



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:55 PM UTC

PDB ID : 1XG3 / pdb_00001xg3
Title : Crystal structure of the C123S 2-methylisocitrate lyase mutant from Escherichia coli in complex with the reaction product, Mg(II)-pyruvate and succinate
Authors : Liu, S.; Lu, Z.; Han, Y.; Melamud, E.; Dunaway-Mariano, D.; Herzberg, O.
Deposited on : 2004-09-16
Resolution : 1.90 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

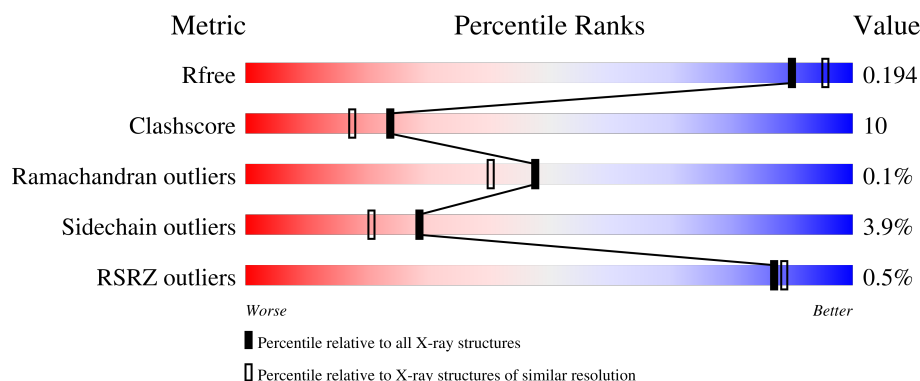
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	295	 78% 17% . .
1	B	295	 76% 18% . .
1	C	295	 74% 21% . .
1	D	295	 77% 17% . .

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 9815 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 Probable methylisocitrate lyase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	288	Total	C	N	O	S	0	0	0
			2190	1375	383	423	9			
1	B	286	Total	C	N	O	S	0	0	0
			2176	1366	381	420	9			
1	C	287	Total	C	N	O	S	0	0	0
			2184	1372	382	421	9			
1	D	283	Total	C	N	O	S	0	0	0
			2152	1352	377	414	9			

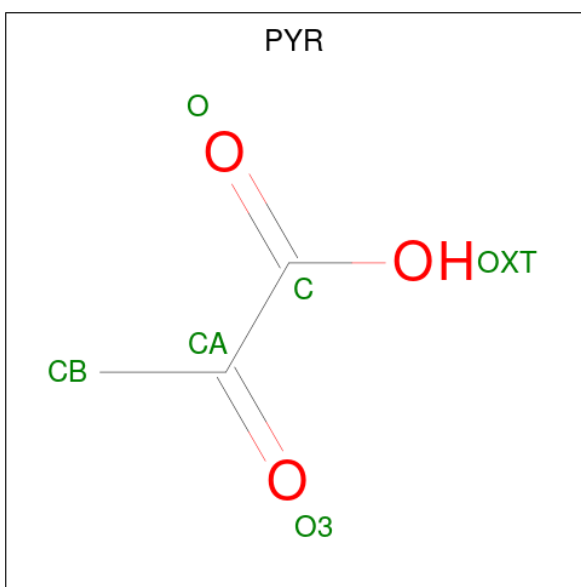
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	123	SER	CYS	engineered mutation	UNP P77541
B	123	SER	CYS	engineered mutation	UNP P77541
C	123	SER	CYS	engineered mutation	UNP P77541
D	123	SER	CYS	engineered mutation	UNP P77541

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

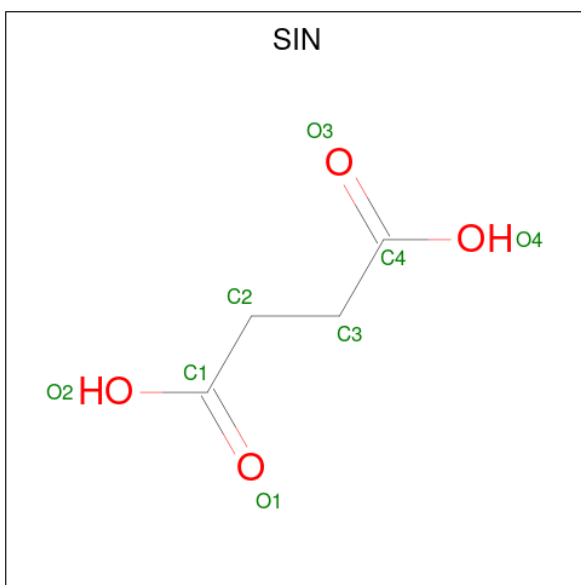
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		
2	B	1	Total	Mg	0	0
			1	1		
2	C	1	Total	Mg	0	0
			1	1		
2	D	1	Total	Mg	0	0
			1	1		

- Molecule 3 is PYRUVIC ACID (CCD ID: PYR) (formula: C₃H₄O₃).



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

- Molecule 4 is SUCCINIC ACID (CCD ID: SIN) (formula: C₄H₆O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			8	4	4		
4	B	1	Total	C	O	0	0
			8	4	4		
4	D	1	Total	C	O	0	0
			8	4	4		

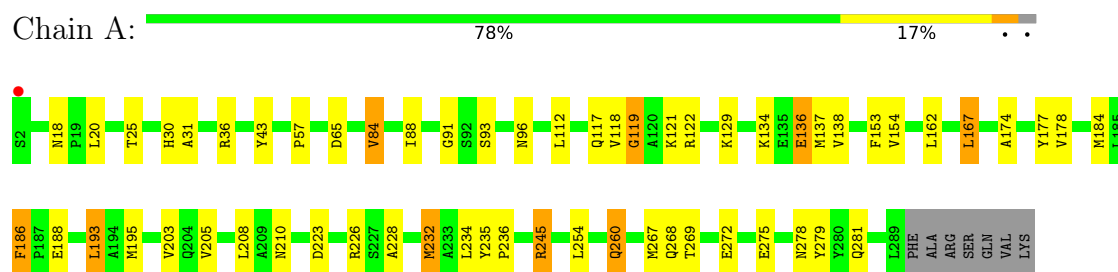
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	250	Total	O	0	0
			250	250		
5	B	277	Total	O	0	0
			277	277		
5	C	267	Total	O	0	0
			267	267		
5	D	267	Total	O	0	0
			267	267		

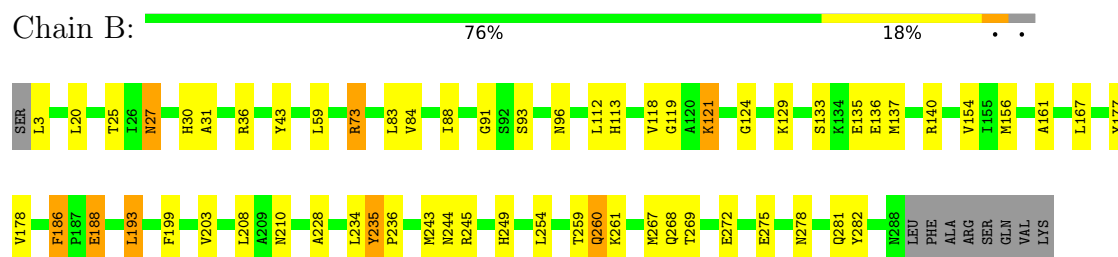
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

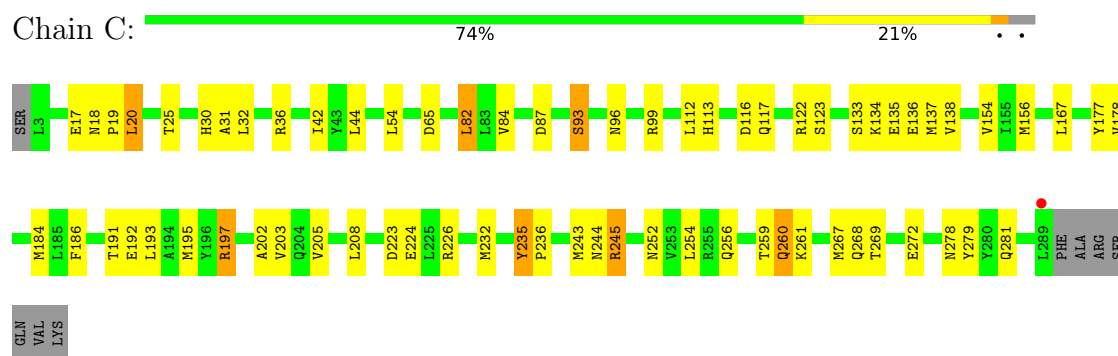
- Molecule 1: Probable methylisocitrate lyase



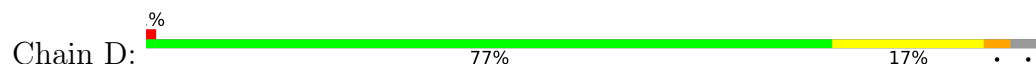
- Molecule 1: Probable methylisocitrate lyase

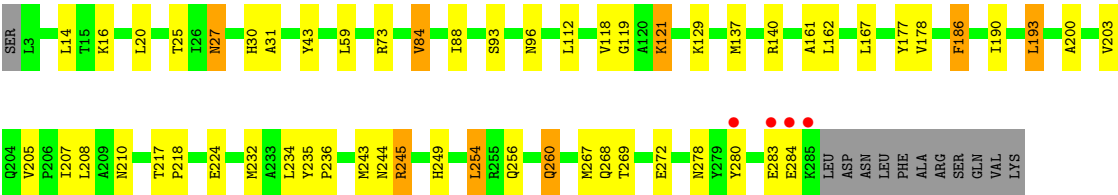


- Molecule 1: Probable methylisocitrate lyase



- Molecule 1: Probable methylisocitrate lyase





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	159.36Å 85.99Å 99.60Å 90.00° 108.28° 90.00°	Depositor
Resolution (Å)	25.85 – 1.90 25.85 – 1.90	Depositor EDS
% Data completeness (in resolution range)	88.3 (25.85-1.90) 88.5 (25.85-1.90)	Depositor EDS
R_{merge}	0.03	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.98 (at 1.91Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.161 , 0.195 0.160 , 0.194	Depositor DCC
R_{free} test set	3591 reflections (4.03%)	wwPDB-VP
Wilson B-factor (Å ²)	17.5	Xtriage
Anisotropy	0.085	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 44.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9815	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.65% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PYR, MG, SIN

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	0.47	1/2224 (0.0%)	0.94	12/3016 (0.4%)
1	B	0.49	1/2210 (0.0%)	0.94	9/2997 (0.3%)
1	C	0.49	0/2218	0.97	12/3008 (0.4%)
1	D	0.50	1/2186 (0.0%)	0.94	7/2964 (0.2%)
All	All	0.49	3/8838 (0.0%)	0.95	40/11985 (0.3%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	210	ASN	CG-OD1	5.41	1.33	1.23
1	D	210	ASN	CG-OD1	5.12	1.33	1.23
1	A	210	ASN	CG-OD1	5.11	1.33	1.23

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	236	PRO	N-CA-C	7.20	123.54	113.53
1	A	268	GLN	N-CA-C	-7.04	100.03	110.23
1	C	93	SER	N-CA-C	6.77	118.40	108.86
1	C	279	TYR	N-CA-C	6.60	118.13	111.07
1	A	279	TYR	N-CA-C	6.41	118.80	111.11
1	B	161	ALA	N-CA-C	6.38	120.32	112.54
1	D	167	LEU	N-CA-C	6.29	117.80	111.07
1	C	268	GLN	N-CA-C	-6.26	100.57	109.96
1	A	236	PRO	N-CA-C	6.21	122.16	113.53
1	C	267	MET	N-CA-C	6.16	119.86	109.76
1	A	205	VAL	N-CA-C	-6.13	102.08	109.01
1	C	236	PRO	N-CA-C	6.12	122.87	113.75
1	D	267	MET	N-CA-C	6.08	119.72	109.76
1	B	154	VAL	N-CA-C	6.05	117.01	108.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	236	PRO	N-CA-C	5.94	122.60	113.75
1	A	267	MET	N-CA-C	5.83	119.32	109.76
1	B	235	TYR	CA-C-N	5.72	125.84	119.32
1	B	235	TYR	C-N-CA	5.72	125.84	119.32
1	B	267	MET	N-CA-C	5.58	119.34	109.96
1	C	235	TYR	CA-C-N	5.57	125.67	119.32
1	C	235	TYR	C-N-CA	5.57	125.67	119.32
1	C	202	ALA	N-CA-C	5.57	117.43	111.36
1	A	188	GLU	N-CA-C	5.56	118.71	110.48
1	A	154	VAL	N-CA-C	5.46	116.17	108.36
1	B	188	GLU	N-CA-C	5.40	118.12	110.50
1	D	161	ALA	N-CA-C	5.40	119.13	112.54
1	D	268	GLN	N-CA-C	-5.36	102.45	110.23
1	D	205	VAL	N-CA-C	-5.34	102.05	108.45
1	A	153	PHE	N-CA-C	-5.31	103.12	110.35
1	B	167	LEU	N-CA-C	5.27	116.72	110.97
1	A	84	VAL	N-CA-C	5.26	116.06	108.48
1	C	205	VAL	N-CA-C	-5.26	102.14	108.45
1	C	116	ASP	N-CA-C	5.25	119.38	112.92
1	C	154	VAL	N-CA-C	5.18	115.77	108.36
1	A	235	TYR	CA-C-N	5.14	124.76	119.05
1	A	235	TYR	C-N-CA	5.14	124.76	119.05
1	C	84	VAL	N-CA-C	5.11	115.84	108.48
1	D	84	VAL	N-CA-C	5.07	115.77	108.48
1	A	167	LEU	N-CA-C	5.04	116.85	111.36
1	B	268	GLN	N-CA-C	-5.02	102.43	109.96

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	2190	0	2189	45	0
1	B	2176	0	2173	46	0
1	C	2184	0	2184	49	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	2152	0	2152	47	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	6	0	0	0	0
3	B	6	0	0	0	0
3	C	6	0	0	0	0
3	D	6	0	0	0	0
4	A	8	0	4	0	0
4	B	8	0	4	0	0
4	D	8	0	4	0	0
5	A	250	0	0	7	0
5	B	277	0	0	8	0
5	C	267	0	0	6	0
5	D	267	0	0	8	0
All	All	9815	0	8710	176	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 10.

All (176) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:260:GLN:HE21	1:A:260:GLN:H	1.08	0.97
1:B:260:GLN:H	1:B:260:GLN:HE21	1.01	0.96
1:C:260:GLN:HE21	1:C:260:GLN:H	1.03	0.96
1:D:260:GLN:HE21	1:D:260:GLN:H	0.98	0.91
1:A:117:GLN:HE22	1:A:122:ARG:H	1.27	0.80
1:B:88:ILE:HD13	1:B:119:GLY:HA2	1.64	0.79
1:C:93:SER:H	1:C:96:ASN:HD22	1.30	0.77
1:D:260:GLN:HE21	1:D:260:GLN:N	1.81	0.77
1:B:93:SER:H	1:B:96:ASN:HD22	1.34	0.76
1:D:93:SER:H	1:D:96:ASN:HD22	1.32	0.76
1:B:260:GLN:HE21	1:B:260:GLN:N	1.82	0.74
1:C:260:GLN:HE21	1:C:260:GLN:N	1.85	0.72
5:C:2333:HOH:O	1:D:73:ARG:HD2	1.90	0.71
1:D:260:GLN:H	1:D:260:GLN:NE2	1.82	0.69
1:B:36:ARG:HD3	5:B:2426:HOH:O	1.93	0.69
1:C:245:ARG:HD3	5:D:2473:HOH:O	1.91	0.69
1:A:93:SER:H	1:A:96:ASN:HD22	1.40	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:88:ILE:CD1	1:B:119:GLY:HA2	2.23	0.68
1:D:280:TYR:HA	1:D:283:GLU:HG2	1.76	0.68
5:A:2171:HOH:O	1:B:245:ARG:HD3	1.93	0.67
1:B:27:ASN:HD22	1:B:27:ASN:C	2.04	0.65
1:C:197:ARG:HG3	5:C:2459:HOH:O	1.98	0.64
1:B:118:VAL:HG12	1:B:119:GLY:N	2.11	0.64
1:D:112:LEU:HD12	1:D:112:LEU:C	2.23	0.64
1:A:167:LEU:HD23	1:A:195:MET:CE	2.27	0.64
1:C:42:ILE:HG13	1:C:82:LEU:HD23	1.80	0.63
1:D:200:ALA:HB2	1:D:207:ILE:HD13	1.81	0.63
1:A:245:ARG:HD3	5:A:2147:HOH:O	1.99	0.63
5:C:2449:HOH:O	1:D:245:ARG:HD3	1.99	0.62
1:B:129:LYS:HG2	5:B:2252:HOH:O	1.99	0.62
1:D:88:ILE:HD13	1:D:119:GLY:HA2	1.82	0.61
1:B:140:ARG:HD3	5:B:2268:HOH:O	2.01	0.61
1:C:254:LEU:HD21	1:D:235:TYR:CG	2.36	0.60
1:C:197:ARG:HG3	1:C:197:ARG:HH11	1.66	0.60
1:D:16:LYS:HE2	5:D:2566:HOH:O	2.02	0.60
1:B:259:THR:OG1	1:B:261:LYS:HG3	2.02	0.59
1:D:140:ARG:HD3	5:D:2633:HOH:O	2.01	0.59
1:D:193:LEU:HD11	1:D:224:GLU:HB3	1.84	0.59
1:A:117:GLN:NE2	1:A:122:ARG:H	1.99	0.59
1:B:278:ASN:HD21	1:B:281:GLN:NE2	2.00	0.59
1:C:184:MET:SD	1:C:232:MET:HE1	2.43	0.59
1:C:112:LEU:HD12	1:C:112:LEU:C	2.28	0.58
1:A:184:MET:SD	1:A:232:MET:HE1	2.43	0.58
1:D:88:ILE:CD1	1:D:119:GLY:HA2	2.34	0.58
1:A:260:GLN:HE21	1:A:260:GLN:N	1.89	0.58
1:B:27:ASN:ND2	1:B:30:HIS:H	2.02	0.56
1:C:65:ASP:HB2	5:C:2450:HOH:O	2.05	0.56
1:C:117:GLN:HE22	1:C:123:SER:H	1.53	0.56
1:D:118:VAL:HG12	1:D:119:GLY:N	2.19	0.56
1:A:167:LEU:HD23	1:A:195:MET:HE3	1.88	0.56
1:B:112:LEU:C	1:B:112:LEU:HD12	2.31	0.56
1:B:260:GLN:H	1:B:260:GLN:NE2	1.86	0.56
1:A:118:VAL:HG12	1:A:119:GLY:N	2.20	0.56
1:C:223:ASP:OD1	1:C:226:ARG:NH2	2.40	0.55
1:D:27:ASN:C	1:D:27:ASN:HD22	2.14	0.55
1:C:252:ASN:O	1:C:256:GLN:HG3	2.07	0.54
1:C:87:ASP:HB2	5:C:2522:HOH:O	2.08	0.54
1:D:27:ASN:ND2	1:D:30:HIS:H	2.06	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:269:THR:OG1	1:D:272:GLU:HG3	2.08	0.54
1:D:186:PHE:CD1	1:D:186:PHE:C	2.87	0.53
1:A:167:LEU:C	1:A:167:LEU:HD13	2.34	0.53
1:C:192:GLU:H	1:C:195:MET:HE3	1.72	0.53
1:A:129:LYS:HG2	5:A:2244:HOH:O	2.08	0.53
1:B:118:VAL:HG12	1:B:119:GLY:H	1.74	0.53
1:C:32:LEU:O	1:C:36:ARG:HG3	2.09	0.52
1:B:30:HIS:HD2	5:B:2235:HOH:O	1.92	0.52
1:C:254:LEU:HD21	1:D:235:TYR:CD2	2.45	0.52
1:D:118:VAL:HG12	1:D:119:GLY:H	1.73	0.52
1:D:278:ASN:ND2	5:D:2567:HOH:O	2.40	0.52
1:D:283:GLU:HG3	1:D:284:GLU:N	2.25	0.52
1:A:223:ASP:OD1	1:A:226:ARG:NH2	2.42	0.51
5:A:2339:HOH:O	1:C:99:ARG:HD3	2.10	0.51
1:C:25:THR:HG21	1:C:31:ALA:HA	1.93	0.51
1:A:30:HIS:HE1	1:B:244:ASN:OD1	1.94	0.51
1:B:124:GLY:HA3	1:B:188:GLU:HG2	1.93	0.51
1:C:278:ASN:HD21	1:C:281:GLN:CD	2.19	0.50
1:D:162:LEU:HD22	1:D:190:ILE:HD13	1.93	0.50
1:C:244:ASN:OD1	1:D:30:HIS:HE1	1.94	0.50
1:D:25:THR:HG21	1:D:31:ALA:HA	1.93	0.50
1:B:135:GLU:H	1:B:135:GLU:CD	2.20	0.50
1:B:199:PHE:O	1:B:203:VAL:HG22	2.11	0.50
1:C:208:LEU:C	1:C:208:LEU:HD23	2.37	0.50
1:A:36:ARG:NH2	1:A:275:GLU:HG2	2.27	0.49
1:A:186:PHE:CD1	1:A:186:PHE:C	2.89	0.49
1:A:137:MET:HG2	1:A:177:TYR:CZ	2.48	0.49
1:C:113:HIS:HA	1:C:156:MET:O	2.11	0.49
1:A:25:THR:HG21	1:A:31:ALA:HA	1.95	0.49
1:A:178:VAL:HG21	1:A:203:VAL:HB	1.95	0.49
1:C:259:THR:OG1	1:C:261:LYS:HG3	2.13	0.48
1:D:186:PHE:C	1:D:186:PHE:HD1	2.21	0.48
1:B:118:VAL:CG1	1:B:119:GLY:N	2.76	0.48
1:B:137:MET:HG2	1:B:177:TYR:CZ	2.48	0.48
1:A:208:LEU:C	1:A:208:LEU:HD23	2.38	0.48
1:B:178:VAL:HG21	1:B:203:VAL:HB	1.96	0.48
1:A:254:LEU:HD21	1:B:235:TYR:CG	2.49	0.47
1:B:186:PHE:C	1:B:186:PHE:CD1	2.91	0.47
1:D:59:LEU:HA	1:D:121:LYS:O	2.14	0.47
1:A:18:ASN:O	1:A:226:ARG:HD3	2.15	0.47
1:A:88:ILE:CD1	1:A:119:GLY:HA2	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:134:LYS:O	1:A:138:VAL:HG23	2.14	0.47
1:B:25:THR:HG21	1:B:31:ALA:HA	1.96	0.47
1:C:186:PHE:C	1:C:186:PHE:CD1	2.92	0.47
1:B:275:GLU:HG2	5:B:2446:HOH:O	2.15	0.46
1:C:93:SER:H	1:C:96:ASN:ND2	2.07	0.46
1:C:191:THR:H	1:C:195:MET:CE	2.28	0.46
1:C:269:THR:OG1	1:C:272:GLU:HG3	2.15	0.46
1:A:118:VAL:HG12	1:A:119:GLY:H	1.80	0.46
1:C:117:GLN:HE21	1:C:122:ARG:H	1.63	0.46
1:A:269:THR:OG1	1:A:272:GLU:HG3	2.15	0.46
1:D:84:VAL:O	1:D:112:LEU:HA	2.16	0.46
1:D:137:MET:HG2	1:D:177:TYR:CZ	2.50	0.46
1:D:129:LYS:HG2	5:D:2493:HOH:O	2.16	0.46
1:A:91:GLY:HA3	1:A:96:ASN:CB	2.46	0.46
1:A:112:LEU:C	1:A:112:LEU:HD12	2.41	0.46
1:C:178:VAL:HG21	1:C:203:VAL:HB	1.97	0.46
5:A:2150:HOH:O	1:B:73:ARG:HD2	2.15	0.46
1:C:191:THR:H	1:C:195:MET:HE1	1.80	0.45
1:B:113:HIS:HA	1:B:156:MET:O	2.17	0.45
1:A:43:TYR:CG	1:A:234:LEU:HD11	2.52	0.45
1:A:186:PHE:C	1:A:186:PHE:HD1	2.24	0.45
1:A:84:VAL:O	1:A:112:LEU:HA	2.16	0.45
1:A:57:PRO:HG2	1:B:282:TYR:CD1	2.52	0.45
1:B:249:HIS:HE1	5:B:2414:HOH:O	1.99	0.45
1:C:30:HIS:HE1	1:D:244:ASN:OD1	1.99	0.45
1:D:30:HIS:HD2	5:D:2411:HOH:O	1.99	0.45
1:A:65:ASP:OD1	1:C:65:ASP:HB3	2.17	0.45
1:C:193:LEU:HD11	1:C:224:GLU:HB3	1.99	0.45
1:D:178:VAL:HG21	1:D:203:VAL:HB	2.00	0.44
1:A:88:ILE:HD13	1:A:119:GLY:HA2	1.98	0.44
1:B:84:VAL:O	1:B:112:LEU:HA	2.17	0.44
1:A:162:LEU:CD2	1:A:195:MET:HE1	2.47	0.44
1:B:186:PHE:C	1:B:186:PHE:HD1	2.25	0.44
1:D:249:HIS:HD2	5:D:2572:HOH:O	2.00	0.44
1:C:54:LEU:HD23	1:D:73:ARG:HG2	1.99	0.44
1:B:118:VAL:CG1	1:B:119:GLY:H	2.30	0.44
1:A:260:GLN:H	1:A:260:GLN:NE2	1.93	0.44
1:D:14:LEU:HD21	1:D:232:MET:CG	2.47	0.44
1:C:260:GLN:H	1:C:260:GLN:NE2	1.88	0.44
1:C:134:LYS:O	1:C:138:VAL:HG23	2.18	0.43
1:B:269:THR:OG1	1:B:272:GLU:HG3	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:18:ASN:HA	1:C:19:PRO:HA	1.90	0.43
1:C:30:HIS:HD2	5:C:2314:HOH:O	2.00	0.43
1:C:32:LEU:O	1:C:36:ARG:CG	2.66	0.43
1:B:135:GLU:CD	1:B:135:GLU:N	2.77	0.43
1:D:112:LEU:C	1:D:112:LEU:CD1	2.92	0.43
1:B:249:HIS:HD2	5:B:2347:HOH:O	2.01	0.43
1:C:235:TYR:CG	1:D:254:LEU:HD11	2.54	0.43
1:C:192:GLU:H	1:C:195:MET:CE	2.32	0.42
1:A:162:LEU:CD1	1:A:195:MET:HE1	2.50	0.42
1:D:280:TYR:HD2	1:D:283:GLU:OE2	2.02	0.42
1:B:91:GLY:HA3	1:B:96:ASN:CG	2.44	0.42
1:C:137:MET:HG2	1:C:177:TYR:CZ	2.55	0.42
1:D:208:LEU:C	1:D:208:LEU:HD23	2.45	0.42
1:D:256:GLN:HG2	5:D:2644:HOH:O	2.19	0.42
1:C:17:GLU:HG2	1:C:20:LEU:HA	2.01	0.42
1:C:186:PHE:C	1:C:186:PHE:HD1	2.28	0.42
1:A:30:HIS:HD2	5:A:2112:HOH:O	2.02	0.42
1:A:136:GLU:HG3	5:A:2310:HOH:O	2.19	0.42
1:B:3:LEU:N	5:B:2377:HOH:O	2.53	0.42
1:B:43:TYR:CG	1:B:234:LEU:HD11	2.55	0.42
1:D:217:THR:HA	1:D:218:PRO:HD3	1.97	0.42
1:A:193:LEU:HD22	1:A:228:ALA:HB2	2.02	0.41
1:D:200:ALA:HB2	1:D:207:ILE:CD1	2.47	0.41
1:A:278:ASN:HD21	1:A:281:GLN:CD	2.28	0.41
1:B:193:LEU:HD22	1:B:228:ALA:HB2	2.02	0.41
1:C:133:SER:OG	1:C:136:GLU:HG3	2.20	0.41
1:B:59:LEU:HA	1:B:121:LYS:O	2.20	0.41
1:B:208:LEU:HD23	1:B:208:LEU:C	2.45	0.41
1:B:133:SER:OG	1:B:136:GLU:HG3	2.20	0.41
1:A:174:ALA:O	1:A:178:VAL:HG23	2.21	0.41
1:C:254:LEU:HD11	1:D:235:TYR:CE1	2.56	0.41
1:A:118:VAL:CG1	1:A:119:GLY:N	2.84	0.41
1:A:167:LEU:HD23	1:A:195:MET:HE2	1.99	0.41
1:A:278:ASN:HD21	1:A:281:GLN:NE2	2.20	0.40
1:D:43:TYR:CG	1:D:234:LEU:HD11	2.56	0.40
1:C:135:GLU:CD	1:C:135:GLU:H	2.29	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	286/295 (97%)	279 (98%)	6 (2%)	1 (0%)	36	29
1	B	284/295 (96%)	277 (98%)	7 (2%)	0	100	100
1	C	285/295 (97%)	279 (98%)	6 (2%)	0	100	100
1	D	281/295 (95%)	271 (96%)	10 (4%)	0	100	100
All	All	1136/1180 (96%)	1106 (97%)	29 (3%)	1 (0%)	48	40

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	119	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	226/232 (97%)	218 (96%)	8 (4%)	32	24
1	B	224/232 (97%)	214 (96%)	10 (4%)	24	17
1	C	225/232 (97%)	217 (96%)	8 (4%)	31	23
1	D	221/232 (95%)	212 (96%)	9 (4%)	27	19
All	All	896/928 (97%)	861 (96%)	35 (4%)	28	21

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	LEU
1	A	121	LYS
1	A	136	GLU
1	A	186	PHE
1	A	193	LEU
1	A	232	MET
1	A	245	ARG
1	A	260	GLN
1	B	20	LEU
1	B	27	ASN
1	B	73	ARG
1	B	83	LEU
1	B	121	LYS
1	B	186	PHE
1	B	193	LEU
1	B	243	MET
1	B	254	LEU
1	B	260	GLN
1	C	20	LEU
1	C	44	LEU
1	C	82	LEU
1	C	167	LEU
1	C	197	ARG
1	C	243	MET
1	C	245	ARG
1	C	260	GLN
1	D	20	LEU
1	D	27	ASN
1	D	121	LYS
1	D	186	PHE
1	D	193	LEU
1	D	243	MET
1	D	245	ARG
1	D	254	LEU
1	D	260	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (34) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	30	HIS
1	A	96	ASN
1	A	117	GLN
1	A	204	GLN

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Mol	Chain	Res	Type
1	A	229	HIS
1	A	244	ASN
1	A	260	GLN
1	A	268	GLN
1	A	278	ASN
1	B	27	ASN
1	B	30	HIS
1	B	35	GLN
1	B	96	ASN
1	B	244	ASN
1	B	249	HIS
1	B	256	GLN
1	B	260	GLN
1	B	278	ASN
1	B	281	GLN
1	C	30	HIS
1	C	35	GLN
1	C	40	GLN
1	C	96	ASN
1	C	117	GLN
1	C	260	GLN
1	C	278	ASN
1	C	281	GLN
1	D	27	ASN
1	D	30	HIS
1	D	96	ASN
1	D	204	GLN
1	D	249	HIS
1	D	260	GLN
1	D	278	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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 11 ligands modelled in this entry, 4 are monoatomic - leaving 7 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	PYR	B	2205	2	5,5,5	1.57	1 (20%)	3,6,6	1.93	1 (33%)
4	SIN	A	2106	-	7,7,7	1.13	0	8,8,8	1.13	0
4	SIN	B	2206	-	7,7,7	1.16	0	8,8,8	1.18	0
3	PYR	C	2305	2	5,5,5	1.61	1 (20%)	3,6,6	1.78	1 (33%)
4	SIN	D	2406	-	7,7,7	1.25	0	8,8,8	1.19	0
3	PYR	D	2405	2	5,5,5	1.50	1 (20%)	3,6,6	1.88	1 (33%)
3	PYR	A	2105	2	5,5,5	1.37	1 (20%)	3,6,6	1.94	1 (33%)

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	PYR	B	2205	2	-	0/4/4/4	-
4	SIN	A	2106	-	-	1/5/5/5	-
4	SIN	B	2206	-	-	2/5/5/5	-
3	PYR	C	2305	2	-	0/4/4/4	-
4	SIN	D	2406	-	-	1/5/5/5	-
3	PYR	D	2405	2	-	0/4/4/4	-
3	PYR	A	2105	2	-	0/4/4/4	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	2205	PYR	OXT-C	-2.31	1.24	1.30
3	C	2305	PYR	O-C	2.20	1.28	1.22
3	D	2405	PYR	O-C	2.06	1.27	1.22
3	A	2105	PYR	CB-CA	2.02	1.54	1.50

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	2105	PYR	OXT-C-CA	3.11	122.21	113.59
3	B	2205	PYR	OXT-C-CA	3.10	122.19	113.59
3	D	2405	PYR	OXT-C-CA	3.01	121.93	113.59
3	C	2305	PYR	OXT-C-CA	2.79	121.34	113.59

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	2106	SIN	C1-C2-C3-C4
4	D	2406	SIN	C1-C2-C3-C4
4	B	2206	SIN	C1-C2-C3-C4
4	B	2206	SIN	O2-C1-C2-C3

There are no ring outliers.

No monomer is involved in short contacts.

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	288/295 (97%)	-0.41	1 (0%) 90 91	10, 19, 31, 41	0
1	B	286/295 (96%)	-0.50	0 100 100	11, 18, 27, 35	0
1	C	287/295 (97%)	-0.51	1 (0%) 90 91	10, 17, 29, 39	0
1	D	283/295 (95%)	-0.48	4 (1%) 73 76	9, 16, 29, 64	0
All	All	1144/1180 (96%)	-0.47	6 (0%) 87 89	9, 18, 29, 64	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	284	GLU	3.1
1	D	285	LYS	3.1
1	D	283	GLU	2.7
1	D	280	TYR	2.6
1	C	289	LEU	2.3
1	A	2	SER	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
4	SIN	D	2406	8/8	0.88	0.10	18,22,26,29	0
4	SIN	B	2206	8/8	0.89	0.11	22,26,28,30	0
4	SIN	A	2106	8/8	0.90	0.10	30,37,38,40	0
3	PYR	A	2105	6/6	0.97	0.05	14,16,17,17	0
3	PYR	C	2305	6/6	0.97	0.05	16,18,18,19	0
3	PYR	D	2405	6/6	0.97	0.04	12,13,14,15	0
3	PYR	B	2205	6/6	0.98	0.04	13,16,16,16	0
2	MG	B	2201	1/1	0.98	0.04	15,15,15,15	0
2	MG	C	2301	1/1	0.99	0.06	17,17,17,17	0
2	MG	D	2401	1/1	0.99	0.01	14,14,14,14	0
2	MG	A	2101	1/1	0.99	0.03	15,15,15,15	0

6.5 Other polymers [i](#)

There are no such residues in this entry.