Full wwPDB X-ray Structure Validation Report ⓘ

Sep 26, 2023 – 03:51 AM EDT

PDB ID	:	6BOM
Title	:	Crystal structure of mutant 2-methylcitrate synthase mcsAG306A from Aspergillus fumigatus.
Authors	:	Schlachter, C.; Chruszcz, M.
Deposited on	:	2017-11-20
Resolution	:	2.05 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:

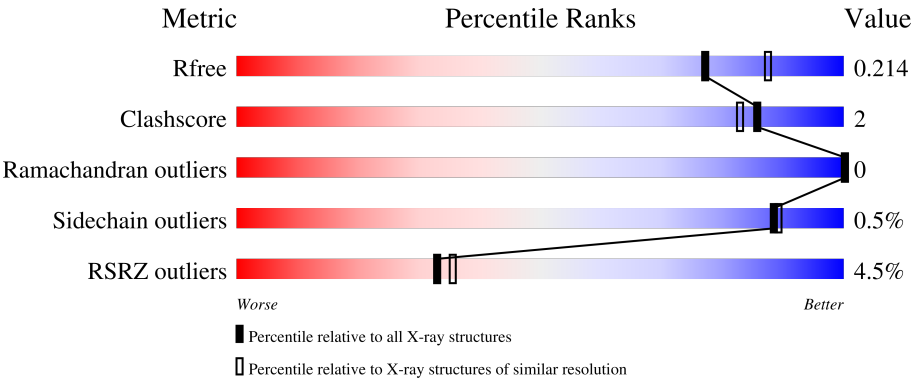
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

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 2.05 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	1692 (2.04-2.04)
Clashscore	141614	1773 (2.04-2.04)
Ramachandran outliers	138981	1752 (2.04-2.04)
Sidechain outliers	138945	1752 (2.04-2.04)
RSRZ outliers	127900	1672 (2.04-2.04)

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	441	3% 93% 7%
1	B	441	2% 93% 6%
1	C	441	% 90% 8%
1	D	441	7% 91% 7%
1	E	441	8% 90% 8%

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Mol	Chain	Length	Quality of chain
1	F	441	5% 92% 6%

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	TRS	B	501	-	X	-	-

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 21359 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 2-methylcitrate synthase, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	440	Total	C	N	O	S	0	1	0
			3426	2195	592	630	9			
1	B	435	Total	C	N	O	S	0	2	0
			3408	2186	588	625	9			
1	C	435	Total	C	N	O	S	0	0	0
			3367	2158	578	622	9			
1	D	436	Total	C	N	O	S	0	0	0
			3367	2158	577	623	9			
1	E	434	Total	C	N	O	S	0	0	0
			3364	2157	578	620	9			
1	F	434	Total	C	N	O	S	0	0	0
			3360	2154	577	620	9			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	25	SER	-	expression tag	UNP B0YD89
A	26	GLY	-	expression tag	UNP B0YD89
A	27	SER	-	expression tag	UNP B0YD89
A	28	GLY	-	expression tag	UNP B0YD89
A	306	ALA	GLY	engineered mutation	UNP B0YD89
B	25	SER	-	expression tag	UNP B0YD89
B	26	GLY	-	expression tag	UNP B0YD89
B	27	SER	-	expression tag	UNP B0YD89
B	28	GLY	-	expression tag	UNP B0YD89
B	306	ALA	GLY	engineered mutation	UNP B0YD89
C	25	SER	-	expression tag	UNP B0YD89
C	26	GLY	-	expression tag	UNP B0YD89
C	27	SER	-	expression tag	UNP B0YD89
C	28	GLY	-	expression tag	UNP B0YD89
C	306	ALA	GLY	engineered mutation	UNP B0YD89
D	25	SER	-	expression tag	UNP B0YD89
D	26	GLY	-	expression tag	UNP B0YD89

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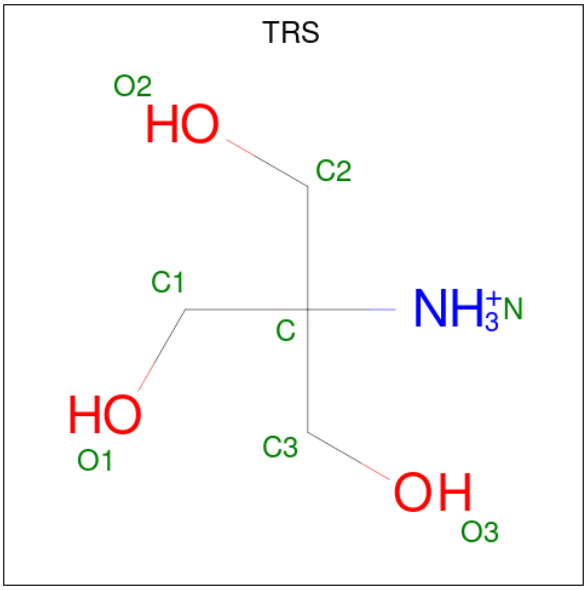
Chain	Residue	Modelled	Actual	Comment	Reference
D	27	SER	-	expression tag	UNP B0YD89
D	28	GLY	-	expression tag	UNP B0YD89
D	306	ALA	GLY	engineered mutation	UNP B0YD89
E	25	SER	-	expression tag	UNP B0YD89
E	26	GLY	-	expression tag	UNP B0YD89
E	27	SER	-	expression tag	UNP B0YD89
E	28	GLY	-	expression tag	UNP B0YD89
E	306	ALA	GLY	engineered mutation	UNP B0YD89
F	25	SER	-	expression tag	UNP B0YD89
F	26	GLY	-	expression tag	UNP B0YD89
F	27	SER	-	expression tag	UNP B0YD89
F	28	GLY	-	expression tag	UNP B0YD89
F	306	ALA	GLY	engineered mutation	UNP B0YD89

- Molecule 2 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula: $C_4H_{12}NO_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	N	O	0	0
			8	4	1	3		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	232	Total	O	0	0
			232	232		

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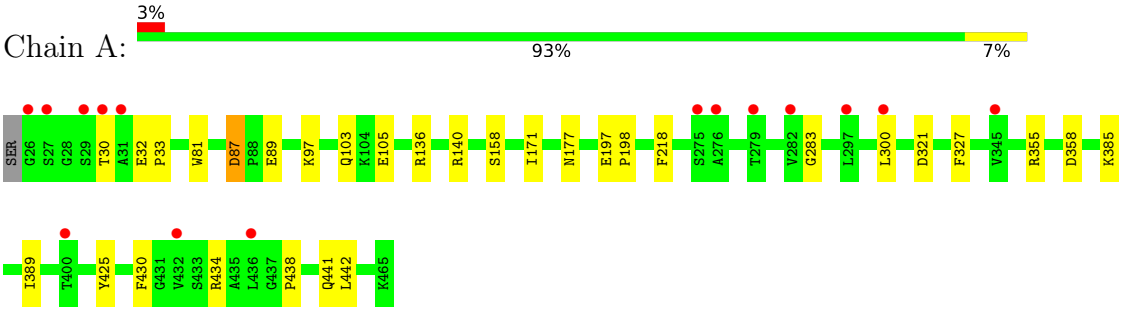
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	258	Total O 258 258	0	0
3	C	182	Total O 182 182	0	0
3	D	171	Total O 171 171	0	0
3	E	110	Total O 110 110	0	0
3	F	106	Total O 106 106	0	0

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.

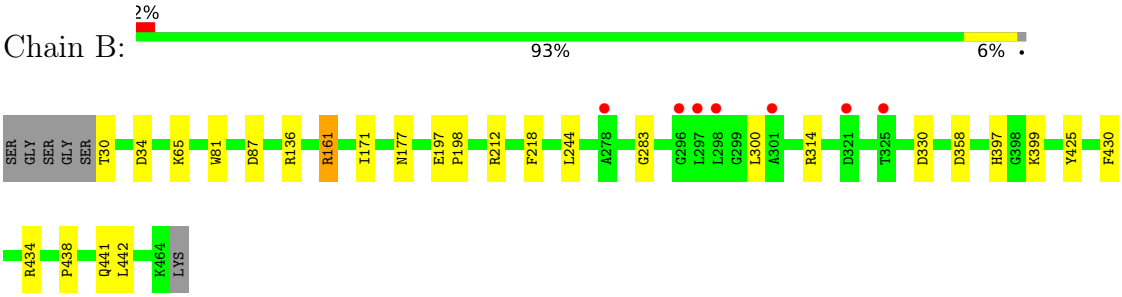
• Molecule 1: 2-methylcitrate synthase, mitochondrial

Chain A:



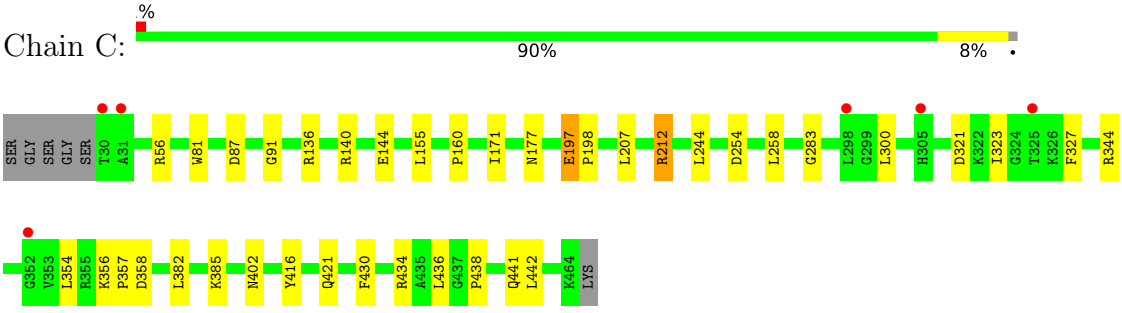
• Molecule 1: 2-methylcitrate synthase, mitochondrial

Chain B:



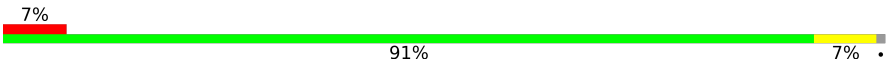
• Molecule 1: 2-methylcitrate synthase, mitochondrial

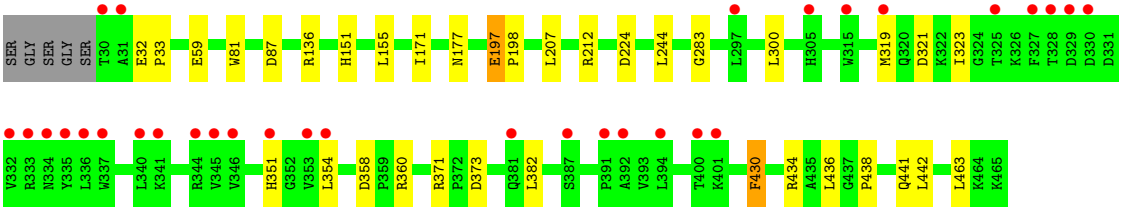
Chain C:



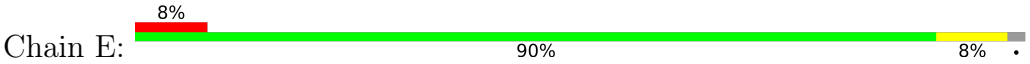
• Molecule 1: 2-methylcitrate synthase, mitochondrial

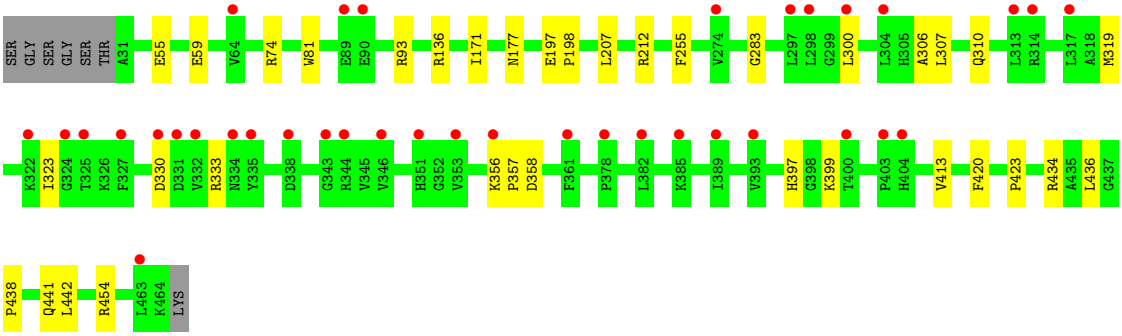
Chain D:



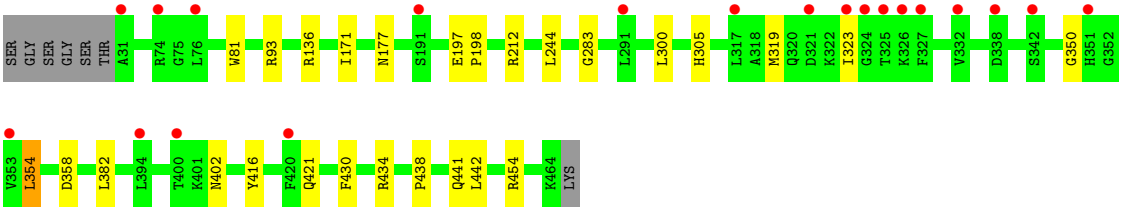


• Molecule 1: 2-methylcitrate synthase, mitochondrial





• Molecule 1: 2-methylcitrate synthase, mitochondrial



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	82.02Å 128.86Å 258.46Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.05 39.87 – 2.05	Depositor EDS
% Data completeness (in resolution range)	99.7 (40.00-2.05) 99.7 (39.87-2.05)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$< I/\sigma(I) >$ [1]	1.95 (at 2.05Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.181 , 0.212 0.187 , 0.214	Depositor DCC
R_{free} test set	8687 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å^2)	40.4	Xtriage
Anisotropy	0.108	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.29 , 38.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.46$, $< L^2 > = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	21359	wwPDB-VP
Average B, all atoms (Å^2)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.35% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: TRS

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	0.80	1/3509 (0.0%)	0.84	5/4760 (0.1%)
1	B	0.82	0/3491	0.85	8/4737 (0.2%)
1	C	0.83	0/3448	0.90	13/4684 (0.3%)
1	D	0.79	3/3449 (0.1%)	0.86	9/4688 (0.2%)
1	E	0.78	2/3445 (0.1%)	0.85	6/4678 (0.1%)
1	F	0.71	0/3441	0.82	3/4674 (0.1%)
All	All	0.79	6/20783 (0.0%)	0.86	44/28221 (0.2%)

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

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	55	GLU	CD-OE1	6.01	1.32	1.25
1	D	59	GLU	CD-OE2	5.62	1.31	1.25
1	D	197	GLU	CD-OE1	5.41	1.31	1.25
1	D	430	PHE	CG-CD1	-5.24	1.30	1.38
1	A	158	SER	CB-OG	5.20	1.49	1.42
1	E	59	GLU	CD-OE2	5.03	1.31	1.25

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	87	ASP	CB-CG-OD2	-8.82	110.36	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	136	ARG	NE-CZ-NH1	-8.45	116.07	120.30
1	A	87	ASP	CB-CG-OD2	8.05	125.55	118.30
1	C	136	ARG	NE-CZ-NH2	7.69	124.14	120.30
1	F	136	ARG	NE-CZ-NH2	7.41	124.01	120.30
1	E	136	ARG	NE-CZ-NH2	7.25	123.92	120.30
1	E	136	ARG	NE-CZ-NH1	-6.92	116.84	120.30
1	B	34	ASP	CB-CG-OD2	-6.89	112.10	118.30
1	E	358	ASP	CB-CG-OD1	6.83	124.45	118.30
1	B	136	ARG	NE-CZ-NH2	6.80	123.70	120.30
1	D	371	ARG	NE-CZ-NH2	-6.79	116.90	120.30
1	D	136	ARG	NE-CZ-NH2	6.79	123.69	120.30
1	A	136	ARG	NE-CZ-NH2	6.62	123.61	120.30
1	D	373	ASP	CB-CG-OD2	-6.60	112.36	118.30
1	B	330	ASP	CB-CG-OD2	6.52	124.17	118.30
1	A	358	ASP	CB-CG-OD1	6.48	124.13	118.30
1	C	321	ASP	CB-CG-OD1	6.36	124.03	118.30
1	C	321	ASP	CB-CG-OD2	-6.23	112.69	118.30
1	C	344	ARG	NE-CZ-NH2	-6.23	117.18	120.30
1	F	358	ASP	CB-CG-OD1	6.13	123.81	118.30
1	C	344	ARG	NE-CZ-NH1	6.06	123.33	120.30
1	C	358	ASP	CB-CG-OD1	6.04	123.73	118.30
1	A	140	ARG	NE-CZ-NH1	5.98	123.29	120.30
1	B	330	ASP	CB-CG-OD1	-5.96	112.94	118.30
1	D	87	ASP	CB-CG-OD1	5.82	123.54	118.30
1	B	358	ASP	CB-CG-OD1	5.74	123.47	118.30
1	C	258	LEU	CB-CG-CD2	5.71	120.71	111.00
1	C	212	ARG	NE-CZ-NH2	-5.68	117.46	120.30
1	C	87	ASP	CB-CG-OD1	5.66	123.39	118.30
1	B	314	ARG	NE-CZ-NH1	5.62	123.11	120.30
1	C	140	ARG	NE-CZ-NH2	5.61	123.10	120.30
1	D	321	ASP	CB-CG-OD1	5.60	123.34	118.30
1	B	87	ASP	CB-CG-OD1	5.54	123.28	118.30
1	E	413	VAL	CG1-CB-CG2	-5.52	102.07	110.90
1	A	327	PHE	CB-CG-CD1	5.41	124.59	120.80
1	B	87	ASP	CB-CG-OD2	-5.32	113.51	118.30
1	C	254	ASP	CB-CG-OD2	-5.27	113.56	118.30
1	C	197	GLU	OE1-CD-OE2	5.25	129.60	123.30
1	E	93	ARG	NE-CZ-NH2	-5.24	117.68	120.30
1	D	224	ASP	CB-CG-OD1	5.20	122.98	118.30
1	D	360	ARG	NE-CZ-NH2	-5.19	117.70	120.30
1	F	93	ARG	NE-CZ-NH2	-5.14	117.73	120.30
1	E	212	ARG	NE-CZ-NH2	-5.09	117.75	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	358	ASP	CB-CG-OD1	5.04	122.84	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	87	ASP	Sidechain

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	3426	0	3435	14	0
1	B	3408	0	3423	13	0
1	C	3367	0	3360	20	0
1	D	3367	0	3345	16	0
1	E	3364	0	3364	16	0
1	F	3360	0	3353	19	0
2	B	8	0	12	0	0
3	A	232	0	0	1	0
3	B	258	0	0	2	0
3	C	182	0	0	1	0
3	D	171	0	0	0	0
3	E	110	0	0	0	0
3	F	106	0	0	0	0
All	All	21359	0	20292	89	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 2.

All (89) 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:C:171:ILE:HG23	1:D:171:ILE:HG23	1.67	0.75
1:B:218[B]:PHE:CZ	1:B:425:TYR:HB3	2.23	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:354:LEU:HD12	1:F:354:LEU:O	1.92	0.69
1:E:171:ILE:HG23	1:F:171:ILE:HG23	1.74	0.69
1:F:350:GLY:CA	1:F:402:ASN:O	2.41	0.69
1:D:351:HIS:ND1	1:D:354:LEU:HD13	2.09	0.68
1:A:171:ILE:HG23	1:B:171:ILE:HG23	1.76	0.67
1:E:300:LEU:O	1:E:306:ALA:HB2	1.95	0.67
1:F:350:GLY:HA3	1:F:402:ASN:O	1.95	0.67
1:A:103:GLN:NE2	3:A:501:HOH:O	2.29	0.65
1:B:197:GLU:HB2	1:B:198:PRO:HD3	1.78	0.65
1:C:197:GLU:HB2	1:C:198:PRO:HD3	1.79	0.64
1:F:319:MET:SD	1:F:382:LEU:HD22	2.37	0.64
1:C:207:LEU:HD11	1:C:436:LEU:HD23	1.81	0.62
1:B:397:HIS:CD2	1:B:399:LYS:HG2	2.34	0.62
1:C:327:PHE:HE1	1:C:385:LYS:HD2	1.65	0.61
1:A:197:GLU:HB2	1:A:198:PRO:HD3	1.82	0.61
1:D:197:GLU:HB3	1:D:198:PRO:HD3	1.82	0.61
1:E:197:GLU:HB2	1:E:198:PRO:HD3	1.83	0.60
1:C:56:ARG:NH2	1:D:463:LEU:O	2.35	0.60
1:E:454:ARG:HH12	1:F:305:HIS:CD2	2.20	0.59
1:D:323:ILE:HD11	1:D:382:LEU:HD11	1.83	0.59
1:F:197:GLU:HB2	1:F:198:PRO:HD3	1.86	0.58
1:F:350:GLY:HA2	1:F:402:ASN:O	2.04	0.57
1:D:351:HIS:ND1	1:D:354:LEU:CD1	2.69	0.56
1:A:218[A]:PHE:CZ	1:A:425:TYR:HB3	2.41	0.55
1:D:319:MET:HG2	1:D:382:LEU:HD21	1.90	0.54
1:A:283:GLY:HA3	1:A:442:LEU:HD11	1.90	0.53
1:F:283:GLY:HA3	1:F:442:LEU:HD11	1.89	0.53
1:F:319:MET:O	1:F:323:ILE:HG12	2.09	0.52
1:E:356:LYS:HB2	1:E:357:PRO:HD2	1.92	0.52
1:A:30:THR:HG22	1:A:30:THR:O	2.10	0.51
1:F:323:ILE:HD11	1:F:382:LEU:HD21	1.93	0.51
1:B:30:THR:HB	3:C:529:HOH:O	2.10	0.51
1:C:283:GLY:HA3	1:C:442:LEU:HD11	1.93	0.50
1:E:283:GLY:HA3	1:E:442:LEU:HD11	1.93	0.50
1:A:385:LYS:HD2	1:A:389:ILE:HD11	1.94	0.49
1:D:283:GLY:HA3	1:D:442:LEU:HD11	1.94	0.49
1:C:160:PRO:HG3	1:D:151:HIS:CG	2.47	0.48
1:B:283:GLY:HA3	1:B:442:LEU:HD11	1.94	0.48
1:B:300:LEU:HD21	1:B:430:PHE:HD2	1.79	0.48
1:E:434:ARG:O	1:E:438:PRO:HD2	2.15	0.47
1:D:434:ARG:O	1:D:438:PRO:HD2	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:354:LEU:O	1:C:402:ASN:HB3	2.15	0.47
1:F:81:TRP:CE3	1:F:441:GLN:HG2	2.50	0.47
1:B:161:ARG:HA	1:B:161:ARG:HD3	1.52	0.46
1:E:454:ARG:NH1	1:F:305:HIS:CD2	2.83	0.46
1:F:416:TYR:CD1	1:F:421:GLN:HG2	2.50	0.46
1:A:97:LYS:HE2	1:A:105:GLU:OE2	2.15	0.46
1:C:416:TYR:CD1	1:C:421:GLN:HG2	2.51	0.46
1:E:81:TRP:CE3	1:E:441:GLN:HG2	2.51	0.46
1:E:310:GLN:HG3	1:E:423:PRO:HB2	1.98	0.46
1:F:434:ARG:O	1:F:438:PRO:HD2	2.15	0.46
1:B:434:ARG:O	1:B:438:PRO:HD2	2.15	0.45
1:F:212:ARG:HA	1:F:244:LEU:HD13	1.98	0.45
1:C:327:PHE:CZ	1:C:382:LEU:HD13	2.52	0.45
1:C:434:ARG:O	1:C:438:PRO:HD2	2.15	0.45
1:E:397:HIS:CE1	1:E:399:LYS:HB3	2.52	0.45
1:B:81:TRP:CE3	1:B:441:GLN:HG2	2.51	0.45
1:B:397:HIS:HE1	3:B:822:HOH:O	2.00	0.45
1:C:327:PHE:CE1	1:C:385:LYS:HD2	2.49	0.44
1:A:81:TRP:CE3	1:A:441:GLN:HG2	2.53	0.44
1:D:81:TRP:CE3	1:D:441:GLN:HG2	2.52	0.44
1:C:81:TRP:CE3	1:C:441:GLN:HG2	2.52	0.44
1:C:300:LEU:HD21	1:C:430:PHE:HD2	1.83	0.44
1:E:319:MET:O	1:E:323:ILE:HG12	2.17	0.44
1:A:32:GLU:HA	1:A:33:PRO:HD3	1.92	0.43
1:D:300:LEU:HD21	1:D:430:PHE:HD2	1.83	0.43
1:D:207:LEU:HD11	1:D:436:LEU:HD23	2.00	0.43
1:A:434:ARG:O	1:A:438:PRO:HD2	2.17	0.43
1:E:74:ARG:HH21	1:F:454:ARG:HD2	1.83	0.43
1:F:300:LEU:HD11	1:F:430:PHE:CD2	2.54	0.42
1:A:385:LYS:HD2	1:A:389:ILE:CD1	2.48	0.42
1:C:323:ILE:HD11	1:C:382:LEU:HD11	2.01	0.42
1:A:89:GLU:OE2	1:A:355:ARG:NE	2.46	0.42
1:C:300:LEU:HD11	1:C:430:PHE:CD2	2.55	0.42
1:B:65:LYS:HE3	3:B:824:HOH:O	2.19	0.42
1:D:212:ARG:HA	1:D:244:LEU:HD13	2.01	0.41
1:E:330:ASP:OD1	1:E:333:ARG:NH2	2.54	0.41
1:F:300:LEU:HD21	1:F:430:PHE:HD2	1.84	0.41
1:C:212:ARG:HA	1:C:244:LEU:HD13	2.03	0.41
1:E:255:PHE:CZ	1:E:420:PHE:CE1	3.09	0.41
1:C:91:GLY:HA2	1:C:354:LEU:HD21	2.03	0.41
1:C:155:LEU:HD11	1:D:155:LEU:HD11	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:300:LEU:HD21	1:A:430:PHE:HD2	1.86	0.40
1:D:32:GLU:HB3	1:D:33:PRO:HD2	2.03	0.40
1:B:212:ARG:HA	1:B:244:LEU:HD13	2.03	0.40
1:C:356:LYS:HB2	1:C:357:PRO:HD2	2.02	0.40
1:E:207:LEU:HD11	1:E:436:LEU:HD23	2.04	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	439/441 (100%)	435 (99%)	4 (1%)	0	100	100
1	B	435/441 (99%)	429 (99%)	6 (1%)	0	100	100
1	C	433/441 (98%)	428 (99%)	5 (1%)	0	100	100
1	D	434/441 (98%)	428 (99%)	6 (1%)	0	100	100
1	E	432/441 (98%)	427 (99%)	5 (1%)	0	100	100
1	F	432/441 (98%)	427 (99%)	5 (1%)	0	100	100
All	All	2605/2646 (98%)	2574 (99%)	31 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	364/365 (100%)	362 (100%)	2 (0%)	88 89
1	B	363/365 (100%)	361 (99%)	2 (1%)	86 87
1	C	356/365 (98%)	354 (99%)	2 (1%)	86 87
1	D	355/365 (97%)	354 (100%)	1 (0%)	92 93
1	E	356/365 (98%)	354 (99%)	2 (1%)	86 87
1	F	355/365 (97%)	353 (99%)	2 (1%)	86 87
All	All	2149/2190 (98%)	2138 (100%)	11 (0%)	88 89

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	177	ASN
1	A	321	ASP
1	B	161	ARG
1	B	177	ASN
1	C	144	GLU
1	C	177	ASN
1	D	177	ASN
1	E	177	ASN
1	E	307	LEU
1	F	177	ASN
1	F	354	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	F	305	HIS

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)

1 ligand is 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	TRS	B	501	-	7,7,7	0.95	0	9,9,9	3.79	7 (77%)

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
2	TRS	B	501	-	-	5/9/9/9	-

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	501	TRS	C2-C-N	-6.58	88.35	107.98
2	B	501	TRS	C1-C-N	-5.57	91.35	107.98
2	B	501	TRS	C3-C-N	-4.07	95.83	107.98
2	B	501	TRS	C2-C-C1	3.84	122.72	110.81
2	B	501	TRS	O3-C3-C	2.92	120.25	111.00
2	B	501	TRS	C3-C-C2	2.76	119.36	110.81
2	B	501	TRS	C3-C-C1	2.20	117.63	110.81

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	501	TRS	C2-C-C1-O1
2	B	501	TRS	C3-C-C2-O2
2	B	501	TRS	C1-C-C3-O3
2	B	501	TRS	C2-C-C3-O3
2	B	501	TRS	C1-C-C2-O2

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

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, 95^{th} 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	440/441 (99%)	-0.00	15 (3%) 45 49	29, 40, 63, 112	0
1	B	435/441 (98%)	-0.07	7 (1%) 72 74	28, 40, 61, 79	0
1	C	435/441 (98%)	-0.08	6 (1%) 75 78	29, 43, 72, 99	0
1	D	436/441 (98%)	0.25	32 (7%) 15 16	30, 46, 94, 121	0
1	E	434/441 (98%)	0.34	37 (8%) 10 11	32, 54, 92, 119	0
1	F	434/441 (98%)	0.23	20 (4%) 32 35	33, 54, 97, 118	0
All	All	2614/2646 (98%)	0.11	117 (4%) 33 35	28, 45, 84, 121	0

All (117) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	344	ARG	7.2
1	A	30	THR	6.1
1	D	340	LEU	5.7
1	D	400	THR	5.1
1	D	328	THR	4.9
1	D	332	VAL	4.9
1	E	325	THR	4.5
1	D	337	TRP	4.3
1	F	332	VAL	4.3
1	E	351	HIS	4.3
1	E	343	GLY	4.1
1	D	325	THR	4.0
1	E	346	VAL	4.0
1	F	325	THR	3.9
1	D	329	ASP	3.9
1	E	389	ILE	3.9
1	E	331	ASP	3.8
1	D	327	PHE	3.8
1	F	351	HIS	3.8

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Mol	Chain	Res	Type	RSRZ
1	D	401	LYS	3.7
1	F	400	THR	3.7
1	D	336	LEU	3.7
1	E	404	HIS	3.7
1	E	353	VAL	3.7
1	D	346	VAL	3.6
1	F	324	GLY	3.6
1	E	330	ASP	3.5
1	E	334	ASN	3.4
1	D	330	ASP	3.4
1	D	353	VAL	3.4
1	D	319	MET	3.3
1	C	352	GLY	3.3
1	F	31	ALA	3.3
1	D	335	TYR	3.3
1	D	333	ARG	3.3
1	D	341	LYS	3.2
1	A	297	LEU	3.2
1	E	300	LEU	3.2
1	C	30	THR	3.2
1	D	31	ALA	3.1
1	E	304	LEU	3.1
1	E	322	LYS	3.1
1	E	338	ASP	3.1
1	A	31	ALA	3.1
1	D	351	HIS	3.0
1	E	393	VAL	3.0
1	F	420	PHE	3.0
1	E	335	TYR	3.0
1	E	400	THR	3.0
1	A	29	SER	2.9
1	A	26	GLY	2.9
1	A	282	VAL	2.9
1	D	30	THR	2.9
1	F	342	SER	2.9
1	D	345	VAL	2.9
1	F	321	ASP	2.9
1	A	300	LEU	2.9
1	D	387	SER	2.8
1	B	297	LEU	2.8
1	D	297	LEU	2.8
1	E	344	ARG	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	394	LEU	2.8
1	A	345	VAL	2.8
1	A	436	LEU	2.7
1	E	90	GLU	2.7
1	E	274	VAL	2.6
1	D	315	TRP	2.6
1	E	403	PRO	2.6
1	C	325	THR	2.6
1	D	392	ALA	2.6
1	E	298	LEU	2.6
1	A	27	SER	2.6
1	F	191	SER	2.6
1	B	298	LEU	2.5
1	E	317	LEU	2.5
1	E	314	ARG	2.5
1	A	432	VAL	2.5
1	F	327	PHE	2.5
1	B	278	ALA	2.4
1	F	353	VAL	2.4
1	F	394	LEU	2.4
1	E	324	GLY	2.4
1	F	338	ASP	2.4
1	A	276	ALA	2.4
1	E	378	PRO	2.3
1	D	305	HIS	2.3
1	B	296	GLY	2.3
1	C	31	ALA	2.3
1	E	313	LEU	2.3
1	A	400	THR	2.3
1	C	305	HIS	2.3
1	E	385	LYS	2.3
1	E	297	LEU	2.3
1	F	74	ARG	2.3
1	E	64	VAL	2.2
1	E	356	LYS	2.2
1	E	463	LEU	2.2
1	E	327	PHE	2.2
1	E	332	VAL	2.2
1	B	301	ALA	2.2
1	E	382	LEU	2.2
1	E	361	PHE	2.2
1	A	275	SER	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	325	THR	2.2
1	A	279	THR	2.1
1	F	323	ILE	2.1
1	C	298	LEU	2.1
1	D	354	LEU	2.1
1	D	381	GLN	2.1
1	F	291	LEU	2.1
1	D	334	ASN	2.0
1	F	317	LEU	2.0
1	E	89	GLU	2.0
1	B	321	ASP	2.0
1	D	391	PRO	2.0
1	F	326	LYS	2.0
1	F	76	LEU	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 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, 95^{th} 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	TRS	B	501	8/8	0.94	0.10	41,44,46,47	0

6.5 Other polymers (i)

There are no such residues in this entry.