GOL	B	606	-	5,5,5	0.06	0	5,5,5	0.29	0
3	GOL	H	606	-	5,5,5	0.18	0	5,5,5	0.66	0
3	GOL	G	608	-	5,5,5	0.11	0	5,5,5	0.31	0
2	SO4	F	601	-	4,4,4	0.21	0	6,6,6	0.29	0
2	SO4	C	603	-	4,4,4	0.39	0	6,6,6	0.21	0
2	SO4	G	602	-	4,4,4	0.25	0	6,6,6	0.24	0
2	SO4	G	606	-	4,4,4	0.39	0	6,6,6	0.12	0
3	GOL	C	606	-	5,5,5	0.14	0	5,5,5	0.46	0
2	SO4	D	605	-	4,4,4	0.27	0	6,6,6	0.64	0
3	GOL	H	607	-	5,5,5	0.27	0	5,5,5	0.89	0
2	SO4	B	603	-	4,4,4	0.37	0	6,6,6	0.22	0
3	GOL	A	605	-	5,5,5	0.20	0	5,5,5	0.63	0
2	SO4	A	601	-	4,4,4	0.36	0	6,6,6	0.33	0
2	SO4	D	608	-	4,4,4	0.26	0	6,6,6	0.04	0
2	SO4	F	604	-	4,4,4	0.60	0	6,6,6	0.42	0
2	SO4	F	602	-	4,4,4	0.39	0	6,6,6	0.07	0
3	GOL	B	608	-	5,5,5	0.47	0	5,5,5	0.59	0
2	SO4	G	603	-	4,4,4	0.40	0	6,6,6	0.06	0
2	SO4	A	604	-	4,4,4	0.31	0	6,6,6	0.25	0
3	GOL	F	608	-	5,5,5	0.27	0	5,5,5	0.29	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	C	609	-	-	4/4/4/4	-
3	GOL	E	606	-	-	2/4/4/4	-
3	GOL	E	608	-	-	3/4/4/4	-
3	GOL	F	609	-	-	4/4/4/4	-
3	GOL	H	605	-	-	3/4/4/4	-
3	GOL	G	609	-	-	0/4/4/4	-
3	GOL	C	607	-	-	3/4/4/4	-
3	GOL	C	608	-	-	1/4/4/4	-
3	GOL	A	606	-	-	2/4/4/4	-
3	GOL	D	609	-	-	3/4/4/4	-
3	GOL	F	610	-	-	4/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	B	607	-	-	2/4/4/4	-
3	GOL	D	610	-	-	4/4/4/4	-
3	GOL	E	607	-	-	0/4/4/4	-
3	GOL	B	606	-	-	1/4/4/4	-
3	GOL	H	606	-	-	0/4/4/4	-
3	GOL	G	608	-	-	2/4/4/4	-
3	GOL	C	606	-	-	4/4/4/4	-
3	GOL	H	607	-	-	2/4/4/4	-
3	GOL	A	605	-	-	0/4/4/4	-
3	GOL	B	608	-	-	4/4/4/4	-
3	GOL	F	608	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (50) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	606	GOL	O1-C1-C2-O2
3	A	606	GOL	O1-C1-C2-C3
3	B	607	GOL	O1-C1-C2-C3
3	C	606	GOL	O1-C1-C2-C3
3	C	606	GOL	C1-C2-C3-O3
3	C	607	GOL	O1-C1-C2-C3
3	C	609	GOL	O1-C1-C2-C3
3	C	609	GOL	C1-C2-C3-O3
3	D	610	GOL	O1-C1-C2-C3
3	D	610	GOL	C1-C2-C3-O3
3	D	610	GOL	O2-C2-C3-O3
3	E	606	GOL	O1-C1-C2-C3
3	E	608	GOL	C1-C2-C3-O3
3	F	609	GOL	O1-C1-C2-C3
3	F	609	GOL	C1-C2-C3-O3
3	H	607	GOL	O1-C1-C2-C3
3	C	606	GOL	O1-C1-C2-O2
3	C	606	GOL	O2-C2-C3-O3
3	C	607	GOL	O1-C1-C2-O2
3	C	609	GOL	O1-C1-C2-O2
3	F	609	GOL	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
3	B	608	GOL	O1-C1-C2-C3
3	B	608	GOL	C1-C2-C3-O3
3	D	609	GOL	O1-C1-C2-C3
3	F	608	GOL	C1-C2-C3-O3
3	F	610	GOL	O1-C1-C2-C3
3	G	608	GOL	O1-C1-C2-C3
3	H	605	GOL	C1-C2-C3-O3
3	B	607	GOL	O1-C1-C2-O2
3	B	608	GOL	O1-C1-C2-O2
3	B	608	GOL	O2-C2-C3-O3
3	E	606	GOL	O1-C1-C2-O2
3	E	608	GOL	O2-C2-C3-O3
3	F	609	GOL	O2-C2-C3-O3
3	G	608	GOL	O1-C1-C2-O2
3	H	605	GOL	O2-C2-C3-O3
3	C	609	GOL	O2-C2-C3-O3
3	D	610	GOL	O1-C1-C2-O2
3	H	607	GOL	O1-C1-C2-O2
3	C	608	GOL	O2-C2-C3-O3
3	D	609	GOL	O1-C1-C2-O2
3	B	606	GOL	O1-C1-C2-O2
3	D	609	GOL	O2-C2-C3-O3
3	H	605	GOL	O1-C1-C2-O2
3	F	610	GOL	C1-C2-C3-O3
3	F	608	GOL	O2-C2-C3-O3
3	F	610	GOL	O2-C2-C3-O3
3	F	610	GOL	O1-C1-C2-O2
3	C	607	GOL	C1-C2-C3-O3
3	E	608	GOL	O1-C1-C2-O2

There are no ring outliers.

20 monomers are involved in 38 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	H	601	SO4	1	0
3	C	609	GOL	1	0
3	E	606	GOL	1	0
3	E	608	GOL	4	0
3	F	609	GOL	2	0
2	D	602	SO4	1	0
3	H	605	GOL	1	0
3	C	607	GOL	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	G	609	GOL	1	0
3	A	606	GOL	3	0
3	C	608	GOL	2	0
3	B	607	GOL	3	0
3	F	610	GOL	2	0
3	E	607	GOL	4	0
3	B	606	GOL	1	0
2	C	603	SO4	1	0
3	H	607	GOL	4	0
2	B	603	SO4	1	0
3	A	605	GOL	2	0
2	A	604	SO4	2	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 ⓘ

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	450/465 (96%)	0.24	19 (4%)	40	48	9, 20, 47, 79	23 (5%)
1	B	453/465 (97%)	0.06	12 (2%)	57	65	11, 21, 44, 88	15 (3%)
1	C	448/465 (96%)	0.11	15 (3%)	49	57	11, 19, 45, 66	15 (3%)
1	D	448/465 (96%)	0.15	19 (4%)	40	48	11, 20, 44, 65	15 (3%)
1	E	451/465 (96%)	0.12	14 (3%)	51	60	11, 20, 44, 80	17 (3%)
1	F	453/465 (97%)	0.07	8 (1%)	67	76	11, 20, 42, 98	19 (4%)
1	G	448/465 (96%)	0.13	14 (3%)	51	60	11, 20, 43, 61	15 (3%)
1	H	448/465 (96%)	0.20	17 (3%)	44	52	12, 20, 42, 90	16 (3%)
All	All	3599/3720 (96%)	0.13	118 (3%)	49	57	9, 20, 44, 98	135 (3%)

All (118) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	76	PHE	8.7
1	H	69	VAL	6.9
1	G	69	VAL	6.8
1	H	73	VAL	6.1
1	H	68	SER	6.1
1	A	452	VAL	5.6
1	H	70	THR	5.5
1	A	455	GLY	5.3
1	H	77	THR	5.0
1	C	69	VAL	4.8
1	D	453	SER	4.7
1	H	515	THR	4.4
1	D	69	VAL	4.4
1	E	68	SER	4.4
1	H	75	LYS	4.4
1	G	453	SER	4.4

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Mol	Chain	Res	Type	RSRZ
1	C	338	GLN	4.3
1	C	335	ILE	4.3
1	F	63	GLY	4.2
1	E	437	PHE	4.2
1	G	452	VAL	3.9
1	A	453	SER	3.9
1	C	68	SER	3.9
1	G	515	THR	3.9
1	C	340	LYS	3.8
1	D	68	SER	3.7
1	A	437	PHE	3.7
1	A	69	VAL	3.7
1	A	456	CYS	3.7
1	C	437	PHE	3.6
1	E	418	ALA	3.6
1	C	337	PRO	3.6
1	C	515	THR	3.6
1	H	74	GLU	3.6
1	A	517	THR	3.5
1	B	63	GLY	3.5
1	D	515	THR	3.5
1	F	65	PHE	3.5
1	D	201	LYS	3.5
1	D	455	GLY	3.4
1	C	70	THR	3.3
1	F	515	THR	3.2
1	H	71	GLU	3.2
1	E	69	VAL	3.2
1	H	72	LYS	3.1
1	G	68	SER	3.1
1	B	515	THR	3.1
1	E	70	THR	3.1
1	E	517	THR	3.0
1	F	64	ARG	3.0
1	H	135[A]	THR	3.0
1	A	449	GLY	3.0
1	D	70	THR	3.0
1	B	418	ALA	3.0
1	C	334	SER	3.0
1	E	338	GLN	2.9
1	D	73	VAL	2.8
1	A	67	GLU	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	452	VAL	2.7
1	E	76	PHE	2.7
1	A	68	SER	2.7
1	H	453	SER	2.7
1	A	338	GLN	2.7
1	B	338	GLN	2.7
1	G	338	GLN	2.6
1	E	129	ASN	2.6
1	D	450	VAL	2.6
1	B	453	SER	2.6
1	A	128	ALA	2.5
1	B	450	VAL	2.5
1	D	74	GLU	2.5
1	G	72	LYS	2.5
1	C	339	LYS	2.5
1	C	336	MET	2.5
1	D	452	VAL	2.5
1	G	451	CYS	2.4
1	E	67	GLU	2.4
1	B	64	ARG	2.4
1	A	413	LYS	2.4
1	A	451	CYS	2.4
1	H	334	SER	2.4
1	B	438[A]	ARG	2.4
1	D	334	SER	2.4
1	E	452	VAL	2.3
1	E	454	LYS	2.3
1	A	333	SER	2.3
1	A	450	VAL	2.3
1	H	337	PRO	2.3
1	A	151	LEU	2.3
1	D	160	HIS	2.3
1	D	449	GLY	2.3
1	D	75	LYS	2.3
1	A	438	ARG	2.2
1	E	424[A]	THR	2.2
1	A	155	PRO	2.2
1	D	72	LYS	2.2
1	D	451	CYS	2.2
1	C	135	THR	2.2
1	G	117	ARG	2.2
1	F	160	HIS	2.2

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Mol	Chain	Res	Type	RSRZ
1	G	128	ALA	2.1
1	D	117	ARG	2.1
1	G	71	GLU	2.1
1	C	377	PHE	2.1
1	G	70	THR	2.1
1	F	74	GLU	2.1
1	B	335	ILE	2.1
1	F	412	ILE	2.1
1	F	453	SER	2.1
1	E	499	ASN	2.1
1	B	411	ARG	2.1
1	G	450	VAL	2.0
1	B	142	MET	2.0
1	C	333	SER	2.0
1	G	334	SER	2.0
1	D	468	LYS	2.0
1	H	117	ARG	2.0
1	H	442	ASP	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	SO4	F	604	5/5	0.73	0.16	23,28,39,44	5
3	GOL	B	607	6/6	0.75	0.22	25,35,43,46	0
3	GOL	G	609	6/6	0.75	0.26	56,60,68,70	0
2	SO4	B	605	5/5	0.79	0.13	65,70,91,95	0
2	SO4	D	604	5/5	0.80	0.20	24,32,37,38	5

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	GOL	C	609	6/6	0.81	0.18	22,28,42,43	0
3	GOL	F	609	6/6	0.81	0.19	46,54,59,60	0
2	SO4	H	601	5/5	0.81	0.13	48,64,77,92	0
3	GOL	G	608	6/6	0.82	0.25	54,62,63,64	0
2	SO4	G	603	5/5	0.83	0.15	41,48,50,52	5
3	GOL	E	607	6/6	0.84	0.18	28,30,40,42	0
3	GOL	A	606	6/6	0.84	0.17	23,32,35,41	0
3	GOL	H	607	6/6	0.84	0.16	44,51,58,61	0
3	GOL	E	608	6/6	0.85	0.17	23,26,43,47	0
3	GOL	D	610	6/6	0.85	0.16	45,54,72,96	0
3	GOL	F	610	6/6	0.85	0.17	25,36,38,40	0
3	GOL	C	607	6/6	0.86	0.17	38,60,63,63	0
2	SO4	H	604	5/5	0.86	0.18	29,38,48,50	5
2	SO4	D	602	5/5	0.86	0.11	51,57,77,79	0
2	SO4	F	605	5/5	0.86	0.16	20,35,42,43	5
3	GOL	B	608	6/6	0.86	0.16	30,36,43,61	0
2	SO4	G	607	5/5	0.87	0.12	51,74,83,86	0
2	SO4	F	607	5/5	0.87	0.15	43,43,47,50	5
3	GOL	F	608	6/6	0.87	0.15	31,46,50,50	0
3	GOL	H	605	6/6	0.87	0.15	24,36,42,51	0
2	SO4	D	607	5/5	0.87	0.14	30,43,54,60	5
2	SO4	F	602	5/5	0.88	0.13	34,48,53,58	5
3	GOL	D	609	6/6	0.88	0.18	46,53,59,60	0
2	SO4	A	604	5/5	0.88	0.15	44,52,57,58	5
3	GOL	C	608	6/6	0.88	0.17	20,41,50,51	0
2	SO4	D	605	5/5	0.89	0.13	31,38,39,47	5
2	SO4	G	606	5/5	0.89	0.14	19,30,31,35	5
3	GOL	A	605	6/6	0.89	0.15	29,43,43,53	0
2	SO4	D	608	5/5	0.89	0.12	80,83,85,95	0
2	SO4	D	606	5/5	0.90	0.11	23,30,32,36	5
2	SO4	A	603	5/5	0.90	0.10	47,59,64,67	0
3	GOL	E	606	6/6	0.90	0.14	26,36,40,41	0
3	GOL	C	606	6/6	0.90	0.16	29,40,47,48	0
2	SO4	C	604	5/5	0.91	0.13	22,33,37,42	5
2	SO4	B	604	5/5	0.91	0.12	41,44,47,48	5
2	SO4	H	602	5/5	0.91	0.10	23,31,34,35	5
2	SO4	F	603	5/5	0.91	0.12	37,38,43,45	5
2	SO4	G	605	5/5	0.91	0.14	28,37,41,45	5
3	GOL	H	606	6/6	0.91	0.15	36,44,51,53	0
2	SO4	C	603	5/5	0.91	0.10	37,38,45,46	5
2	SO4	D	603	5/5	0.92	0.13	52,61,67,69	0
3	GOL	B	606	6/6	0.92	0.14	19,37,45,49	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	H	603	5/5	0.92	0.09	36,38,42,44	5
2	SO4	F	606	5/5	0.93	0.11	23,34,46,52	5
2	SO4	B	603	5/5	0.94	0.13	32,34,44,53	5
2	SO4	E	605	5/5	0.94	0.09	38,40,47,51	5
2	SO4	G	602	5/5	0.94	0.10	36,46,47,51	0
2	SO4	G	604	5/5	0.95	0.09	47,49,57,66	0
2	SO4	C	605	5/5	0.95	0.09	43,45,52,57	5
2	SO4	B	602	5/5	0.95	0.10	43,44,48,61	0
2	SO4	A	602	5/5	0.95	0.10	17,21,30,31	5
2	SO4	F	601	5/5	0.95	0.10	24,28,33,35	5
5	NA	C	610	1/1	0.95	0.08	27,27,27,27	0
2	SO4	E	604	5/5	0.96	0.10	27,35,40,44	5
4	CL	E	601	1/1	0.96	0.12	24,24,24,24	0
4	CL	E	602	1/1	0.96	0.11	25,25,25,25	0
2	SO4	B	601	5/5	0.96	0.08	23,27,29,29	5
2	SO4	E	603	5/5	0.97	0.08	21,30,33,33	0
4	CL	C	601	1/1	0.98	0.06	15,15,15,15	0
2	SO4	A	601	5/5	0.98	0.06	19,23,24,28	5
4	CL	D	601	1/1	0.99	0.06	16,16,16,16	0
4	CL	G	601	1/1	0.99	0.09	16,16,16,16	0
4	CL	C	602	1/1	0.99	0.07	16,16,16,16	0

6.5 Other polymers [i](#)

There are no such residues in this entry.