



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 3, 2017 – 08:09 PM EDT

PDB ID : 5VBF
Title : Crystal structure of succinate semialdehyde dehydrogenase from *Burkholderia vietnamiensis*
Authors : Seattle Structural Genomics Center for Infectious Disease (SSGCID)
Deposited on : 2017-03-29
Resolution : 2.35 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
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20029077
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-20029077

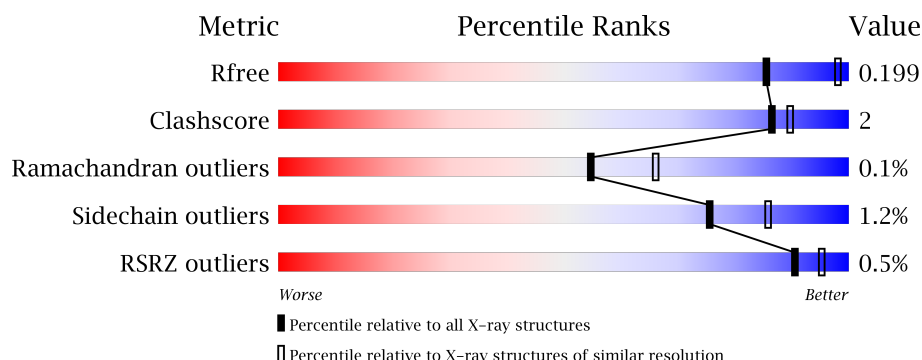
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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.



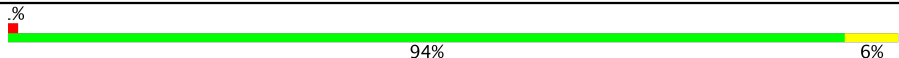
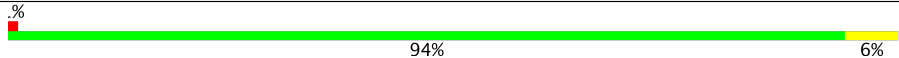
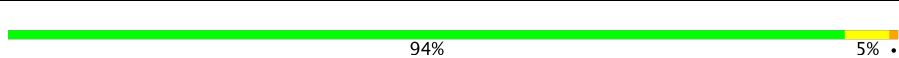
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	100719	1522 (2.38-2.34)
Clashscore	112137	1626 (2.38-2.34)
Ramachandran outliers	110173	1605 (2.38-2.34)
Sidechain outliers	110143	1606 (2.38-2.34)
RSRZ outliers	101464	1528 (2.38-2.34)

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	478	<div> <div>92%</div> <div>8%</div> </div>
1	B	478	<div> <div>94%</div> <div>6%</div> </div>
1	C	478	<div> <div>92%</div> <div>6%</div> </div>
1	D	478	<div> <div>94%</div> <div>5%</div> </div>
1	E	478	<div> <div>93%</div> <div>6%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	478	
1	G	478	
1	H	478	

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	MES	A	601	-	-	-	X
2	MES	B	601	-	-	-	X
2	MES	E	601	-	-	-	X
2	MES	F	601	-	-	-	X
2	MES	G	601	-	-	-	X
2	MES	H	601	-	-	-	X
4	FMT	B	603	-	-	-	X
4	FMT	E	603	-	-	-	X
5	UNX	A	604	-	-	-	X
5	UNX	B	604	-	-	-	X
5	UNX	C	603	-	-	-	X
5	UNX	D	603	-	-	-	X
5	UNX	E	604	-	-	-	X
5	UNX	F	603	-	-	-	X
5	UNX	G	603	-	-	-	X
5	UNX	H	604	-	-	-	X

2 Entry composition

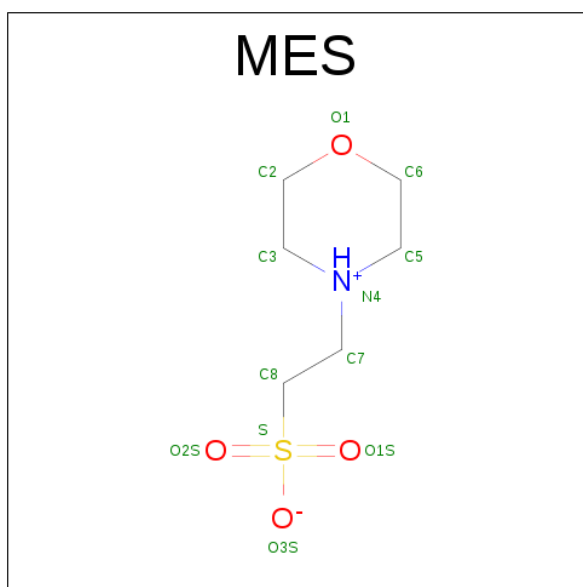
There are 6 unique types of molecules in this entry. The entry contains 31891 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 NAD-dependent succinate-semialdehyde dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	478	Total	C	N	O	S	0	7	0
			3619	2288	655	655	21			
1	B	478	Total	C	N	O	S	0	8	0
			3643	2301	665	656	21			
1	C	476	Total	C	N	O	S	0	2	0
			3584	2266	650	648	20			
1	D	477	Total	C	N	O	S	0	5	0
			3591	2273	644	653	21			
1	E	476	Total	C	N	O	S	0	5	0
			3584	2268	641	654	21			
1	F	476	Total	C	N	O	S	0	6	0
			3596	2277	646	652	21			
1	G	477	Total	C	N	O	S	0	3	0
			3582	2265	643	653	21			
1	H	477	Total	C	N	O	S	0	5	0
			3605	2279	651	654	21			

- Molecule 2 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: C₆H₁₃NO₄S).



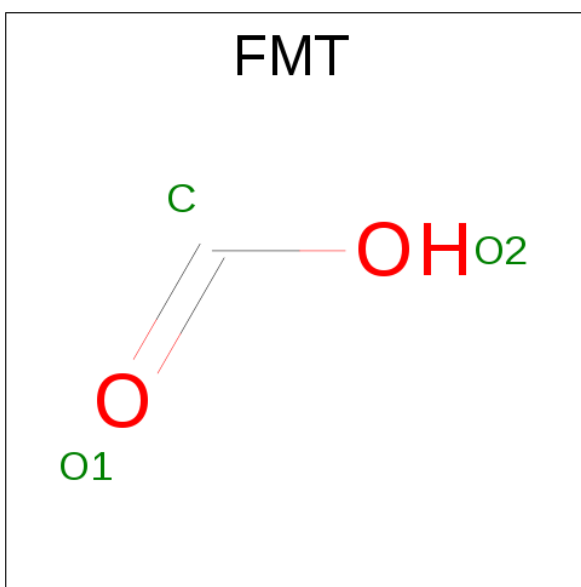
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
2	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
2	C	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
2	D	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
2	E	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
2	F	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
2	G	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
2	H	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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		
3	E	1	Total	C	O	0	0
			6	3	3		
3	F	1	Total	C	O	0	0
			6	3	3		
3	G	1	Total	C	O	0	0
			6	3	3		
3	H	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is FORMIC ACID (three-letter code: FMT) (formula: CH₂O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			3	1	2		
4	B	1	Total	C	O	0	0
			3	1	2		
4	E	1	Total	C	O	0	0
			3	1	2		
4	H	1	Total	C	O	0	0
			3	1	2		

- Molecule 5 is UNKNOWN ATOM OR ION (three-letter code: UNX) (formula: X).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	G	1	Total	X	0	0
			1	1		
5	D	1	Total	X	0	0
			1	1		
5	E	1	Total	X	0	0
			1	1		
5	H	1	Total	X	0	0
			1	1		
5	B	1	Total	X	0	0
			1	1		
5	C	1	Total	X	0	0
			1	1		
5	A	1	Total	X	0	0
			1	1		
5	F	1	Total	X	0	0
			1	1		

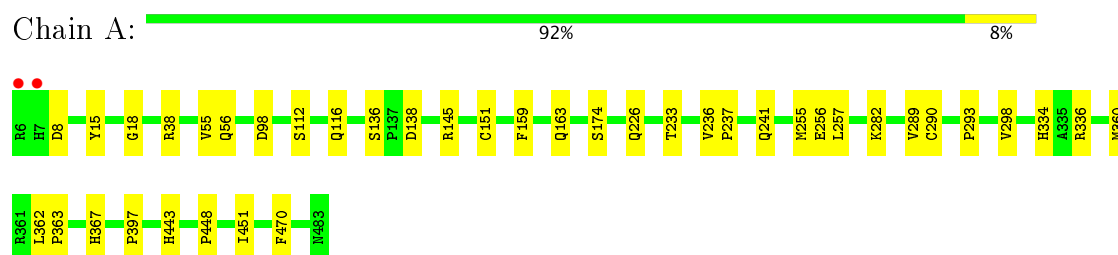
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	378	Total 378	O 378	0	0
6	B	441	Total 441	O 441	0	1
6	C	391	Total 391	O 391	0	0
6	D	371	Total 371	O 371	0	0
6	E	377	Total 378	O 378	0	1
6	F	297	Total 298	O 298	0	2
6	G	320	Total 320	O 320	0	0
6	H	345	Total 346	O 346	0	1

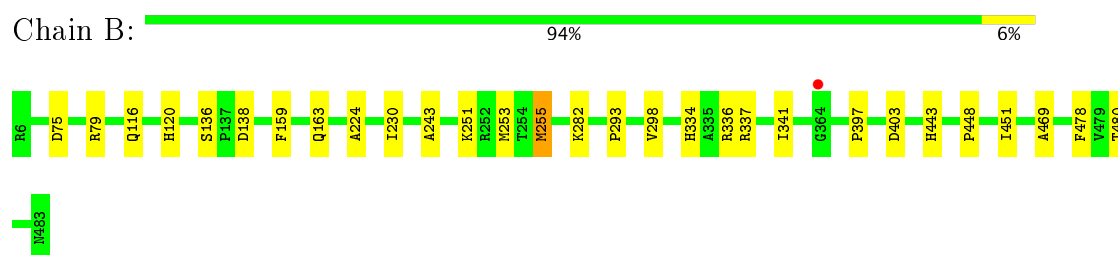
3 Residue-property plots [i](#)

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.

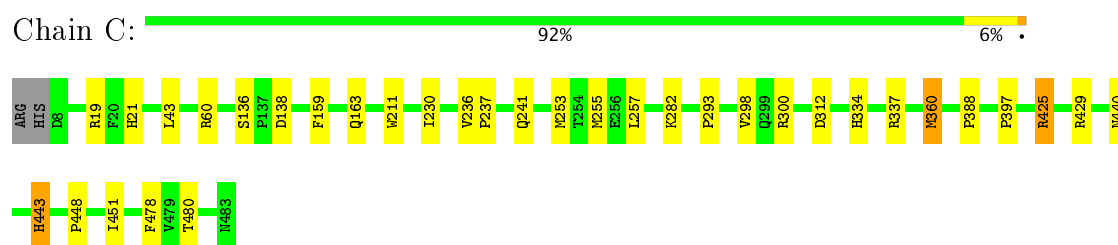
- Molecule 1: NAD-dependent succinate-semialdehyde dehydrogenase



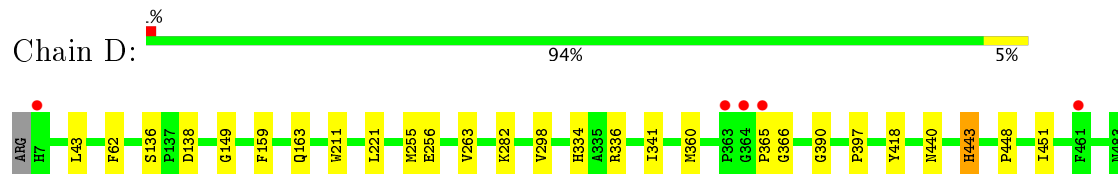
- Molecule 1: NAD-dependent succinate-semialdehyde dehydrogenase



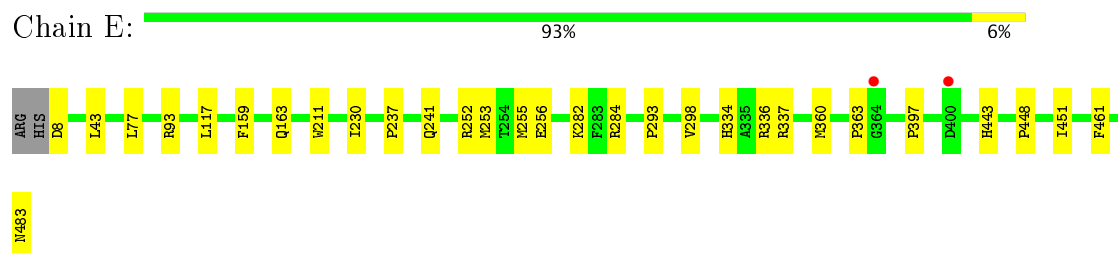
- Molecule 1: NAD-dependent succinate-semialdehyde dehydrogenase



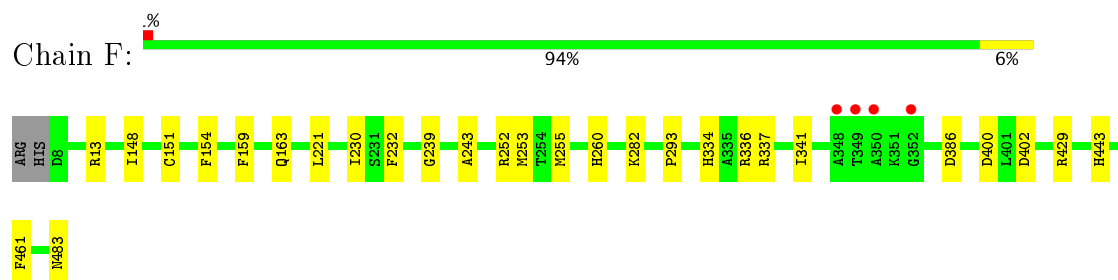
- Molecule 1: NAD-dependent succinate-semialdehyde dehydrogenase



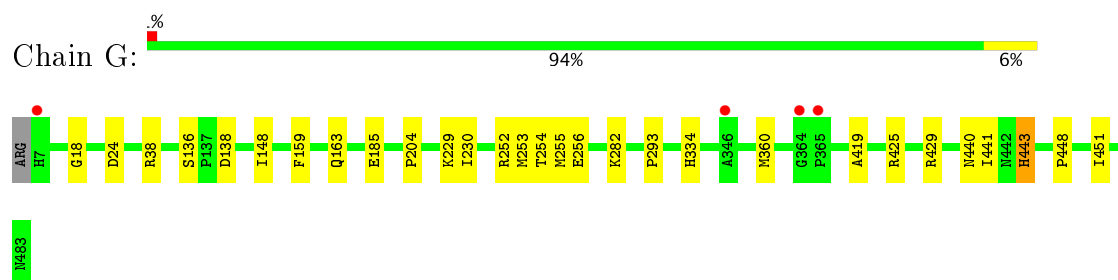
- Molecule 1: NAD-dependent succinate-semialdehyde dehydrogenase



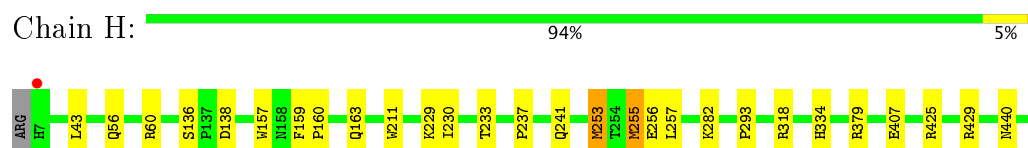
- Molecule 1: NAD-dependent succinate-semialdehyde dehydrogenase



- Molecule 1: NAD-dependent succinate-semialdehyde dehydrogenase



- Molecule 1: NAD-dependent succinate-semialdehyde dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	106.62Å 191.76Å 219.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.17 – 2.35 48.84 – 2.35	Depositor EDS
% Data completeness (in resolution range)	98.7 (48.17-2.35) 98.7 (48.84-2.35)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.84 (at 2.34Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.147 , 0.196 0.153 , 0.199	Depositor DCC
R_{free} test set	1984 reflections (1.08%)	DCC
Wilson B-factor (Å ²)	16.7	Xtriage
Anisotropy	0.289	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 49.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	31891	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 6.21% 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: UNX, GOL, FMT, MES

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	0.40	0/3725	0.55	1/5064 (0.0%)
1	B	0.40	0/3747	0.56	0/5090
1	C	0.41	0/3674	0.56	0/4994
1	D	0.39	0/3689	0.56	0/5018
1	E	0.39	0/3683	0.55	0/5009
1	F	0.38	0/3697	0.55	0/5027
1	G	0.38	0/3675	0.55	0/4998
1	H	0.41	0/3705	0.55	0/5037
All	All	0.39	0/29595	0.55	1/40237 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	98	ASP	CB-CG-OD1	5.16	122.94	118.30

There are no chirality outliers.

There are no planarity outliers.

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	3619	0	3581	24	0
1	B	3643	0	3617	19	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	3584	0	3551	18	0
1	D	3591	0	3538	15	0
1	E	3584	0	3537	17	0
1	F	3596	0	3553	21	0
1	G	3582	0	3532	14	0
1	H	3605	0	3566	16	0
2	A	12	0	12	2	0
2	B	12	0	12	0	0
2	C	12	0	12	0	0
2	D	12	0	12	0	0
2	E	12	0	12	0	0
2	F	12	0	12	1	0
2	G	12	0	12	1	0
2	H	12	0	12	0	0
3	A	6	0	8	0	0
3	B	6	0	8	0	0
3	C	6	0	8	0	0
3	D	6	0	8	0	0
3	E	6	0	8	0	0
3	F	6	0	8	0	0
3	G	6	0	8	0	0
3	H	6	0	8	0	0
4	A	3	0	1	0	0
4	B	3	0	1	0	0
4	E	3	0	1	1	0
4	H	3	0	1	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
5	C	1	0	0	0	0
5	D	1	0	0	0	0
5	E	1	0	0	0	0
5	F	1	0	0	0	0
5	G	1	0	0	0	0
5	H	1	0	0	0	0
6	A	378	0	0	4	0
6	B	441	0	0	3	0
6	C	391	0	0	1	0
6	D	371	0	0	1	0
6	E	378	0	0	3	0
6	F	298	0	0	1	0
6	G	320	0	0	1	0
6	H	346	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	31891	0	28639	143	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 (143) 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:F:230:ILE:HB	1:F:253[B]:MET:HG3	1.49	0.93
1:B:334[B]:HIS:HD2	1:B:336:ARG:HG2	1.41	0.86
1:B:334[B]:HIS:CD2	1:B:336:ARG:HG2	2.12	0.84
1:G:136:SER:OG	1:G:138:ASP:OD1	1.99	0.80
1:D:334[B]:HIS:HD2	1:D:336:ARG:HG2	1.47	0.78
1:D:136:SER:OG	1:D:138:ASP:OD1	2.03	0.76
1:C:312:ASP:HB3	1:E:363:PRO:HB3	1.68	0.74
1:H:379:ARG:NH1	1:H:407:GLU:OE1	2.22	0.72
1:C:136:SER:OG	1:C:138:ASP:OD1	2.07	0.71
1:H:255:MET:HG2	1:H:257:LEU:HD21	1.74	0.67
1:B:334[B]:HIS:CD2	1:B:337:ARG:HG2	2.30	0.65
1:A:334:HIS:ND1	1:A:336:ARG:HG2	2.12	0.64
1:F:239:GLY:O	1:F:255:MET:HE1	1.98	0.64
1:E:334:HIS:ND1	1:E:336:ARG:HG2	2.13	0.63
1:C:19:ARG:HD2	1:C:21:HIS:CE1	2.34	0.63
1:E:8:ASP:O	1:E:93[A]:ARG:NH2	2.25	0.63
1:D:334[B]:HIS:CD2	1:D:336:ARG:HG2	2.31	0.63
1:B:403:ASP:HB3	6:B:807:HOH:O	1.99	0.63
1:A:136:SER:OG	1:A:138:ASP:OD1	2.16	0.62
1:B:136:SER:OG	1:B:138:ASP:OD1	2.14	0.62
1:H:425:ARG:HH11	1:H:429[A]:ARG:CZ	2.13	0.61
1:H:136:SER:OG	1:H:138:ASP:OD1	2.18	0.60
1:F:402:ASP:OD1	1:F:429[B]:ARG:NH2	2.35	0.60
1:B:334[B]:HIS:ND1	6:B:702:HOH:O	2.25	0.60
1:A:145:ARG:NH2	6:A:703:HOH:O	2.36	0.59
1:A:289:VAL:HG13	2:A:601:MES:H81	1.84	0.59
1:A:112:SER:O	1:A:116[B]:GLN:HG2	2.03	0.58
1:B:116:GLN:O	1:B:120:HIS:HD2	1.86	0.57
1:F:243:ALA:HB2	1:F:255:MET:HE1	1.86	0.57
1:B:243:ALA:HB2	1:B:255:MET:HE2	1.87	0.56
1:G:18:GLY:HA3	1:G:204:PRO:HB3	1.87	0.56
1:G:159:PHE:CE1	2:G:601:MES:H81	2.42	0.55
1:F:243:ALA:HB2	1:F:255:MET:CE	2.37	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:425:ARG:HE	1:C:429:ARG:NH2	2.05	0.54
1:C:60[B]:ARG:NH1	6:C:704:HOH:O	2.41	0.54
1:G:256:GLU:HB3	6:G:950:HOH:O	2.08	0.54
1:A:55:VAL:HG12	1:A:226[B]:GLN:HE21	1.73	0.53
1:E:159:PHE:HB2	1:E:163:GLN:HG2	1.92	0.52
1:D:159:PHE:HB2	1:D:163:GLN:HG2	1.92	0.52
1:E:334:HIS:HB2	6:E:978:HOH:O	2.11	0.51
1:E:256:GLU:HB3	6:E:955:HOH:O	2.11	0.51
1:C:448:PRO:O	1:C:451:ILE:HG13	2.11	0.51
1:H:159:PHE:HB2	1:H:163:GLN:HG2	1.93	0.50
1:E:334:HIS:CE1	1:E:336:ARG:HG2	2.47	0.50
1:A:56:GLN:HA	1:A:226[B]:GLN:HE22	1.76	0.49
1:F:334[A]:HIS:ND1	1:F:336:ARG:HB3	2.27	0.49
1:F:148:ILE:HD12	1:F:252:ARG:HD3	1.93	0.49
1:A:256:GLU:HB3	6:A:973:HOH:O	2.12	0.49
1:A:8:ASP:OD1	1:A:8:ASP:N	2.42	0.48
1:C:237:PRO:O	1:C:241:GLN:HG3	2.13	0.48
1:E:237:PRO:O	1:E:241:GLN:HG3	2.13	0.48
1:C:236:VAL:HA	1:C:257:LEU:HD13	1.96	0.47
2:F:601:MES:H51	2:F:601:MES:H82	1.49	0.47
1:F:253[B]:MET:SD	1:F:255:MET:CE	3.03	0.47
1:B:337:ARG:O	1:B:341:ILE:HG12	2.14	0.47
1:G:159:PHE:HB2	1:G:163:GLN:HG2	1.95	0.47
1:G:282:LYS:HD3	1:G:293:PRO:O	2.15	0.47
1:G:230:ILE:O	1:G:253:MET:HA	2.15	0.47
1:H:440:ASN:HB3	1:H:443:HIS:O	2.15	0.47
1:H:157:TRP:O	1:H:160:PRO:HD3	2.15	0.47
1:G:229:LYS:NZ	1:G:254:THR:OG1	2.48	0.47
1:D:43:LEU:HD11	1:D:211:TRP:CE2	2.51	0.46
1:F:334[A]:HIS:CE1	1:F:336:ARG:HE	2.33	0.46
1:B:159:PHE:HB2	1:B:163:GLN:HG2	1.97	0.46
1:G:425:ARG:NH2	1:G:429:ARG:HD2	2.31	0.46
1:B:478:PHE:CZ	1:B:480:THR:HB	2.51	0.46
1:A:151:CYS:HB3	6:A:872:HOH:O	2.16	0.45
1:C:337:ARG:NH1	1:C:388:PRO:O	2.49	0.45
1:D:282:LYS:HD3	1:D:282:LYS:O	2.17	0.45
1:E:448:PRO:O	1:E:451:ILE:HG13	2.17	0.45
1:B:224:ALA:O	1:B:251:LYS:NZ	2.49	0.45
1:F:159:PHE:HB2	1:F:163:GLN:HG2	1.99	0.45
1:C:478:PHE:CZ	1:C:480:THR:HB	2.52	0.44
1:D:62:PHE:CZ	1:D:149:GLY:HA2	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:282:LYS:HD3	1:H:293:PRO:O	2.18	0.44
1:A:15:TYR:OH	1:A:18:GLY:HA2	2.17	0.44
1:B:282:LYS:HD3	1:B:293:PRO:O	2.17	0.44
1:A:282:LYS:HD3	1:A:293:PRO:O	2.17	0.44
1:D:440:ASN:HB3	1:D:443:HIS:O	2.17	0.44
1:E:43:LEU:HD11	1:E:211:TRP:CE2	2.52	0.44
1:G:148:ILE:CD1	1:G:252:ARG:HG3	2.48	0.44
1:A:362:LEU:HB3	1:A:363:PRO:HD2	1.98	0.44
1:E:334:HIS:CE1	1:E:337:ARG:HG2	2.53	0.43
1:F:337:ARG:O	1:F:341:ILE:HG12	2.18	0.43
1:D:256:GLU:HB3	6:D:932:HOH:O	2.19	0.43
1:B:334[B]:HIS:CE1	6:B:702:HOH:O	2.70	0.43
1:E:282:LYS:HD3	1:E:293:PRO:O	2.18	0.43
1:A:290:CYS:HB2	2:A:601:MES:O2S	2.18	0.43
1:D:341:ILE:HD11	1:D:390:GLY:HA3	2.01	0.43
1:H:233:THR:HA	1:H:256:GLU:O	2.19	0.43
1:A:233:THR:HA	1:A:256:GLU:O	2.19	0.43
1:A:174:SER:HB3	1:A:470:PHE:CD1	2.54	0.43
1:C:230:ILE:O	1:C:253:MET:HA	2.20	0.42
1:A:334:HIS:CE1	1:A:336:ARG:HH11	2.37	0.42
1:E:230:ILE:O	1:E:253:MET:HA	2.19	0.42
1:E:483:ASN:HA	4:E:603:FMT:C	2.49	0.42
1:H:230:ILE:O	1:H:253:MET:HA	2.20	0.42
1:E:77:LEU:HG	1:E:117:LEU:HD11	2.01	0.42
1:H:229:LYS:HG2	1:H:230:ILE:N	2.33	0.42
1:B:75:ASP:O	1:B:79[A]:ARG:HG3	2.20	0.42
1:F:400:ASP:OD1	1:F:402:ASP:HB2	2.19	0.42
1:B:230:ILE:O	1:B:253:MET:HA	2.19	0.42
1:C:298:VAL:O	1:C:397:PRO:HA	2.20	0.42
1:F:253[B]:MET:SD	1:F:255:MET:HE2	2.59	0.42
1:F:253[B]:MET:SD	1:F:255:MET:HE3	2.60	0.42
1:D:298:VAL:O	1:D:397:PRO:HA	2.20	0.42
1:E:298:VAL:O	1:E:397:PRO:HA	2.19	0.42
1:H:425:ARG:NH1	1:H:429[A]:ARG:CZ	2.82	0.42
1:B:448:PRO:O	1:B:451:ILE:HG13	2.20	0.41
1:C:425:ARG:HE	1:C:429:ARG:CZ	2.32	0.41
1:D:282:LYS:C	1:D:282:LYS:HD3	2.41	0.41
1:F:154:PHE:HB2	1:F:232:PHE:HD1	1.85	0.41
1:H:318:ARG:HD2	6:H:993:HOH:O	2.20	0.41
1:E:284:ARG:NE	6:E:706:HOH:O	2.45	0.41
1:F:151:CYS:HB3	6:F:811:HOH:O	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:43:LEU:HD11	1:H:211:TRP:CE2	2.55	0.41
1:A:334:HIS:HB2	6:A:992:HOH:O	2.19	0.41
1:F:483:ASN:HB3	1:G:425:ARG:HA	2.01	0.41
1:C:360:MET:SD	1:C:360:MET:N	2.93	0.41
1:A:448:PRO:O	1:A:451:ILE:HG13	2.20	0.41
1:F:148:ILE:CD1	1:F:252:ARG:HD3	2.50	0.41
1:C:43:LEU:HD11	1:C:211:TRP:CE2	2.55	0.41
1:B:120:HIS:CE1	1:B:469:ALA:HB1	2.56	0.41
1:B:298:VAL:O	1:B:397:PRO:HA	2.20	0.41
1:C:159:PHE:HB2	1:C:163:GLN:HG2	2.03	0.41
1:F:221:LEU:HD23	1:F:221:LEU:HA	1.92	0.41
1:A:236:VAL:HA	1:A:257:LEU:HD13	2.02	0.41
1:D:263:VAL:HG22	1:D:418:TYR:HB2	2.03	0.41
1:A:237:PRO:O	1:A:241:GLN:HG3	2.21	0.41
1:D:221:LEU:HA	1:D:221:LEU:HD23	1.89	0.41
1:D:448:PRO:O	1:D:451:ILE:HG13	2.20	0.41
1:A:334:HIS:HA	1:A:367:HIS:CD2	2.56	0.41
1:C:440:ASN:HB3	1:C:443:HIS:O	2.21	0.41
1:F:260:HIS:NE2	1:F:386:ASP:O	2.54	0.40
1:A:159:PHE:HB2	1:A:163:GLN:HG2	2.03	0.40
1:A:298:VAL:O	1:A:397:PRO:HA	2.22	0.40
1:H:237:PRO:O	1:H:241:GLN:HG3	2.20	0.40
1:F:282:LYS:HD3	1:F:293:PRO:O	2.22	0.40
1:G:440:ASN:HB3	1:G:443:HIS:O	2.21	0.40
1:G:448:PRO:O	1:G:451:ILE:HG13	2.21	0.40
1:C:282:LYS:HD3	1:C:293:PRO:O	2.22	0.40
1:G:419:ALA:O	1:G:441:ILE:HA	2.21	0.40
1:H:56[A]:GLN:HG2	1:H:60[A]:ARG:HH12	1.86	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	483/478 (101%)	473 (98%)	10 (2%)	0	100	100
1	B	484/478 (101%)	476 (98%)	8 (2%)	0	100	100
1	C	476/478 (100%)	465 (98%)	11 (2%)	0	100	100
1	D	480/478 (100%)	469 (98%)	8 (2%)	3 (1%)	28	32
1	E	479/478 (100%)	470 (98%)	9 (2%)	0	100	100
1	F	480/478 (100%)	469 (98%)	11 (2%)	0	100	100
1	G	478/478 (100%)	468 (98%)	9 (2%)	1 (0%)	51	61
1	H	480/478 (100%)	471 (98%)	9 (2%)	0	100	100
All	All	3840/3824 (100%)	3761 (98%)	75 (2%)	4 (0%)	55	66

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	366	GLY
1	D	365[A]	PRO
1	D	365[B]	PRO
1	G	24	ASP

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	364/364 (100%)	360 (99%)	4 (1%)	78	87
1	B	367/364 (101%)	365 (100%)	2 (0%)	91	95
1	C	359/364 (99%)	353 (98%)	6 (2%)	66	77
1	D	360/364 (99%)	357 (99%)	3 (1%)	85	91
1	E	360/364 (99%)	355 (99%)	5 (1%)	71	82
1	F	361/364 (99%)	357 (99%)	4 (1%)	78	87
1	G	359/364 (99%)	352 (98%)	7 (2%)	62	75
1	H	363/364 (100%)	358 (99%)	5 (1%)	71	82
All	All	2893/2912 (99%)	2857 (99%)	36 (1%)	75	86

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	ARG
1	A	255	MET
1	A	360	MET
1	A	443	HIS
1	B	255	MET
1	B	443	HIS
1	C	255	MET
1	C	300	ARG
1	C	334	HIS
1	C	360	MET
1	C	425	ARG
1	C	443	HIS
1	D	255	MET
1	D	360	MET
1	D	443	HIS
1	E	252	ARG
1	E	255	MET
1	E	360	MET
1	E	443	HIS
1	E	461	PHE
1	F	13[A]	ARG
1	F	13[B]	ARG
1	F	443	HIS
1	F	461	PHE
1	G	38	ARG
1	G	185[A]	GLU
1	G	185[B]	GLU
1	G	255	MET
1	G	334	HIS
1	G	360	MET
1	G	443	HIS
1	H	253	MET
1	H	255	MET
1	H	334[A]	HIS
1	H	334[B]	HIS
1	H	443	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	120	HIS

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Mol	Chain	Res	Type
1	C	226	GLN
1	C	260	HIS
1	C	367	HIS

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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 28 ligands modelled in this entry, 8 are unknown - leaving 20 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	MES	A	601	-	12,12,12	2.11	1 (8%)	14,16,16	3.02	7 (50%)
3	GOL	A	602	-	5,5,5	0.88	0	5,5,5	0.97	0
4	FMT	A	603	-	0,2,2	0.00	-	0,1,1	0.00	-
2	MES	B	601	-	12,12,12	2.11	1 (8%)	14,16,16	2.65	7 (50%)
3	GOL	B	602	-	5,5,5	0.82	0	5,5,5	1.10	0
4	FMT	B	603	-	0,2,2	0.00	-	0,1,1	0.00	-
2	MES	C	601	-	12,12,12	2.06	1 (8%)	14,16,16	2.59	6 (42%)
3	GOL	C	602	-	5,5,5	0.93	0	5,5,5	1.03	0
2	MES	D	601	-	12,12,12	2.11	1 (8%)	14,16,16	2.63	7 (50%)
3	GOL	D	602	-	5,5,5	1.14	0	5,5,5	0.92	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	MES	E	601	-	12,12,12	2.10	1 (8%)	14,16,16	2.90	7 (50%)
3	GOL	E	602	-	5,5,5	0.79	0	5,5,5	1.04	0
4	FMT	E	603	-	0,2,2	0.00	-	0,1,1	0.00	-
2	MES	F	601	-	12,12,12	2.24	1 (8%)	14,16,16	2.44	5 (35%)
3	GOL	F	602	-	5,5,5	1.00	0	5,5,5	0.78	0
2	MES	G	601	-	12,12,12	2.05	1 (8%)	14,16,16	2.42	6 (42%)
3	GOL	G	602	-	5,5,5	0.75	0	5,5,5	1.14	0
2	MES	H	601	-	12,12,12	2.16	1 (8%)	14,16,16	2.93	7 (50%)
3	GOL	H	602	-	5,5,5	0.96	0	5,5,5	0.97	0
4	FMT	H	603	-	0,2,2	0.00	-	0,1,1	0.00	-

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	MES	A	601	-	-	0/6/14/14	0/1/1/1
3	GOL	A	602	-	-	0/4/4/4	0/0/0/0
4	FMT	A	603	-	-	0/0/0/0	0/0/0/0
2	MES	B	601	-	-	0/6/14/14	0/1/1/1
3	GOL	B	602	-	-	0/4/4/4	0/0/0/0
4	FMT	B	603	-	-	0/0/0/0	0/0/0/0
2	MES	C	601	-	-	0/6/14/14	0/1/1/1
3	GOL	C	602	-	-	0/4/4/4	0/0/0/0
2	MES	D	601	-	-	0/6/14/14	0/1/1/1
3	GOL	D	602	-	-	0/4/4/4	0/0/0/0
2	MES	E	601	-	-	0/6/14/14	0/1/1/1
3	GOL	E	602	-	-	0/4/4/4	0/0/0/0
4	FMT	E	603	-	-	0/0/0/0	0/0/0/0
2	MES	F	601	-	-	0/6/14/14	0/1/1/1
3	GOL	F	602	-	-	0/4/4/4	0/0/0/0
2	MES	G	601	-	-	0/6/14/14	0/1/1/1
3	GOL	G	602	-	-	0/4/4/4	0/0/0/0
2	MES	H	601	-	-	0/6/14/14	0/1/1/1
3	GOL	H	602	-	-	0/4/4/4	0/0/0/0
4	FMT	H	603	-	-	0/0/0/0	0/0/0/0

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	601	MES	C8-S	-7.52	1.66	1.77
2	H	601	MES	C8-S	-7.14	1.66	1.77
2	B	601	MES	C8-S	-7.00	1.67	1.77
2	A	601	MES	C8-S	-7.00	1.67	1.77
2	D	601	MES	C8-S	-6.97	1.67	1.77
2	E	601	MES	C8-S	-6.96	1.67	1.77
2	G	601	MES	C8-S	-6.83	1.67	1.77
2	C	601	MES	C8-S	-6.80	1.67	1.77

All (52) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	MES	C2-C3-N4	-5.06	103.02	110.11
2	H	601	MES	C2-C3-N4	-4.67	103.56	110.11
2	G	601	MES	C2-C3-N4	-4.24	104.17	110.11
2	H	601	MES	C6-C5-N4	-3.89	104.66	110.11
2	F	601	MES	C6-C5-N4	-3.86	104.69	110.11
2	E	601	MES	C2-C3-N4	-3.74	104.86	110.11
2	E	601	MES	C6-C5-N4	-3.71	104.91	110.11
2	D	601	MES	C2-C3-N4	-3.66	104.98	110.11
2	C	601	MES	C2-C3-N4	-3.28	105.52	110.11
2	C	601	MES	C6-C5-N4	-2.94	105.98	110.11
2	D	601	MES	C6-C5-N4	-2.93	106.00	110.11
2	B	601	MES	C2-C3-N4	-2.79	106.19	110.11
2	B	601	MES	C6-C5-N4	-2.02	107.28	110.11
2	A	601	MES	C6-O1-C2	2.13	117.09	109.89
2	D	601	MES	O2S-S-C8	2.36	108.82	106.79
2	C	601	MES	C7-N4-C5	2.58	117.88	111.26
2	E	601	MES	C7-N4-C3	2.69	118.15	111.26
2	B	601	MES	O3S-S-C8	2.84	109.55	106.06
2	B	601	MES	C7-N4-C3	2.84	118.53	111.26
2	H	601	MES	O3S-S-C8	2.87	109.59	106.06
2	D	601	MES	C7-N4-C3	2.89	118.65	111.26
2	G	601	MES	C7-N4-C3	2.91	118.71	111.26
2	G	601	MES	O3S-S-C8	3.01	109.76	106.06
2	F	601	MES	C7-N4-C5	3.06	119.09	111.26
2	G	601	MES	C7-N4-C5	3.11	119.22	111.26
2	B	601	MES	C7-N4-C5	3.20	119.47	111.26
2	C	601	MES	C7-N4-C3	3.25	119.59	111.26
2	D	601	MES	O3S-S-C8	3.26	110.06	106.06
2	F	601	MES	C5-N4-C3	3.27	116.28	108.87
2	A	601	MES	C7-N4-C5	3.29	119.69	111.26
2	H	601	MES	C7-N4-C5	3.30	119.71	111.26

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	601	MES	C7-N4-C5	3.34	119.83	111.26
2	A	601	MES	C7-N4-C3	3.35	119.83	111.26
2	H	601	MES	C7-N4-C3	3.43	120.05	111.26
2	E	601	MES	C7-N4-C5	3.53	120.32	111.26
2	G	601	MES	C5-N4-C3	3.67	117.19	108.87
2	G	601	MES	O2S-S-C8	3.83	110.08	106.79
2	E	601	MES	O3S-S-C8	3.85	110.80	106.06
2	A	601	MES	O3S-S-C8	4.02	111.00	106.06
2	E	601	MES	O2S-S-C8	4.14	110.35	106.79
2	F	601	MES	O1S-S-C8	4.20	110.40	106.79
2	C	601	MES	C5-N4-C3	4.22	118.42	108.87
2	B	601	MES	O2S-S-C8	4.40	110.57	106.79
2	F	601	MES	O2S-S-C8	4.55	110.70	106.79
2	H	601	MES	C5-N4-C3	4.69	119.48	108.87
2	H	601	MES	O1S-S-C8	4.75	110.88	106.79
2	A	601	MES	C5-N4-C3	4.99	120.17	108.87
2	A	601	MES	O1S-S-C8	5.11	111.18	106.79
2	C	601	MES	O1S-S-C8	5.47	111.49	106.79
2	D	601	MES	C5-N4-C3	5.54	121.42	108.87
2	E	601	MES	C5-N4-C3	5.59	121.54	108.87
2	B	601	MES	C5-N4-C3	5.74	121.87	108.87

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	MES	2	0
4	E	603	FMT	1	0
2	F	601	MES	1	0
2	G	601	MES	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	478/478 (100%)	-0.63	2 (0%) 92 96	10, 17, 29, 60	0
1	B	478/478 (100%)	-0.72	1 (0%) 94 97	9, 15, 28, 47	0
1	C	476/478 (99%)	-0.66	0 100 100	9, 15, 29, 52	0
1	D	477/478 (99%)	-0.41	5 (1%) 82 90	11, 18, 33, 66	0
1	E	476/478 (99%)	-0.53	2 (0%) 92 96	11, 18, 31, 66	0
1	F	476/478 (99%)	-0.40	4 (0%) 86 92	11, 21, 38, 58	0
1	G	477/478 (99%)	-0.51	4 (0%) 86 92	11, 19, 34, 69	0
1	H	477/478 (99%)	-0.58	1 (0%) 94 97	9, 16, 30, 57	0
All	All	3815/3824 (99%)	-0.55	19 (0%) 90 95	9, 17, 33, 69	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	7	HIS	6.5
1	A	7	HIS	3.7
1	D	7	HIS	3.6
1	F	350	ALA	3.0
1	D	364	GLY	2.8
1	F	348	ALA	2.6
1	F	349	THR	2.5
1	D	365[A]	PRO	2.5
1	E	400	ASP	2.5
1	G	365	PRO	2.4
1	A	6	ARG	2.4
1	F	352	GLY	2.3
1	G	364	GLY	2.3
1	E	364	GLY	2.3
1	H	7	HIS	2.2
1	D	363	PRO	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	461	PHE	2.2
1	G	346	ALA	2.0
1	B	364	GLY	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
5	UNX	C	603	1/1	0.91	0.71	41.58	30,30,30,30	0
5	UNX	H	604	1/1	0.91	0.63	35.39	30,30,30,30	0
5	UNX	A	604	1/1	0.95	0.55	33.03	30,30,30,30	0
5	UNX	B	604	1/1	0.88	0.45	29.58	30,30,30,30	0
5	UNX	F	603	1/1	0.73	0.53	26.07	30,30,30,30	0
5	UNX	D	603	1/1	0.81	0.45	16.61	30,30,30,30	0
5	UNX	E	604	1/1	0.86	0.41	8.44	30,30,30,30	0
4	FMT	E	603	3/3	0.84	0.24	5.87	35,35,37,40	0
2	MES	H	601	12/12	0.96	0.14	4.92	38,50,55,58	0
4	FMT	B	603	3/3	0.86	0.13	4.86	41,41,42,46	0
5	UNX	G	603	1/1	0.86	0.31	4.50	30,30,30,30	0
2	MES	G	601	12/12	0.96	0.14	4.18	38,56,63,64	0
2	MES	A	601	12/12	0.97	0.14	3.82	20,55,62,63	0
2	MES	F	601	12/12	0.97	0.13	3.46	34,56,62,62	0
2	MES	E	601	12/12	0.97	0.14	2.68	32,48,63,63	0
2	MES	B	601	12/12	0.96	0.13	2.46	23,44,59,61	0
3	GOL	G	602	6/6	0.96	0.15	1.92	17,32,36,39	0
2	MES	D	601	12/12	0.97	0.17	1.57	32,47,56,57	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	MES	C	601	12/12	0.96	0.12	1.28	33,48,55,56	0
3	GOL	B	602	6/6	0.96	0.12	1.10	16,28,33,34	0
3	GOL	E	602	6/6	0.98	0.12	0.75	12,25,27,29	0
3	GOL	F	602	6/6	0.95	0.15	0.43	23,33,35,39	0
3	GOL	D	602	6/6	0.97	0.12	0.08	20,26,31,34	0
3	GOL	C	602	6/6	0.97	0.12	-0.03	17,32,32,36	0
3	GOL	H	602	6/6	0.97	0.14	-0.29	16,28,29,36	0
3	GOL	A	602	6/6	0.98	0.11	-1.04	12,37,38,41	0
4	FMT	H	603	3/3	0.86	0.14	-	28,28,31,34	0
4	FMT	A	603	3/3	0.87	0.26	-	47,47,48,50	0

6.5 Other polymers [i](#)

There are no such residues in this entry.