



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 7, 2018 – 12:23 PM EST

PDB ID : 5MVI
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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20030736
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20030736

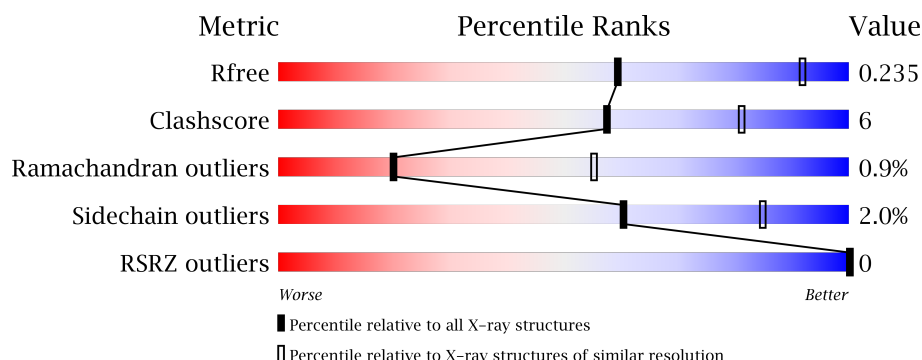
1 Overall quality at a glance

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}	100719	1348 (3.10-3.02)
Clashscore	112137	1462 (3.10-3.02)
Ramachandran outliers	110173	1410 (3.10-3.02)
Sidechain outliers	110143	1410 (3.10-3.02)
RSRZ outliers	101464	1355 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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 63%, a yellow segment representing 5%, and a grey segment representing 31%. The percentages are labeled below the corresponding segments.

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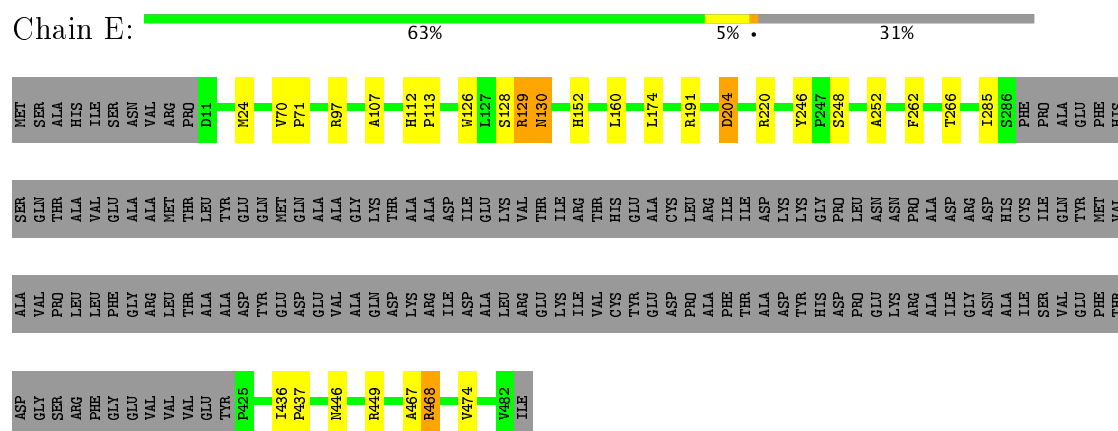
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Chain	Residue	Modelled	Actual	Comment	Reference
F	418	GLY	ASP	conflict	UNP A0A0L5RC52

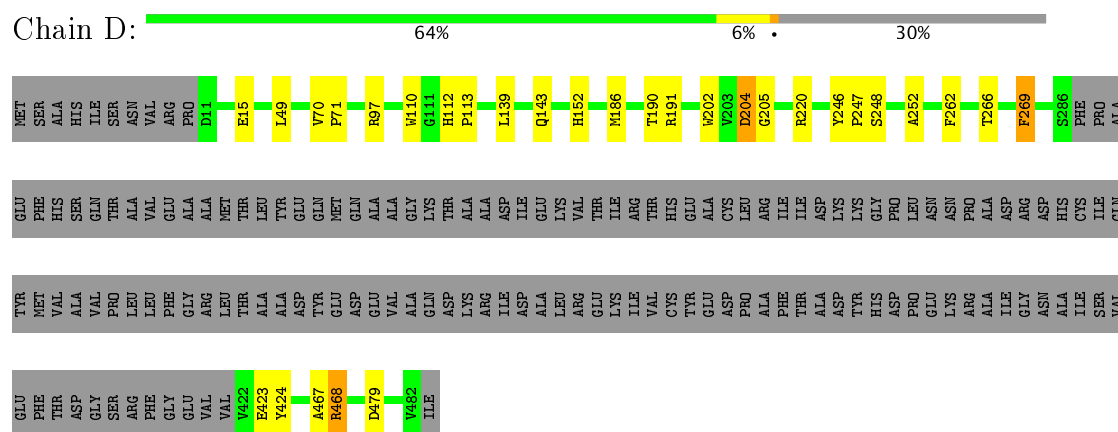
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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.

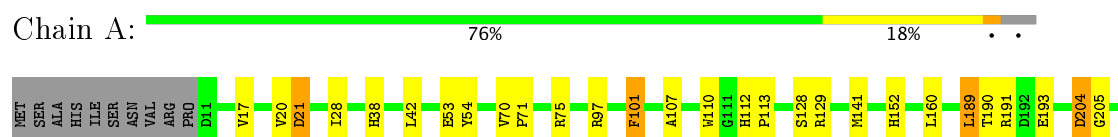
- Molecule 1: 2-methylcitrate dehydratase

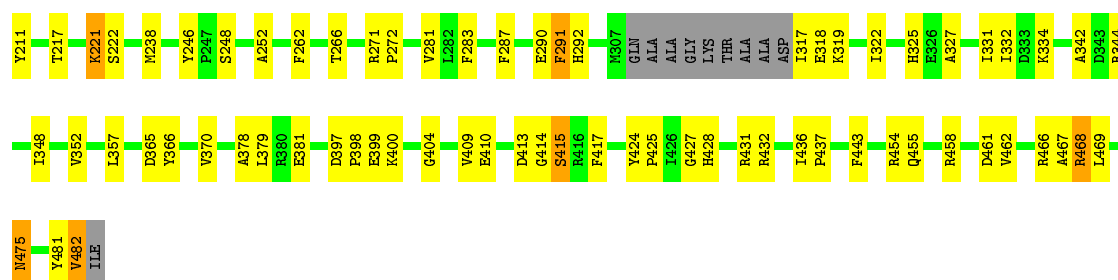


- Molecule 1: 2-methylcitrate dehydratase



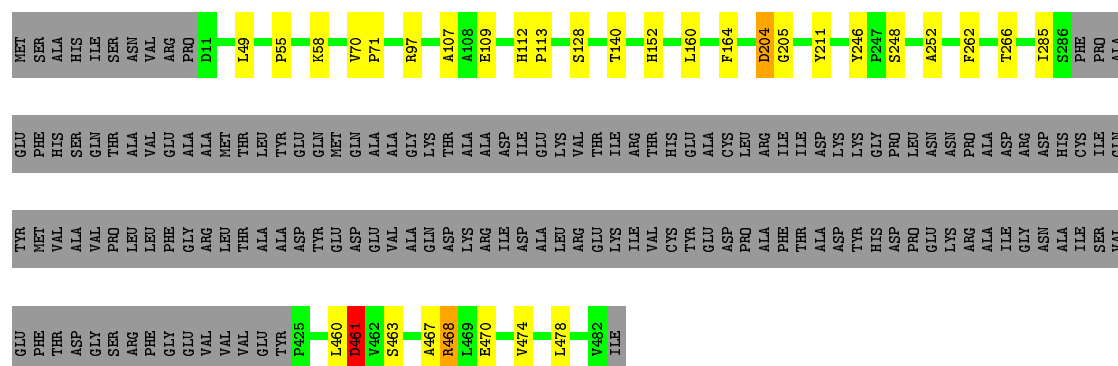
- Molecule 1: 2-methylcitrate dehydratase





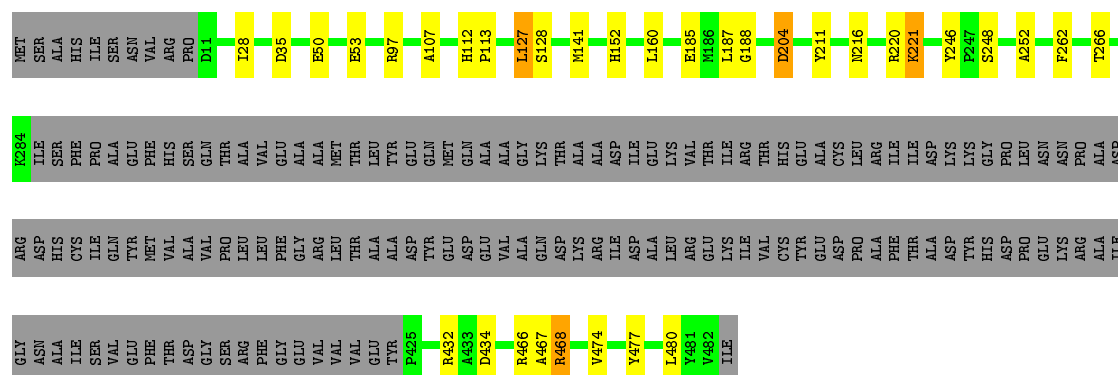
- Molecule 1: 2-methylcitrate dehydratase

Chain B: 63% 6% 31%



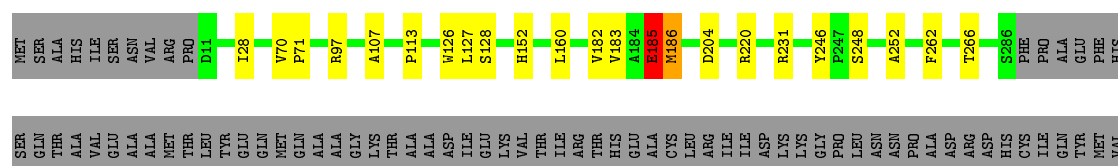
- Molecule 1: 2-methylcitrate dehydratase

Chain C: 62% 6% 31%



- Molecule 1: 2-methylcitrate dehydratase

Chain F: 63% 5% 31%



[illegible]

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 (5.17%)	DCC
Wilson B-factor (Å ²)	25.5	Xtriage
Anisotropy	0.373	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 13.3	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	0.80	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	0.79	1/3352 (0.0%)	0.76	1/4559 (0.0%)
1	B	0.73	0/2406	0.71	2/3272 (0.1%)
1	C	0.66	0/2424	0.71	3/3293 (0.1%)
1	D	0.68	1/2454 (0.0%)	0.72	3/3333 (0.1%)
1	E	0.67	0/2385	0.71	2/3239 (0.1%)
1	F	0.59	1/2385 (0.0%)	0.66	2/3238 (0.1%)
All	All	0.70	3/15406 (0.0%)	0.72	13/20934 (0.1%)

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 (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	481	TYR	CB-CG	11.14	1.68	1.51
1	D	202	TRP	CB-CG	-5.81	1.39	1.50
1	F	126	TRP	CB-CG	5.12	1.59	1.50

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	28	ILE	CG1-CB-CG2	-8.72	92.22	111.40
1	D	49	LEU	CA-CB-CG	6.50	130.26	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	478	LEU	CA-CB-CG	6.50	130.26	115.30
1	C	480	LEU	CA-CB-CG	6.38	129.98	115.30
1	C	127	LEU	CA-CB-CG	6.07	129.26	115.30
1	F	127	LEU	CA-CB-CG	-6.04	101.42	115.30
1	F	231	ARG	NE-CZ-NH1	6.02	123.31	120.30
1	A	189	LEU	CB-CG-CD2	-5.67	101.36	111.00
1	D	186	MET	CA-CB-CG	5.55	122.73	113.30
1	E	129	ARG	N-CA-C	5.25	125.16	111.00
1	E	174	LEU	CA-CB-CG	-5.22	103.29	115.30
1	D	479	ASP	CB-CG-OD2	5.20	122.98	118.30
1	B	49	LEU	CA-CB-CG	-5.07	103.65	115.30

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

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	69	0
1	B	2359	0	2114	24	0
1	C	2375	0	2152	24	0
1	D	2406	0	2172	17	0
1	E	2341	0	2098	20	0
1	F	2342	0	2113	18	0
All	All	15106	0	13539	160	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 (160) 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:460:LEU:O	1:B:461:ASP:O	1.81	0.98
1:E:24:MET:CE	1:E:191:ARG:HA	2.02	0.89
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:E:24:MET:HE3	1:E:191:ARG:HA	1.63	0.78
1:A:291:PHE:O	1:A:291:PHE:HD1	1.66	0.78
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: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:C:141:MET:HE2	1:C:474:VAL:HA	1.81	0.63
1:D:269:PHE:CD2	1:D:269:PHE:N	2.65	0.62
1:C:97:ARG:HG3	1:C:97:ARG:O	2.00	0.62
1:B:204:ASP:OD1	1:B:205:GLY:N	2.29	0.62
1:E:97:ARG:HG3	1:E:97:ARG:O	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:HG3	1:D:97:ARG:O	2.01	0.61
1:A:325:HIS:HD2	1:A:404:GLY:O	1.85	0.60
1:B:97:ARG:HG3	1:B:97:ARG:O	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:365:ASP:HA	1:A:370:VAL:HG21	1.85	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:344:ARG:NH1	1:A:366:TYR:O	2.38	0.57
1:F:467:ALA:O	1:F:468:ARG:CB	2.53	0.57
1:C:467:ALA:O	1:C:468:ARG:CB	2.53	0.56
1:D:467:ALA:O	1:D:468:ARG:CB	2.52	0.56
1:A:20:VAL:O	1:A:21:ASP:CG	2.44	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:C:141:MET:CE	1:C:477:TYR:HB3	2.35	0.55
1:E:24:MET:HE1	1:E:191:ARG:HA	1.83	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:C:141:MET:HE1	1:C:474:VAL:HB	1.88	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:B:460:LEU:C	1:B:461:ASP:O	2.44	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:221:LYS:HD2	1:A:221:LYS:C	2.30	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:C:141:MET:CE	1:C:474:VAL:HA	2.40	0.51
1:A:409:VAL:HB	1:A:417:PHE:CG	2.46	0.51
1:E:113:PRO:HB2	1:E:152:HIS:CE1	2.45	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:357:LEU:HG	1:A:379:LEU:HD23	1.95	0.49
1:D:220:ARG:HB2	1:C:246:TYR:OH	2.12	0.49
1:D:246:TYR:CE1	1:C:211:TYR:HB3	2.48	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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:113:PRO:HB2	1:F:152:HIS:CG	2.48	0.48
1:B:113:PRO:HB2	1:B:152:HIS:CG	2.49	0.47
1:C:141:MET:HE3	1:C:477:TYR:HB3	1.96	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:B:248:SER:O	1:B:252:ALA:HB2	2.14	0.47
1:A:397:ASP:HA	1:A:398:PRO:HD3	1.69	0.47
1:A:399:GLU:CG	1:A:399:GLU:O	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:B:55:PRO:HA	1:B:58:LYS:HE2	1.98	0.45
1:D:139:LEU:CD2	1:D:143:GLN:OE1	2.64	0.45
1:C:127:LEU:HD21	1:C:187:LEU:HD23	1.97	0.45
1:C:141:MET:HE2	1:C:477:TYR:HB3	1.98	0.45
1:B:70:VAL:HA	1:B:71:PRO:HD2	1.82	0.45
1:A:141:MET:HB2	1:A:469:LEU:O	2.16	0.45
1:A:128:SER:HB3	1:A:475:ASN:OD1	2.17	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:248:SER:O	1:E:252:ALA:HB2	2.17	0.45
1:A:455:GLN:HE22	1:A:482:VAL:C	2.20	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:A:221:LYS:C	1:A:221:LYS:CD	2.85	0.44
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:A:413:ASP:C	1:A:415:SER:H	2.21	0.44
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:D:70:VAL:HA	1:D:71:PRO:HD2	1.78	0.43
1:A:113:PRO:HB2	1:A:152:HIS:CE1	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:70:VAL:HA	1:E:71:PRO:HD2	1.86	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:C:221:LYS:C	1:C:221:LYS:HD2	2.39	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
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:A:211:TYR:CB	1:B:246:TYR:CE1	3.02	0.42
1:E:436:ILE:N	1:E:437:PRO:CD	2.82	0.42
1:A:53:GLU:HB2	1:A:54:TYR:CD1	2.54	0.42
1:E:128:SER:OG	1:E:474:VAL:HG22	2.20	0.42
1:B:461:ASP:O	1:B:463:SER:N	2.53	0.42
1:D:15:GLU:CG	1:D:269:PHE:HZ	2.33	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:A:28:ILE:HG12	1:A:28:ILE:H	1.70	0.41
1:B:107:ALA:HB1	1:B:160:LEU:HA	2.03	0.41
1:D:247:PRO:HD2	1:C:216:ASN:OD1	2.20	0.41
1:A:70:VAL:HA	1:A:71:PRO:HD2	1.83	0.41
1:B:58:LYS:HB2	1:B:58:LYS:HE2	1.90	0.41
1:A:292:HIS:CD2	1:A:327:ALA:HB3	2.54	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:O	1:A:454:ARG:HD2	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:HG12	1:F:28:ILE:H	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%)	17	50
1	B	330/483 (68%)	319 (97%)	8 (2%)	3 (1%)	20	55
1	C	328/483 (68%)	319 (97%)	7 (2%)	2 (1%)	28	64
1	D	333/483 (69%)	323 (97%)	7 (2%)	3 (1%)	20	55
1	E	330/483 (68%)	317 (96%)	10 (3%)	3 (1%)	20	55
1	F	330/483 (68%)	320 (97%)	6 (2%)	4 (1%)	15	47
All	All	2110/2898 (73%)	2029 (96%)	61 (3%)	20 (1%)	20	55

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

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Mol	Chain	Res	Type
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%)	43	76
1	B	196/395 (50%)	194 (99%)	2 (1%)	80	92
1	C	207/395 (52%)	201 (97%)	6 (3%)	48	78
1	D	207/395 (52%)	203 (98%)	4 (2%)	62	85
1	E	194/395 (49%)	191 (98%)	3 (2%)	70	89
1	F	196/395 (50%)	194 (99%)	2 (1%)	80	92
All	All	1270/2370 (54%)	1244 (98%)	26 (2%)	60	85

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	204	ASP
1	E	262	PHE
1	E	266	THR
1	D	110	TRP
1	D	262	PHE
1	D	266	THR
1	D	269	PHE
1	A	101	PHE
1	A	110	TRP
1	A	221	LYS
1	A	238	MET
1	A	262	PHE
1	A	266	THR
1	A	291	PHE
1	A	475	ASN

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Mol	Chain	Res	Type
1	A	482	VAL
1	B	262	PHE
1	B	266	THR
1	C	204	ASP
1	C	221	LYS
1	C	262	PHE
1	C	266	THR
1	C	432	ARG
1	C	434	ASP
1	F	262	PHE
1	F	266	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	280	ASN
1	A	325	HIS
1	A	455	GLN
1	C	280	ASN
1	F	280	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates 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%)	-0.79	0 100 100	2, 2, 24, 41	0
1	B	334/483 (69%)	-0.78	0 100 100	2, 2, 23, 43	0
1	C	332/483 (68%)	-0.71	0 100 100	2, 4, 26, 53	0
1	D	337/483 (69%)	-0.72	0 100 100	2, 2, 37, 69	0
1	E	334/483 (69%)	-0.73	0 100 100	2, 2, 31, 52	0
1	F	334/483 (69%)	-0.67	0 100 100	2, 11, 32, 53	0
All	All	2134/2898 (73%)	-0.74	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 carbohydrates 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.