



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 27, 2026 – 05:47 PM EDT

PDB ID : 5MVI / pdb_00005mvi
Title : Crystal structure of 2-methylcitrate dehydratase (PrpD) from *Salmonella enterica*
Authors : Race, P.R.; Baker, G.E.
Deposited on : 2017-01-16
Resolution : 3.05 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	:	4-5-2 with Phenix2.0
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

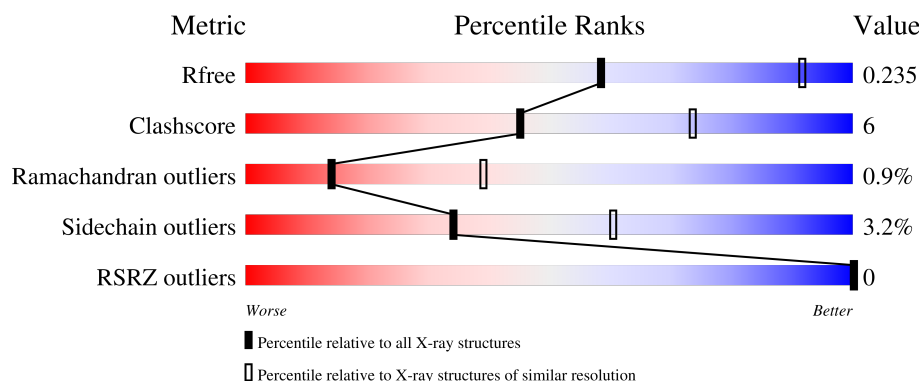
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.05 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}	180053	2469 (3.10-3.02)
Clashscore	190562	2569 (3.10-3.02)
Ramachandran outliers	187476	2424 (3.10-3.02)
Sidechain outliers	187428	2423 (3.10-3.02)
RSRZ outliers	180081	2469 (3.10-3.02)

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	483	
1	B	483	
1	C	483	
1	D	483	
1	E	483	

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Mol	Chain	Length	Quality of chain
1	F	483	 A horizontal bar chart showing the quality of chain F. The bar is divided into three segments: a green segment representing 62%, a yellow segment representing 6%, and a grey segment representing 31%. A small black dot is located at the end of the yellow segment.

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 15106 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 dehydratase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	E	334	Total	C	N	O	S	0	0	0
			2341	1491	410	428	12			
1	D	337	Total	C	N	O	S	0	0	0
			2406	1534	417	444	11			
1	A	463	Total	C	N	O	S	0	0	0
			3283	2100	566	604	13			
1	B	334	Total	C	N	O	S	0	0	0
			2359	1508	409	433	9			
1	C	332	Total	C	N	O	S	0	0	0
			2375	1517	409	437	12			
1	F	334	Total	C	N	O	S	0	0	0
			2342	1493	407	429	13			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	8	VAL	ILE	conflict	UNP A0A0L5RC52
E	123	THR	ILE	conflict	UNP A0A0L5RC52
E	418	GLY	ASP	conflict	UNP A0A0L5RC52
D	8	VAL	ILE	conflict	UNP A0A0L5RC52
D	123	THR	ILE	conflict	UNP A0A0L5RC52
D	418	GLY	ASP	conflict	UNP A0A0L5RC52
A	8	VAL	ILE	conflict	UNP A0A0L5RC52
A	123	THR	ILE	conflict	UNP A0A0L5RC52
A	418	GLY	ASP	conflict	UNP A0A0L5RC52
B	8	VAL	ILE	conflict	UNP A0A0L5RC52
B	123	THR	ILE	conflict	UNP A0A0L5RC52
B	418	GLY	ASP	conflict	UNP A0A0L5RC52
C	8	VAL	ILE	conflict	UNP A0A0L5RC52
C	123	THR	ILE	conflict	UNP A0A0L5RC52
C	418	GLY	ASP	conflict	UNP A0A0L5RC52
F	8	VAL	ILE	conflict	UNP A0A0L5RC52
F	123	THR	ILE	conflict	UNP A0A0L5RC52

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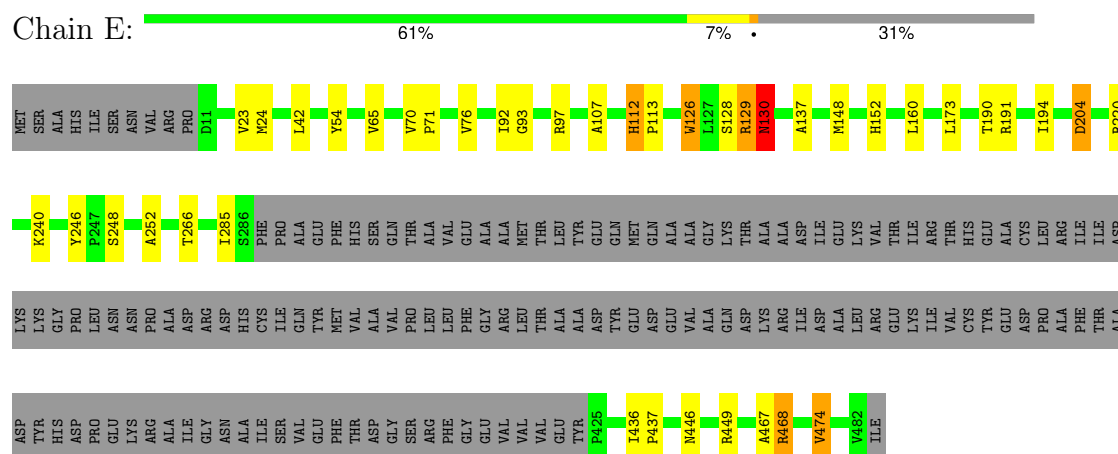
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Chain	Residue	Modelled	Actual	Comment	Reference
F	418	GLY	ASP	conflict	UNP A0A0L5RC52

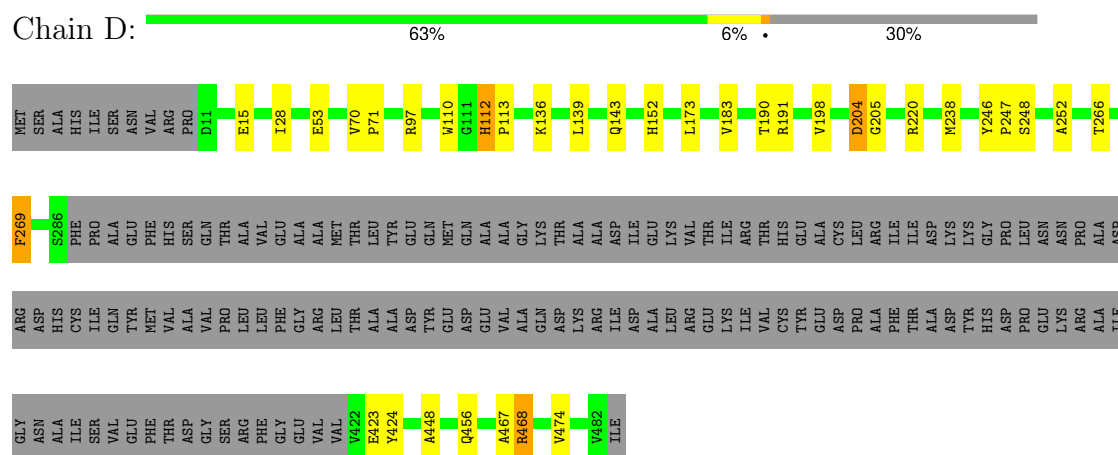
3 Residue-property plots

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.

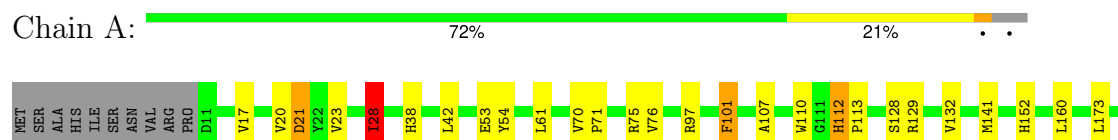
• Molecule 1: 2-methylcitrate dehydratase

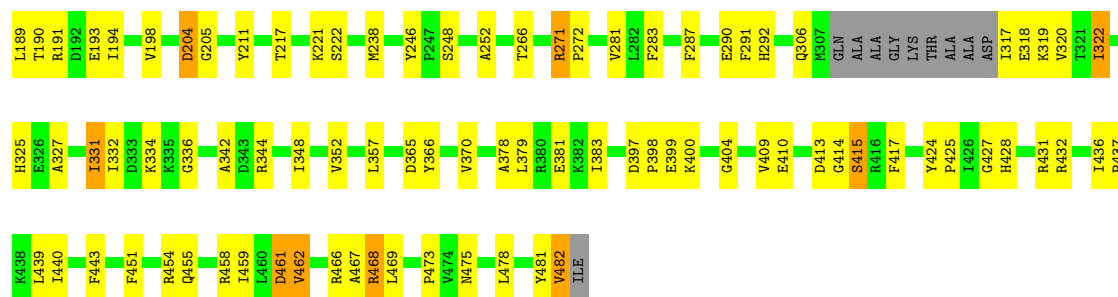


• Molecule 1: 2-methylcitrate dehydratase

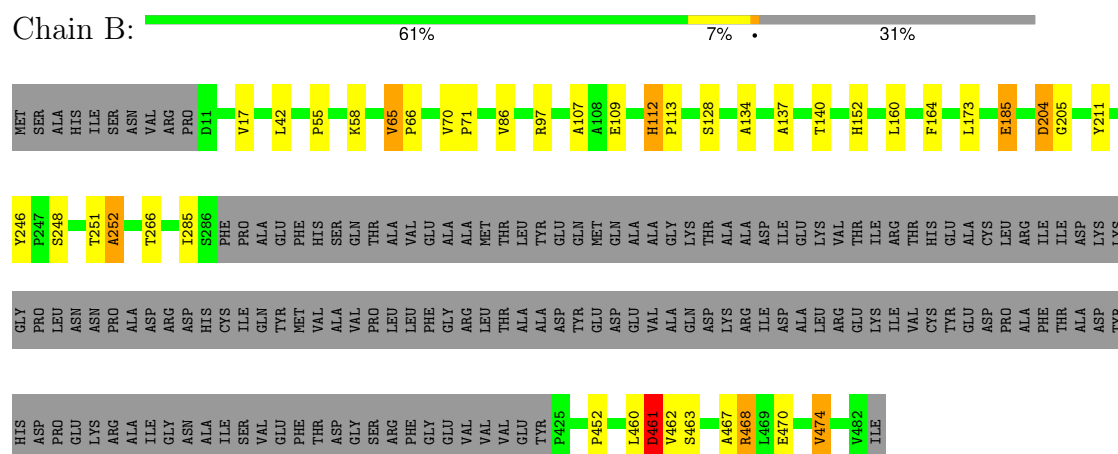


• Molecule 1: 2-methylcitrate dehydratase

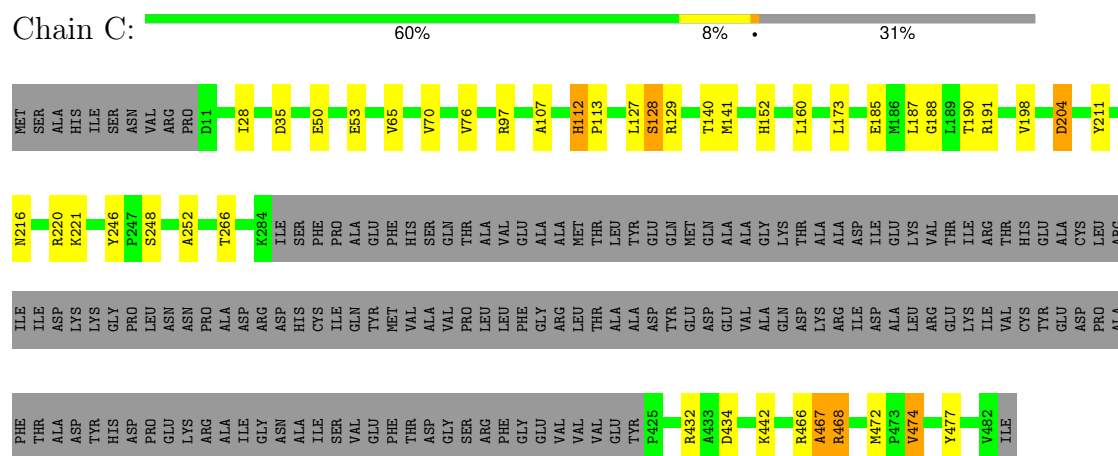




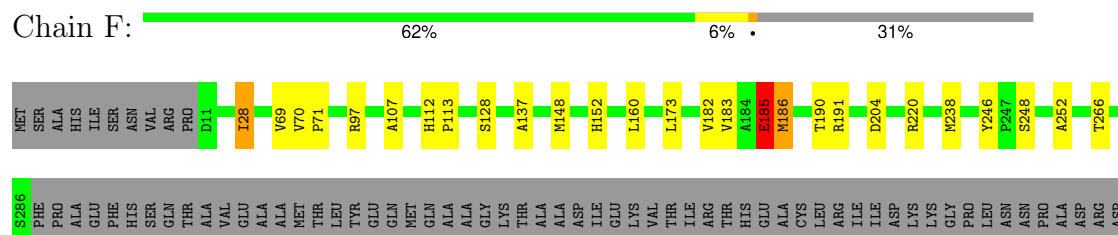
- Molecule 1: 2-methylcitrate dehydratase



- Molecule 1: 2-methylcitrate dehydratase



- Molecule 1: 2-methylcitrate dehydratase



HIS CYS ILE GLN TYR MET VAL ALA VAL PRO LEU PHE GLY ARG LEU THR ALA ALA ASP TYR GLU ASP VAL VAL ALA GLN ASP LYS ARG TLE ASP ALA LEU ARG GLU LYS TLE VAL CYS TYR ASP PRO ALA PHE THR ALA ASP TYR HIS ASP PRO GLU LYS ARG ALA TLE GLY ASN

ALA ILE SER VAL GLU PHE THR ASP GLY SER PHE GLY VAL VAL VAL GLU TYR P425 M446 R449 T453 A467 R468 V474 Y482 ILE

4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	139.66Å 139.66Å 513.48Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	39.14 – 3.05 39.14 – 3.05	Depositor EDS
% Data completeness (in resolution range)	99.5 (39.14-3.05) 99.6 (39.14-3.05)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.71 (at 3.06Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.213 , 0.239 0.211 , 0.235	Depositor DCC
R_{free} test set	3483 reflections (4.89%)	wwPDB-VP
Wilson B-factor (Å ²)	25.5	Xtriage
Anisotropy	0.373	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 30.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.36$, $\langle L^2 \rangle = 0.19$	Xtriage
Estimated twinning fraction	0.387 for -h-k,k,-l	Xtriage
Reported twinning fraction	0.532 for H, K, L 0.468 for K, H, -L	Depositor
Outliers	0 of 70839 reflections	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	15106	wwPDB-VP
Average B, all atoms (Å ²)	9.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.65% 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

5.1 Standard geometry

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	1.12	8/3352 (0.2%)	1.13	19/4559 (0.4%)
1	B	1.05	5/2406 (0.2%)	1.06	10/3272 (0.3%)
1	C	0.97	3/2424 (0.1%)	1.05	14/3293 (0.4%)
1	D	0.99	1/2454 (0.0%)	1.05	5/3333 (0.2%)
1	E	0.97	0/2385	1.09	13/3239 (0.4%)
1	F	0.87	1/2385 (0.0%)	0.99	7/3238 (0.2%)
All	All	1.01	18/15406 (0.1%)	1.07	68/20934 (0.3%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	4
1	B	0	1
1	D	0	1
1	E	0	2
1	F	0	1
All	All	0	9

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	473	PRO	CA-C	8.55	1.62	1.52
1	A	478	LEU	CA-C	7.80	1.64	1.52
1	A	481	TYR	CB-CG	7.60	1.68	1.51
1	D	183	VAL	CA-CB	6.77	1.63	1.54
1	A	440	ILE	CB-CG1	6.75	1.67	1.53
1	C	128	SER	CA-C	6.65	1.61	1.52
1	C	140	THR	CA-C	6.05	1.59	1.52
1	F	69	VAL	N-CA	5.91	1.53	1.46
1	B	251	THR	CA-CB	5.91	1.64	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	194	ILE	CA-CB	5.55	1.61	1.54
1	B	252	ALA	N-CA	5.51	1.52	1.46
1	B	137	ALA	CA-C	5.47	1.58	1.52
1	C	128	SER	C-O	-5.39	1.17	1.24
1	B	462	VAL	N-CA	5.33	1.53	1.46
1	A	451	PHE	CA-C	5.18	1.58	1.53
1	A	194	ILE	CA-C	5.13	1.59	1.52
1	A	439	LEU	N-CA	-5.07	1.39	1.46
1	B	86	VAL	N-CA	5.03	1.52	1.46

All (68) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	129	ARG	N-CA-C	12.73	125.16	111.03
1	A	462	VAL	N-CA-C	-9.61	100.52	111.00
1	A	271	ARG	N-CA-C	8.00	115.19	108.07
1	A	336	GLY	CA-C-N	7.41	127.46	119.90
1	A	336	GLY	C-N-CA	7.41	127.46	119.90
1	E	65	VAL	CA-C-N	-6.85	112.72	119.78
1	E	65	VAL	C-N-CA	-6.85	112.72	119.78
1	A	132	VAL	N-CA-C	-6.66	104.02	110.42
1	A	473	PRO	N-CA-C	6.51	121.06	111.03
1	E	137	ALA	CA-C-N	-6.38	111.86	119.84
1	E	137	ALA	C-N-CA	-6.38	111.86	119.84
1	C	129	ARG	N-CA-C	-6.34	104.24	112.23
1	D	112	HIS	CA-C-N	6.32	126.01	119.56
1	D	112	HIS	C-N-CA	6.32	126.01	119.56
1	F	69	VAL	N-CA-C	6.26	116.64	107.37
1	E	126	TRP	N-CA-C	6.26	118.66	111.02
1	A	331	ILE	N-CA-C	6.25	116.42	110.74
1	D	28	ILE	CB-CA-C	-6.18	102.05	111.69
1	E	194	ILE	N-CA-CB	6.17	119.84	110.58
1	C	28	ILE	CG1-CB-CG2	-6.16	92.22	110.70
1	F	453	THR	CB-CA-C	-6.14	100.65	110.84
1	A	28	ILE	CB-CA-C	-6.12	103.14	111.33
1	E	130	ASN	N-CA-C	6.10	123.80	110.80
1	B	461	ASP	N-CA-C	-5.96	98.11	110.80
1	F	185	GLU	N-CA-C	5.81	123.18	110.80
1	A	461	ASP	N-CA-C	5.79	117.59	111.28
1	C	442	LYS	N-CA-C	5.76	118.03	111.11
1	A	271	ARG	N-CA-CB	-5.74	105.38	111.29
1	C	112	HIS	CA-C-N	5.69	126.07	119.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	112	HIS	C-N-CA	5.69	126.07	119.47
1	A	198	VAL	CB-CA-C	-5.65	104.42	112.22
1	E	112	HIS	CA-C-N	5.64	126.01	119.47
1	E	112	HIS	C-N-CA	5.64	126.01	119.47
1	C	472	MET	CA-C-N	5.63	125.91	119.83
1	C	472	MET	C-N-CA	5.63	125.91	119.83
1	D	198	VAL	CB-CA-C	-5.62	104.56	112.14
1	A	61	LEU	N-CA-C	5.61	118.75	109.94
1	A	383	ILE	N-CA-C	5.58	116.16	108.12
1	C	140	THR	CB-CA-C	5.56	119.55	109.37
1	B	17	VAL	CB-CA-C	-5.53	104.67	112.14
1	F	112	HIS	CA-C-N	5.53	125.89	119.47
1	F	112	HIS	C-N-CA	5.53	125.89	119.47
1	C	198	VAL	CB-CA-C	-5.46	104.68	112.22
1	A	112	HIS	CA-C-N	5.38	125.71	119.47
1	A	112	HIS	C-N-CA	5.38	125.71	119.47
1	C	65	VAL	CA-C-N	-5.36	114.24	119.76
1	C	65	VAL	C-N-CA	-5.36	114.24	119.76
1	B	112	HIS	CA-C-N	5.34	125.66	119.47
1	B	112	HIS	C-N-CA	5.34	125.66	119.47
1	B	452	PRO	CA-C-N	5.31	127.71	120.54
1	B	452	PRO	C-N-CA	5.31	127.71	120.54
1	F	137	ALA	CA-C-N	-5.25	115.36	120.98
1	F	137	ALA	C-N-CA	-5.25	115.36	120.98
1	C	70	VAL	N-CA-C	5.23	113.19	107.60
1	A	320	VAL	N-CA-C	5.20	115.58	107.99
1	E	76	VAL	CB-CA-C	-5.18	104.89	110.85
1	D	136	LYS	N-CA-C	-5.12	102.57	109.95
1	A	76	VAL	N-CA-C	5.08	113.52	107.73
1	B	65	VAL	CA-C-N	-5.06	114.55	119.76
1	B	65	VAL	C-N-CA	-5.06	114.55	119.76
1	A	322	ILE	N-CA-C	5.05	115.18	108.11
1	C	467	ALA	N-CA-C	-5.04	103.87	110.53
1	E	54	TYR	CA-C-N	5.03	124.69	119.56
1	E	54	TYR	C-N-CA	5.03	124.69	119.56
1	C	76	VAL	CB-CA-C	-5.01	105.09	110.85
1	A	459	ILE	CB-CA-C	-5.00	105.56	111.97
1	B	134	ALA	N-CA-C	-5.00	106.77	112.92
1	B	185	GLU	N-CA-C	-5.00	105.74	111.14

There are no chirality outliers.

All (9) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	317	ILE	Peptide
1	A	318	GLU	Peptide
1	A	331	ILE	Peptide
1	A	462	VAL	Peptide
1	B	461	ASP	Mainchain
1	D	424	TYR	Peptide
1	E	129	ARG	Peptide
1	E	285	ILE	Peptide
1	F	185	GLU	Peptide

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3283	0	2890	70	0
1	B	2359	0	2114	25	0
1	C	2375	0	2152	26	0
1	D	2406	0	2172	18	0
1	E	2341	0	2098	22	0
1	F	2342	0	2113	20	0
All	All	15106	0	13539	169	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 6.

All (169) 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:E:24:MET:HE1	1:E:191:ARG:HA	1.08	1.04
1:B:460:LEU:O	1:B:461:ASP:O	1.81	0.98
1:C:141:MET:HE2	1:C:474:VAL:HA	1.45	0.98
1:E:24:MET:HE1	1:E:191:ARG:CA	2.00	0.90
1:E:24:MET:CE	1:E:191:ARG:HA	2.02	0.86
1:C:141:MET:HE3	1:C:477:TYR:HB3	1.55	0.86
1:A:38:HIS:ND1	1:A:466:ARG:CG	2.45	0.80
1:A:283:PHE:O	1:A:427:GLY:HA3	1.83	0.79
1:A:291:PHE:O	1:A:291:PHE:HD1	1.66	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:141:MET:HE3	1:C:477:TYR:CB	2.17	0.73
1:F:182:VAL:O	1:F:186:MET:HB2	1.89	0.73
1:D:246:TYR:OH	1:C:220:ARG:HB2	1.90	0.71
1:A:425:PRO:HD2	1:A:428:HIS:HB2	1.73	0.71
1:A:204:ASP:OD1	1:A:205:GLY:N	2.23	0.70
1:B:140:THR:HG22	1:B:470:GLU:O	1.92	0.69
1:A:319:LYS:HB3	1:A:410:GLU:HB2	1.74	0.69
1:F:183:VAL:HA	1:F:186:MET:HG2	1.74	0.69
1:A:291:PHE:O	1:A:291:PHE:CD1	2.46	0.68
1:C:35:ASP:OD1	1:C:466:ARG:NH1	2.26	0.68
1:C:141:MET:HE2	1:C:474:VAL:CA	2.23	0.68
1:F:183:VAL:O	1:F:186:MET:HB3	1.94	0.67
1:A:467:ALA:O	1:A:468:ARG:HB3	1.94	0.67
1:E:126:TRP:O	1:E:130:ASN:HB2	1.94	0.67
1:D:204:ASP:OD1	1:D:205:GLY:N	2.26	0.66
1:A:97:ARG:O	1:A:97:ARG:HG3	1.96	0.65
1:B:140:THR:CG2	1:B:470:GLU:O	2.45	0.65
1:D:269:PHE:N	1:D:269:PHE:CD2	2.65	0.62
1:C:97:ARG:O	1:C:97:ARG:HG3	2.00	0.62
1:B:204:ASP:OD1	1:B:205:GLY:N	2.29	0.62
1:E:97:ARG:O	1:E:97:ARG:HG3	2.00	0.61
1:A:322:ILE:HD11	1:A:352:VAL:HG11	1.81	0.61
1:F:97:ARG:O	1:F:97:ARG:HG3	1.99	0.61
1:A:467:ALA:O	1:A:468:ARG:CB	2.49	0.61
1:D:97:ARG:O	1:D:97:ARG:HG3	2.01	0.61
1:A:20:VAL:O	1:A:21:ASP:CG	2.44	0.60
1:A:325:HIS:HD2	1:A:404:GLY:O	1.85	0.60
1:B:97:ARG:O	1:B:97:ARG:HG3	2.00	0.60
1:A:378:ALA:O	1:A:381:GLU:CG	2.49	0.60
1:A:281:VAL:O	1:A:432:ARG:NH2	2.35	0.59
1:A:38:HIS:CE1	1:A:466:ARG:CG	2.85	0.59
1:D:246:TYR:OH	1:C:220:ARG:CB	2.50	0.59
1:E:220:ARG:HB2	1:F:246:TYR:OH	2.03	0.59
1:A:365:ASP:HA	1:A:370:VAL:HG21	1.85	0.58
1:B:460:LEU:C	1:B:461:ASP:O	2.44	0.58
1:E:126:TRP:O	1:E:130:ASN:ND2	2.31	0.58
1:A:332:ILE:HD12	1:A:348:ILE:CB	2.33	0.58
1:F:182:VAL:O	1:F:186:MET:CB	2.52	0.57
1:A:217:THR:CG2	1:A:342:ALA:O	2.52	0.57
1:A:221:LYS:C	1:A:221:LYS:HD2	2.30	0.57
1:A:344:ARG:NH1	1:A:366:TYR:O	2.38	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:467:ALA:O	1:F:468:ARG:CB	2.53	0.57
1:D:467:ALA:O	1:D:468:ARG:CB	2.52	0.56
1:C:467:ALA:O	1:C:468:ARG:CB	2.53	0.56
1:E:467:ALA:O	1:E:468:ARG:CB	2.53	0.56
1:B:467:ALA:O	1:B:468:ARG:CB	2.53	0.55
1:A:458:ARG:HA	1:A:461:ASP:HB2	1.88	0.55
1:A:221:LYS:HG2	1:A:342:ALA:HB1	1.89	0.54
1:A:113:PRO:HB2	1:A:152:HIS:CG	2.42	0.54
1:A:20:VAL:O	1:A:21:ASP:CB	2.54	0.54
1:A:211:TYR:HB3	1:B:246:TYR:CE1	2.44	0.53
1:A:379:LEU:O	1:A:379:LEU:HD12	2.09	0.53
1:A:189:LEU:HD23	1:A:193:GLU:HB3	1.90	0.53
1:A:400:LYS:HE2	1:A:424:TYR:OH	2.09	0.52
1:A:42:LEU:HD13	1:A:443:PHE:CD2	2.45	0.52
1:A:75:ARG:CG	1:A:129:ARG:NH1	2.72	0.52
1:C:50:GLU:O	1:C:53:GLU:HG3	2.10	0.52
1:E:113:PRO:HB2	1:E:152:HIS:CE1	2.45	0.51
1:A:409:VAL:HB	1:A:417:PHE:CG	2.46	0.51
1:E:246:TYR:OH	1:F:220:ARG:HB2	2.11	0.50
1:E:113:PRO:HB2	1:E:152:HIS:CG	2.47	0.49
1:A:455:GLN:HE22	1:A:482:VAL:C	2.20	0.49
1:D:220:ARG:HB2	1:C:246:TYR:OH	2.12	0.49
1:A:357:LEU:HG	1:A:379:LEU:HD23	1.95	0.49
1:D:246:TYR:CE1	1:C:211:TYR:HB3	2.48	0.49
1:A:221:LYS:C	1:A:221:LYS:CD	2.85	0.49
1:F:446:ASN:O	1:F:449:ARG:HG2	2.12	0.49
1:A:414:GLY:O	1:A:415:SER:O	2.31	0.48
1:C:141:MET:HE3	1:C:477:TYR:CG	2.47	0.48
1:A:413:ASP:C	1:A:415:SER:H	2.21	0.48
1:A:217:THR:HG21	1:A:342:ALA:O	2.14	0.48
1:A:17:VAL:O	1:A:20:VAL:O	2.31	0.48
1:D:113:PRO:HB2	1:D:152:HIS:CE1	2.47	0.48
1:F:113:PRO:HB2	1:F:152:HIS:CG	2.48	0.48
1:C:141:MET:CE	1:C:477:TYR:HB3	2.35	0.48
1:B:113:PRO:HB2	1:B:152:HIS:CG	2.49	0.47
1:A:436:ILE:N	1:A:437:PRO:HD2	2.29	0.47
1:A:101:PHE:CE1	1:A:287:PHE:O	2.67	0.47
1:B:285:ILE:O	1:B:285:ILE:CG1	2.63	0.47
1:A:204:ASP:CG	1:A:205:GLY:H	2.21	0.47
1:B:248:SER:O	1:B:252:ALA:HB2	2.14	0.47
1:C:221:LYS:HD2	1:C:221:LYS:C	2.39	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:397:ASP:HA	1:A:398:PRO:HD3	1.69	0.47
1:A:399:GLU:O	1:A:399:GLU:CG	2.63	0.47
1:D:113:PRO:HB2	1:D:152:HIS:CG	2.49	0.47
1:E:446:ASN:O	1:E:449:ARG:HG2	2.15	0.46
1:A:38:HIS:CG	1:A:466:ARG:CG	2.98	0.46
1:A:42:LEU:HB2	1:A:443:PHE:CZ	2.51	0.46
1:A:283:PHE:O	1:A:427:GLY:CA	2.60	0.46
1:C:248:SER:O	1:C:252:ALA:HB2	2.16	0.46
1:E:126:TRP:O	1:E:130:ASN:CB	2.64	0.45
1:D:139:LEU:CD2	1:D:143:GLN:OE1	2.64	0.45
1:B:55:PRO:HA	1:B:58:LYS:HE2	1.98	0.45
1:C:127:LEU:HD21	1:C:187:LEU:HD23	1.97	0.45
1:B:70:VAL:HA	1:B:71:PRO:HD2	1.82	0.45
1:A:128:SER:HB3	1:A:475:ASN:OD1	2.17	0.45
1:A:141:MET:HB2	1:A:469:LEU:O	2.16	0.45
1:F:70:VAL:HA	1:F:71:PRO:HD2	1.83	0.45
1:C:113:PRO:HB2	1:C:152:HIS:CG	2.52	0.45
1:E:190:THR:O	1:E:191:ARG:C	2.59	0.45
1:E:248:SER:O	1:E:252:ALA:HB2	2.17	0.45
1:F:248:SER:O	1:F:252:ALA:HB2	2.16	0.45
1:B:467:ALA:O	1:B:468:ARG:HB3	2.17	0.45
1:B:113:PRO:HB2	1:B:152:HIS:CE1	2.53	0.44
1:A:211:TYR:HB3	1:B:246:TYR:CZ	2.52	0.44
1:F:113:PRO:HB2	1:F:152:HIS:CE1	2.53	0.44
1:E:220:ARG:CB	1:F:246:TYR:OH	2.66	0.44
1:A:248:SER:O	1:A:252:ALA:HB2	2.18	0.44
1:D:70:VAL:HA	1:D:71:PRO:HD2	1.78	0.43
1:A:290:GLU:OE1	1:A:292:HIS:HB2	2.18	0.43
1:A:413:ASP:O	1:A:415:SER:N	2.42	0.43
1:C:113:PRO:HB2	1:C:152:HIS:CE1	2.53	0.43
1:E:70:VAL:HA	1:E:71:PRO:HD2	1.86	0.43
1:A:113:PRO:HB2	1:A:152:HIS:CE1	2.52	0.43
1:C:221:LYS:C	1:C:221:LYS:CD	2.92	0.43
1:D:190:THR:HG22	1:D:191:ARG:N	2.33	0.43
1:D:248:SER:O	1:D:252:ALA:HB2	2.19	0.43
1:F:185:GLU:CG	1:F:186:MET:N	2.81	0.43
1:A:217:THR:HG22	1:A:342:ALA:O	2.19	0.43
1:A:246:TYR:CE1	1:B:211:TYR:HB3	2.53	0.43
1:B:112:HIS:HA	1:B:113:PRO:HD2	1.83	0.43
1:E:112:HIS:HA	1:E:113:PRO:HD2	1.81	0.43
1:A:425:PRO:O	1:A:431:ARG:HD3	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:190:THR:O	1:C:191:ARG:C	2.62	0.42
1:F:190:THR:O	1:F:191:ARG:C	2.62	0.42
1:C:112:HIS:HA	1:C:113:PRO:HD2	1.81	0.42
1:A:190:THR:HG22	1:A:191:ARG:N	2.33	0.42
1:F:183:VAL:C	1:F:186:MET:HB3	2.44	0.42
1:E:436:ILE:N	1:E:437:PRO:CD	2.82	0.42
1:A:211:TYR:CB	1:B:246:TYR:CE1	3.02	0.42
1:E:128:SER:OG	1:E:474:VAL:HG22	2.20	0.42
1:A:53:GLU:HB2	1:A:54:TYR:CD1	2.54	0.42
1:D:448:ALA:HA	1:D:456:GLN:CD	2.44	0.42
1:D:15:GLU:CG	1:D:269:PHE:HZ	2.33	0.42
1:B:461:ASP:O	1:B:463:SER:N	2.53	0.42
1:A:221:LYS:HD2	1:A:222:SER:N	2.35	0.42
1:A:271:ARG:HB2	1:A:272:PRO:HD2	2.02	0.42
1:A:357:LEU:HG	1:A:379:LEU:CD2	2.49	0.42
1:C:128:SER:OG	1:C:474:VAL:HG22	2.19	0.42
1:D:247:PRO:HD2	1:C:216:ASN:OD1	2.20	0.41
1:A:28:ILE:H	1:A:28:ILE:HG12	1.70	0.41
1:B:107:ALA:HB1	1:B:160:LEU:HA	2.03	0.41
1:A:70:VAL:HA	1:A:71:PRO:HD2	1.83	0.41
1:B:58:LYS:HE2	1:B:58:LYS:HB2	1.90	0.41
1:E:92:ILE:O	1:E:93:GLY:C	2.63	0.41
1:A:292:HIS:CD2	1:A:327:ALA:HB3	2.54	0.41
1:B:65:VAL:O	1:B:66:PRO:C	2.63	0.41
1:C:107:ALA:HB1	1:C:160:LEU:HA	2.02	0.41
1:A:107:ALA:HB1	1:A:160:LEU:HA	2.03	0.41
1:A:424:TYR:HA	1:A:425:PRO:HD3	1.90	0.41
1:A:454:ARG:HD2	1:A:454:ARG:O	2.20	0.41
1:C:185:GLU:O	1:C:188:GLY:N	2.47	0.41
1:A:112:HIS:HA	1:A:113:PRO:HD2	1.84	0.41
1:B:128:SER:OG	1:B:474:VAL:HG22	2.21	0.41
1:B:109:GLU:OE2	1:B:164:PHE:N	2.50	0.41
1:F:107:ALA:HB1	1:F:160:LEU:HA	2.03	0.40
1:E:107:ALA:HB1	1:E:160:LEU:HA	2.03	0.40
1:F:28:ILE:H	1:F:28:ILE:HG12	1.69	0.40
1:D:112:HIS:HA	1:D:113:PRO:HD2	1.83	0.40
1:F:128:SER:OG	1:F:474:VAL:HG22	2.21	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	459/483 (95%)	431 (94%)	23 (5%)	5 (1%)	11	34
1	B	330/483 (68%)	319 (97%)	8 (2%)	3 (1%)	14	39
1	C	328/483 (68%)	319 (97%)	7 (2%)	2 (1%)	21	48
1	D	333/483 (69%)	323 (97%)	7 (2%)	3 (1%)	14	39
1	E	330/483 (68%)	317 (96%)	10 (3%)	3 (1%)	14	39
1	F	330/483 (68%)	320 (97%)	6 (2%)	4 (1%)	10	32
All	All	2110/2898 (73%)	2029 (96%)	61 (3%)	20 (1%)	14	39

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	21	ASP
1	A	204	ASP
1	A	415	SER
1	B	461	ASP
1	C	204	ASP
1	E	204	ASP
1	D	204	ASP
1	A	334	LYS
1	A	468	ARG
1	F	204	ASP
1	E	130	ASN
1	D	468	ARG
1	B	204	ASP
1	C	468	ARG
1	F	186	MET
1	F	468	ARG
1	E	468	ARG
1	D	423	GLU
1	B	468	ARG
1	F	185	GLU

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	270/395 (68%)	261 (97%)	9 (3%)	33	60
1	B	196/395 (50%)	191 (97%)	5 (3%)	40	65
1	C	207/395 (52%)	201 (97%)	6 (3%)	37	62
1	D	207/395 (52%)	200 (97%)	7 (3%)	32	59
1	E	194/395 (49%)	186 (96%)	8 (4%)	27	55
1	F	196/395 (50%)	190 (97%)	6 (3%)	35	61
All	All	1270/2370 (54%)	1229 (97%)	41 (3%)	34	60

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	23	VAL
1	E	42	LEU
1	E	148	MET
1	E	173	LEU
1	E	204	ASP
1	E	240	LYS
1	E	266	THR
1	E	474	VAL
1	D	53	GLU
1	D	110	TRP
1	D	173	LEU
1	D	238	MET
1	D	266	THR
1	D	269	PHE
1	D	474	VAL
1	A	23	VAL
1	A	28	ILE
1	A	101	PHE
1	A	110	TRP
1	A	173	LEU
1	A	238	MET
1	A	266	THR

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Mol	Chain	Res	Type
1	A	306	GLN
1	A	482	VAL
1	B	42	LEU
1	B	173	LEU
1	B	185	GLU
1	B	266	THR
1	B	474	VAL
1	C	173	LEU
1	C	204	ASP
1	C	266	THR
1	C	432	ARG
1	C	434	ASP
1	C	474	VAL
1	F	28	ILE
1	F	148	MET
1	F	173	LEU
1	F	238	MET
1	F	266	THR
1	F	474	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	165	ASN
1	E	280	ASN
1	E	456	GLN
1	D	280	ASN
1	D	456	GLN
1	A	292	HIS
1	A	306	GLN
1	A	325	HIS
1	A	455	GLN
1	C	165	ASN
1	C	455	GLN
1	F	165	ASN

5.3.3 RNA ⓘ

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](#)

There are no ligands in this entry.

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	463/483 (95%)	-1.73	0 100 100	2, 2, 24, 41	0
1	B	334/483 (69%)	-1.74	0 100 100	2, 2, 23, 43	0
1	C	332/483 (68%)	-1.72	0 100 100	2, 4, 26, 53	0
1	D	337/483 (69%)	-1.72	0 100 100	2, 2, 37, 69	0
1	E	334/483 (69%)	-1.73	0 100 100	2, 2, 31, 52	0
1	F	334/483 (69%)	-1.70	0 100 100	2, 11, 32, 53	0
All	All	2134/2898 (73%)	-1.72	0 100 100	2, 3, 29, 69	0

There are no RSRZ outliers to report.

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.