



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:19 AM GMT

PDB ID : 3HYV  
Title : 3-D X-Ray structure of the sulfide:quinone oxidoreductase from the hyperthermophilic bacterium Aquifex aeolicus  
Authors : Marcia, M.; Ermler, U.; Peng, G.H.; Michel, H.  
Deposited on : 2009-06-23  
Resolution : 2.30 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](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
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) : trunk26865

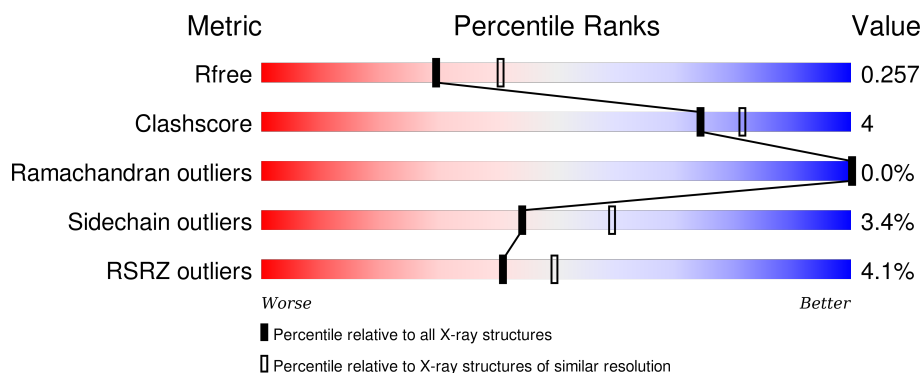
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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))
$R_{free}$	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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  $\geq 3$ , 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	430	<div> <div>3%</div> <div>90%</div> <div>10%</div> </div>
1	B	430	<div> <div>3%</div> <div>90%</div> <div>10%</div> </div>
1	C	430	<div> <div>3%</div> <div>87%</div> <div>12%</div> <div>.</div> </div>
1	D	430	<div> <div>10%</div> <div>89%</div> <div>10%</div> <div>.</div> </div>
1	E	430	<div> <div>4%</div> <div>91%</div> <div>8%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	430	

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
3	LMT	A	600	-	-	-	X
3	LMT	B	600	-	-	-	X
3	LMT	C	600	-	-	-	X
3	LMT	D	600	-	-	-	X
3	LMT	E	600	-	-	-	X
3	LMT	F	600	-	-	-	X
4	H2S	A	700	-	-	-	X
4	H2S	B	700	-	-	-	X
4	H2S	C	700	-	-	-	X
4	H2S	D	700	-	-	-	X
4	H2S	E	700	-	-	-	X
4	H2S	F	700	-	-	-	X
5	PS9	A	800	-	-	X	X
5	PS9	B	800	-	-	X	X
5	PS9	C	800	-	-	-	X
5	PS9	D	802	-	-	-	X
5	PS9	E	800	-	-	-	X
5	PS9	F	800	-	-	-	X
7	SO4	C	432	-	-	X	-

## 2 Entry composition [i](#)

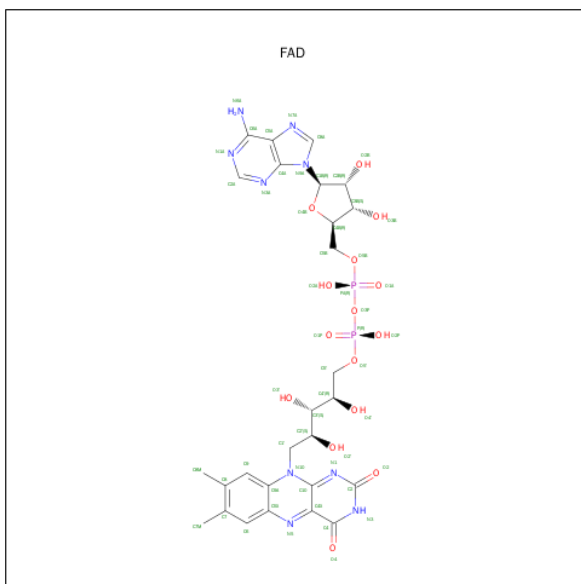
There are 8 unique types of molecules in this entry. The entry contains 21480 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Sulfide-quinone reductase.

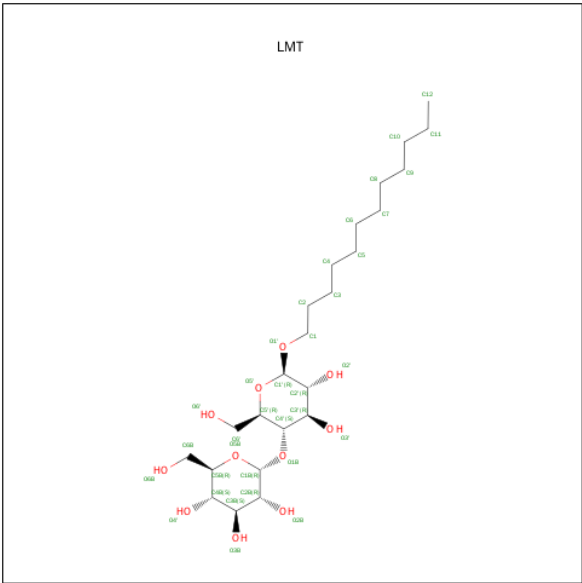
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	429	Total	C	N	O	S	0	8	0
			3370	2182	556	609	23			
1	B	429	Total	C	N	O	S	0	7	0
			3364	2179	556	606	23			
1	C	429	Total	C	N	O	S	0	6	0
			3358	2174	555	606	23			
1	D	429	Total	C	N	O	S	0	6	0
			3357	2175	555	604	23			
1	E	429	Total	C	N	O	S	0	4	0
			3346	2166	555	602	23			
1	F	429	Total	C	N	O	S	0	4	0
			3346	2166	555	602	23			

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula:  $C_{27}H_{33}N_9O_{15}P_2$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	C	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	D	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	E	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	F	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is DODECYL-BETA-D-MALTOSIDE (three-letter code: LMT) (formula: C<sub>24</sub>H<sub>46</sub>O<sub>11</sub>).



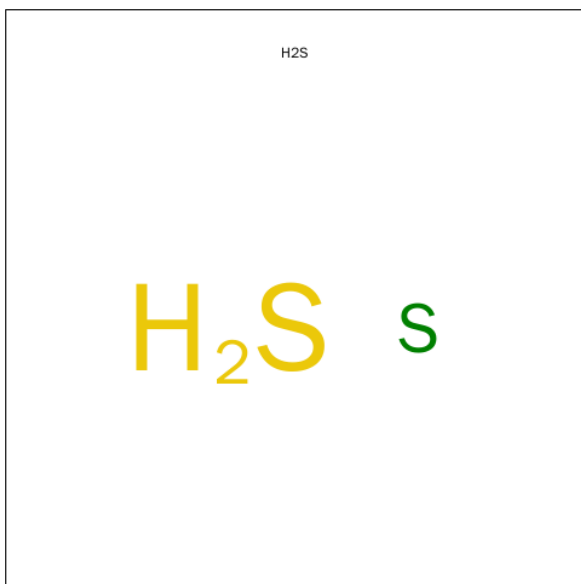
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			35	24	11		
3	B	1	Total	C	O	0	0
			35	24	11		
3	C	1	Total	C	O	0	0
			35	24	11		
3	D	1	Total	C	O	0	0
			35	24	11		
3	E	1	Total	C	O	0	0
			35	24	11		

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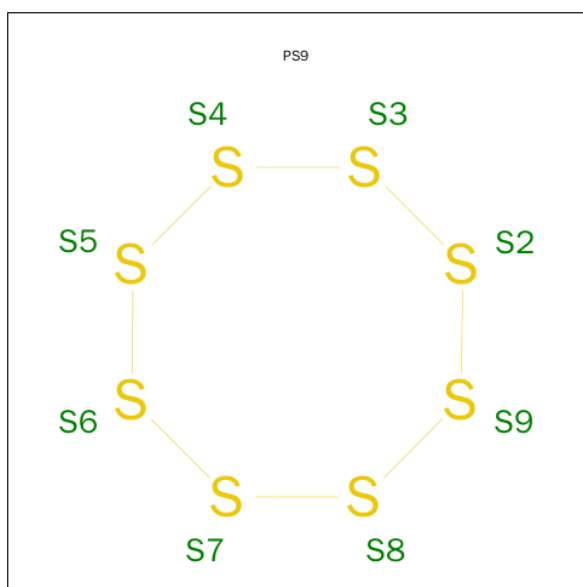
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	F	1	Total	C	O	0	0
			35	24	11		

- Molecule 4 is HYDROSULFURIC ACID (three-letter code: H<sub>2</sub>S) (formula: H<sub>2</sub>S).



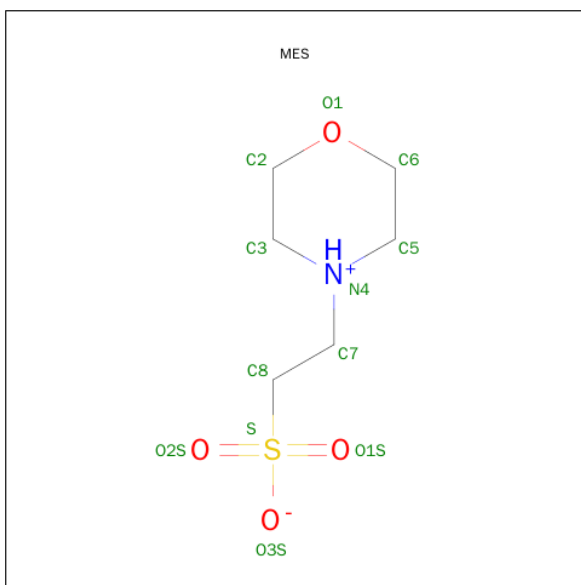
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	S	0	0
			1	1		
4	B	1	Total	S	0	0
			1	1		
4	C	1	Total	S	0	0
			1	1		
4	D	1	Total	S	0	0
			1	1		
4	E	1	Total	S	0	0
			1	1		
4	F	1	Total	S	0	0
			1	1		

- Molecule 5 is OCTATHIOCANES (three-letter code: PS<sub>9</sub>) (formula: S<sub>8</sub>).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total S 8 8	0	0
5	B	1	Total S 6 6	0	0
5	C	1	Total S 6 6	0	0
5	D	1	Total S 1 1	0	0
5	D	1	Total S 5 5	0	0
5	E	1	Total S 8 8	0	0
5	F	1	Total S 8 8	0	0

- Molecule 6 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula:  $C_6H_{13}NO_4S$ ).



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

- Molecule 7 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).





Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	O	S	0	0
			5	4	1		
7	A	1	Total	O	S	0	0
			5	4	1		
7	A	1	Total	O	S	0	0
			5	4	1		
7	A	1	Total	O	S	0	0
			5	4	1		
7	B	1	Total	O	S	0	0
			5	4	1		
7	B	1	Total	O	S	0	0
			5	4	1		
7	B	1	Total	O	S	0	0
			5	4	1		
7	B	1	Total	O	S	0	0
			5	4	1		
7	C	1	Total	O	S	0	0
			5	4	1		
7	C	1	Total	O	S	0	0
			5	4	1		
7	C	1	Total	O	S	0	0
			5	4	1		
7	C	1	Total	O	S	0	0
			5	4	1		
7	D	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	D	1	Total	O	S	0	0
			5	4	1		
7	D	1	Total	O	S	0	0
			5	4	1		
7	E	1	Total	O	S	0	0
			5	4	1		
7	E	1	Total	O	S	0	0
			5	4	1		
7	E	1	Total	O	S	0	0
			5	4	1		
7	E	1	Total	O	S	0	0
			5	4	1		
7	F	1	Total	O	S	0	0
			5	4	1		
7	F	1	Total	O	S	0	0
			5	4	1		
7	F	1	Total	O	S	0	0
			5	4	1		

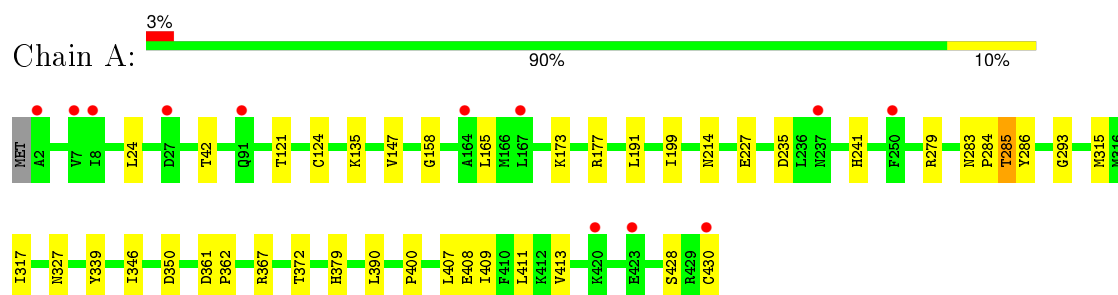
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	97	Total	O	0	0
			97	97		
8	B	107	Total	O	0	0
			107	107		
8	C	105	Total	O	0	0
			105	105		
8	D	57	Total	O	0	0
			57	57		
8	E	92	Total	O	0	0
			92	92		
8	F	113	Total	O	0	0
			113	113		

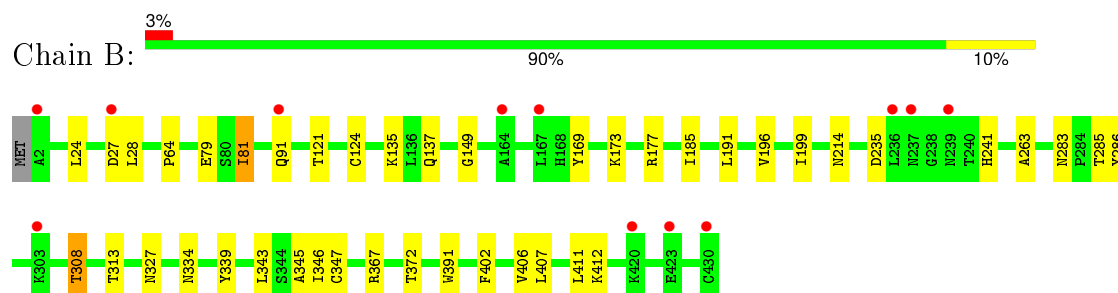
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

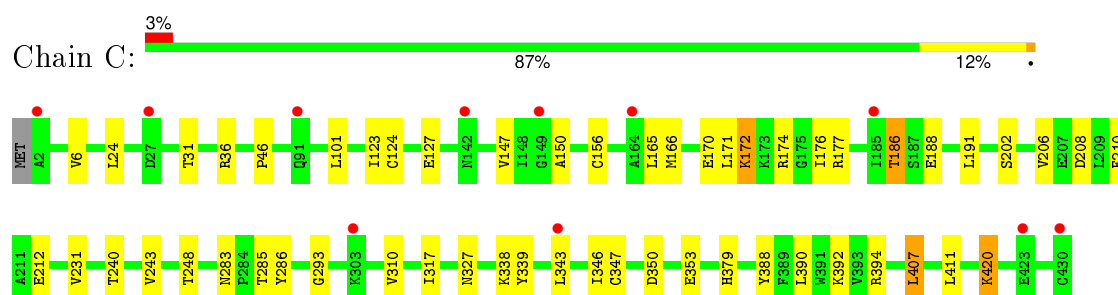
- Molecule 1: Sulfide-quinone reductase



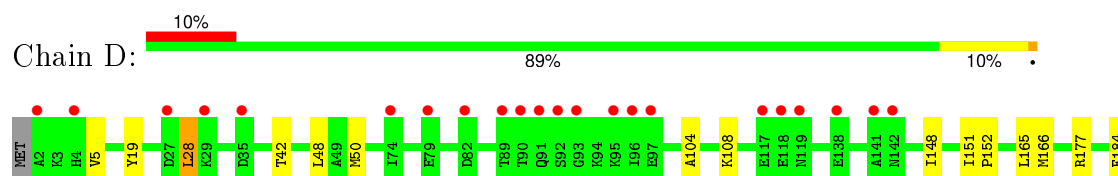
- Molecule 1: Sulfide-quinone reductase

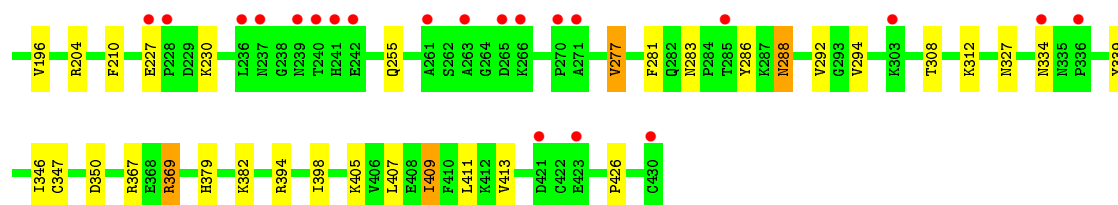


- Molecule 1: Sulfide-quinone reductase

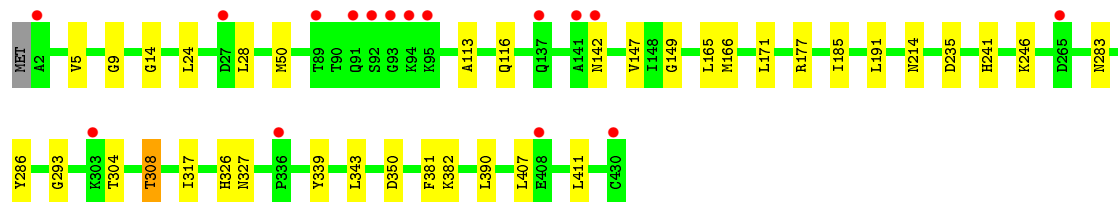


- Molecule 1: Sulfide-quinone reductase

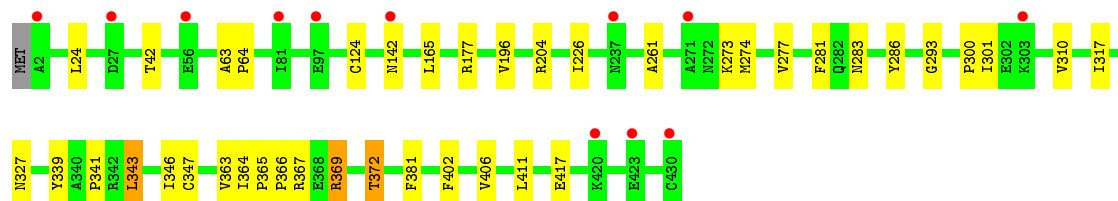
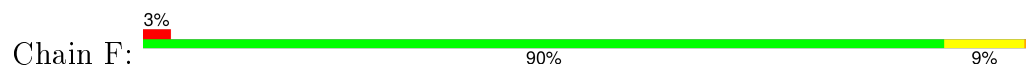




• Molecule 1: Sulfide-quinone reductase



• Molecule 1: Sulfide-quinone reductase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	112.85Å 154.92Å 178.04Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.04 – 2.30 20.04 – 2.30	Depositor EDS
% Data completeness (in resolution range)	97.9 (20.04-2.30) 97.9 (20.04-2.30)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.96 (at 2.30Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.193 , 0.235 0.217 , 0.257	Depositor DCC
$R_{free}$ test set	6821 reflections (5.30%)	DCC
Wilson B-factor (Å <sup>2</sup> )	38.7	Xtriage
Anisotropy	0.124	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 49.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 135604 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	21480	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.18% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: CSS, LMT, H2S, PS9, SO4, MES, FAD

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	0.52	0/3477	0.57	0/4720
1	B	0.50	0/3468	0.58	0/4707
1	C	0.50	0/3459	0.58	0/4696
1	D	0.47	0/3458	0.54	0/4695
1	E	0.48	0/3441	0.57	0/4672
1	F	0.50	0/3441	0.57	0/4672
All	All	0.49	0/20744	0.57	0/28162

There are no bond length outliers.

There are no bond angle outliers.

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	3370	0	3373	25	0
1	B	3364	0	3372	26	0
1	C	3358	0	3359	33	0
1	D	3357	0	3364	31	0
1	E	3346	0	3347	18	0
1	F	3346	0	3347	27	0
2	A	53	0	30	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	53	0	29	0	0
2	C	53	0	29	0	0
2	D	53	0	29	1	0
2	E	53	0	29	1	0
2	F	53	0	29	1	0
3	A	35	0	46	0	0
3	B	35	0	46	0	0
3	C	35	0	46	2	0
3	D	35	0	46	0	0
3	E	35	0	46	8	0
3	F	35	0	46	9	0
4	A	1	0	0	1	0
4	B	1	0	0	1	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
4	E	1	0	0	0	0
4	F	1	0	0	1	0
5	A	8	0	0	2	0
5	B	6	0	0	2	0
5	C	6	0	0	1	0
5	D	6	0	0	1	0
5	E	8	0	0	0	0
5	F	8	0	0	1	0
6	A	12	0	12	0	0
6	B	12	0	12	0	0
6	C	12	0	12	1	0
6	D	12	0	12	0	0
6	E	12	0	12	0	0
6	F	12	0	12	0	0
7	A	20	0	0	0	0
7	B	25	0	0	1	0
7	C	20	0	0	2	0
7	D	15	0	0	1	0
7	E	25	0	0	0	0
7	F	15	0	0	1	0
8	A	97	0	0	2	0
8	B	107	0	0	3	0
8	C	105	0	0	4	0
8	D	57	0	0	0	0
8	E	92	0	0	1	0
8	F	113	0	0	1	0
All	All	21480	0	20685	168	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 4.

All (168) 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:D:347[B]:CYS:SG	5:D:802:PS9:S9	2.39	1.19
1:C:420:LYS:H	1:C:420:LYS:HD3	1.28	0.98
3:E:600:LMT:H1B	3:F:600:LMT:O2B	1.64	0.97
1:D:327:ASN:HD21	1:D:339:TYR:H	1.17	0.92
3:F:600:LMT:O2B	3:F:600:LMT:H4'	1.72	0.89
1:C:327:ASN:HD21	1:C:339:TYR:H	1.24	0.84
1:E:327:ASN:HD21	1:E:339:TYR:H	1.24	0.84
3:E:600:LMT:C1B	3:F:600:LMT:O2B	2.26	0.84
1:C:420:LYS:H	1:C:420:LYS:CD	1.91	0.83
1:C:283:ASN:HD22	1:C:286:TYR:H	1.27	0.82
1:F:283:ASN:HD22	1:F:286:TYR:H	1.28	0.79
1:A:283:ASN:HD22	1:A:286:TYR:H	1.31	0.77
1:E:283:ASN:HD22	1:E:286:TYR:H	1.29	0.77
1:B:283:ASN:HD22	1:B:286:TYR:H	1.29	0.75
1:F:274:MET:HE1	8:F:822:HOH:O	1.87	0.74
1:F:42:THR:HG21	2:F:441:FAD:O4'	1.88	0.73
1:F:327:ASN:HD21	1:F:339:TYR:H	1.35	0.72
1:B:327:ASN:HD21	1:B:339:TYR:H	1.39	0.70
3:E:600:LMT:H2B	3:F:600:LMT:O2B	1.92	0.69
1:C:186:THR:HG21	1:C:188:GLU:OE1	1.93	0.69
3:E:600:LMT:C2B	3:F:600:LMT:O2B	2.41	0.68
1:A:327:ASN:HD21	1:A:339:TYR:H	1.41	0.68
1:A:124[A]:CYS:SG	4:A:700:H2S:S	2.67	0.64
1:B:169:TYR:CE2	1:B:173[B]:LYS:HD2	2.33	0.63
1:F:346:ILE:HD11	1:F:411:LEU:CD1	2.29	0.63
1:B:196:VAL:HG22	1:B:367:ARG:NH1	2.14	0.62
1:B:285:THR:HG23	8:B:451:HOH:O	1.99	0.61
1:A:367:ARG:H	1:E:214:ASN:HD22	1.47	0.61
1:D:346:ILE:HD11	1:D:411:LEU:CD1	2.31	0.61
1:D:288:ASN:H	1:D:288:ASN:HD22	1.47	0.61
1:B:81:ILE:HG23	1:B:263:ALA:HB2	1.82	0.61
1:B:199:ILE:O	1:B:372:THR:HG21	2.00	0.60
1:D:196:VAL:HG13	1:D:367:ARG:NH1	2.16	0.60
1:E:283:ASN:ND2	1:E:286:TYR:H	1.99	0.60
1:D:346:ILE:HD11	1:D:411:LEU:HD13	1.84	0.60
1:A:367:ARG:H	1:E:214:ASN:ND2	2.01	0.59
1:E:411:LEU:HD21	3:E:600:LMT:H62	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:350:ASP:OD1	1:C:379:HIS:HD2	1.88	0.56
1:E:327:ASN:HD21	1:E:339:TYR:N	1.99	0.56
1:A:283:ASN:ND2	1:A:286:TYR:H	2.02	0.56
1:F:402:PHE:O	1:F:406:VAL:HG23	2.06	0.56
1:D:283:ASN:HD22	1:D:286:TYR:H	1.51	0.56
1:A:285:THR:HG23	8:A:438:HOH:O	2.04	0.56
1:F:124[A]:CYS:SG	4:F:700:H2S:S	2.91	0.56
1:D:327:ASN:HD21	1:D:339:TYR:N	1.96	0.56
1:C:147:VAL:HG11	1:C:231:VAL:HG21	1.88	0.56
1:D:350:ASP:OD1	1:D:379:HIS:HD2	1.88	0.55
1:C:166:MET:HE3	1:C:350:ASP:HB3	1.89	0.55
1:A:199:ILE:O	1:A:372:THR:HG21	2.07	0.55
1:C:394:ARG:NH1	7:C:432:SO4:S	2.79	0.55
1:D:196:VAL:CG1	1:D:367:ARG:NH1	2.70	0.55
1:D:19:TYR:CD2	1:D:398:ILE:HG12	2.42	0.54
1:C:394:ARG:NH1	7:C:432:SO4:O1	2.41	0.53
1:F:346:ILE:HD11	1:F:411:LEU:HD11	1.90	0.53
1:C:420:LYS:CD	1:C:420:LYS:N	2.69	0.53
1:A:293:GLY:HA2	1:A:317:ILE:HG12	1.91	0.53
3:F:600:LMT:H2O1	3:F:600:LMT:H4'	1.74	0.53
1:D:19:TYR:HD2	1:D:398:ILE:HG12	1.74	0.52
1:B:124[A]:CYS:SG	4:B:700:H2S:S	2.98	0.52
1:B:346:ILE:HD11	1:B:411:LEU:CD1	2.39	0.52
1:F:196:VAL:HG22	1:F:367:ARG:NH1	2.25	0.52
1:E:293:GLY:HA2	1:E:317:ILE:HG12	1.91	0.52
1:D:108:LYS:HB3	1:D:255:GLN:O	2.10	0.52
1:D:294:VAL:HA	1:D:312:LYS:HD3	1.91	0.52
1:B:196:VAL:HG22	1:B:367:ARG:HH11	1.74	0.51
1:C:293:GLY:HA2	1:C:317:ILE:HG12	1.92	0.51
3:E:600:LMT:H6'1	3:F:600:LMT:H2'	1.92	0.51
1:D:50:MET:HG3	1:D:166:MET:HE2	1.92	0.51
3:E:600:LMT:H2B	3:F:600:LMT:H2O1	1.77	0.50
1:D:394:ARG:NH1	7:D:432:SO4:O3	2.44	0.50
1:B:235:ASP:OD2	1:B:241:HIS:HE1	1.94	0.50
1:E:350:ASP:HB2	1:E:382:LYS:HD3	1.94	0.50
1:F:283:ASN:ND2	1:F:286:TYR:H	2.03	0.50
1:E:327:ASN:ND2	1:E:339:TYR:H	2.02	0.50
1:F:300:PRO:HA	1:F:310:VAL:HG23	1.94	0.50
1:C:6:VAL:HG12	1:C:31:THR:CG2	2.41	0.49
1:D:227:GLU:HB2	1:D:230:LYS:HG2	1.95	0.49
1:F:327:ASN:HD21	1:F:339:TYR:N	2.07	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:121:THR:CG2	1:B:135:LYS:HD3	2.43	0.48
1:B:283:ASN:ND2	1:B:286:TYR:H	2.05	0.48
1:A:121:THR:CG2	1:A:135:LYS:HD3	2.43	0.48
1:D:405:LYS:O	1:D:409:ILE:HG23	2.13	0.48
1:D:196:VAL:CG1	1:D:367:ARG:HH12	2.26	0.48
1:C:283:ASN:ND2	1:C:286:TYR:H	2.05	0.48
1:E:304:THR:OG1	1:E:308:THR:HB	2.13	0.48
1:B:327:ASN:HD21	1:B:339:TYR:N	2.09	0.47
1:E:326:HIS:HD2	8:E:451:HOH:O	1.97	0.47
1:C:346:ILE:HD11	1:C:411:LEU:CD1	2.45	0.47
1:A:158:GLY:HA3	5:A:800:PS9:S7	2.55	0.47
1:C:165:LEU:HD21	1:C:210:PHE:HE1	1.79	0.47
1:C:208:ASP:O	1:C:212:GLU:HG3	2.15	0.47
1:B:313:THR:HG21	1:B:345:ALA:HB1	1.98	0.46
1:D:277:VAL:HG22	1:D:281:PHE:C	2.36	0.46
1:E:235:ASP:OD2	1:E:241:HIS:HE1	1.99	0.46
1:C:202:SER:O	1:C:206:VAL:HG23	2.15	0.46
1:D:104:ALA:HB2	1:D:292:VAL:HG22	1.98	0.46
1:E:50:MET:HG3	1:E:166:MET:HE2	1.98	0.46
1:B:308:THR:HG22	8:B:476:HOH:O	2.16	0.45
1:B:149:GLY:HA3	1:B:185:ILE:O	2.16	0.45
1:C:147:VAL:HG11	1:C:231:VAL:CG2	2.47	0.45
1:D:369:ARG:NH1	7:F:431:SO4:O2	2.50	0.45
1:C:36:ARG:NH1	1:C:127:GLU:OE1	2.49	0.44
1:C:407:LEU:HD13	3:C:600:LMT:H92	1.99	0.44
1:A:350:ASP:OD1	1:A:379:HIS:HD2	2.00	0.44
1:C:353:GLU:CG	6:C:900:MES:H51	2.47	0.44
1:D:165:LEU:HD21	1:D:210:PHE:HE1	1.82	0.44
1:B:137:GLN:NE2	8:B:457:HOH:O	2.51	0.44
1:F:347[A]:CYS:HB3	5:F:800:PS9:S6	2.58	0.44
1:A:346:ILE:HD11	1:A:411:LEU:CD1	2.48	0.44
3:E:600:LMT:H3'	3:F:600:LMT:H6E	2.00	0.43
1:F:196:VAL:HG21	1:F:363:VAL:HG13	1.99	0.43
1:C:6:VAL:HG23	1:C:101:LEU:CD1	2.48	0.43
1:F:293:GLY:HA2	1:F:317:ILE:HG12	1.99	0.43
1:D:204:ARG:CZ	1:F:204:ARG:HD3	2.48	0.43
1:D:350:ASP:HB2	1:D:382:LYS:HD3	2.00	0.43
1:B:347[B]:CYS:SG	5:B:800:PS9:S5	3.15	0.43
1:C:147:VAL:HG12	1:C:248:THR:HG22	1.99	0.43
1:A:121:THR:HG22	1:A:135:LYS:HD3	2.00	0.43
1:D:5:VAL:HG12	1:D:28:LEU:HD22	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:142:ASN:O	1:E:246:LYS:HE2	2.19	0.43
1:F:372:THR:CG2	1:F:372:THR:O	2.66	0.43
1:B:173[A]:LYS:HA	1:B:173[A]:LYS:HD3	1.86	0.43
1:F:341:PRO:HB2	1:F:343:LEU:HD13	2.01	0.42
1:C:124[A]:CYS:HB2	8:C:743:HOH:O	2.19	0.42
1:B:214:ASN:HD21	1:F:366:PRO:HA	1.83	0.42
1:F:261:ALA:HB2	1:F:273:LYS:HD3	2.00	0.42
1:E:113:ALA:HB3	1:E:116:GLN:HB2	2.02	0.42
1:B:402:PHE:O	1:B:406:VAL:HG23	2.19	0.42
1:C:46:PRO:HG3	1:C:123:ILE:HD12	2.01	0.42
1:A:283:ASN:HA	1:A:284:PRO:HD3	1.89	0.42
1:B:347[A]:CYS:HB3	5:B:800:PS9:S5	2.60	0.42
1:D:151:ILE:HB	1:D:152:PRO:HD2	2.02	0.42
1:A:158:GLY:HA3	5:A:800:PS9:S4	2.60	0.42
1:B:27:ASP:OD1	1:B:27:ASP:N	2.53	0.42
7:B:435:SO4:O3	1:F:369:ARG:HD3	2.20	0.42
1:D:50:MET:HG3	1:D:166:MET:CE	2.50	0.42
1:A:315:MET:CE	1:A:346:ILE:HD12	2.50	0.42
1:F:196:VAL:HG23	1:F:301:ILE:HD13	2.02	0.41
1:A:279:ARG:HD2	1:A:428:SER:OG	2.20	0.41
1:C:172:LYS:HE3	8:C:624:HOH:O	2.18	0.41
1:D:277:VAL:O	1:D:426:PRO:HD3	2.19	0.41
2:D:441:FAD:H9	2:D:441:FAD:H1'1	1.87	0.41
1:C:310:VAL:HG12	8:C:784:HOH:O	2.21	0.41
1:A:400:PRO:HB3	1:B:391:TRP:CZ2	2.55	0.41
1:D:148:ILE:O	1:D:184:PHE:HA	2.19	0.41
1:C:170:GLU:O	1:C:174:ARG:HG2	2.19	0.41
1:F:196:VAL:HG22	1:F:367:ARG:HH11	1.84	0.41
1:C:156[A]:CSS:SG	5:C:800:PS9:S2	3.18	0.41
2:A:441:FAD:H1'1	2:A:441:FAD:H9	1.77	0.41
1:A:407:LEU:HB3	1:A:413:VAL:CG1	2.51	0.41
1:A:361:ASP:HA	1:A:362:PRO:HA	1.94	0.41
1:C:124[B]:CYS:HB2	8:C:743:HOH:O	2.21	0.41
2:E:441:FAD:H9	2:E:441:FAD:H1'1	1.85	0.41
1:C:388:TYR:CZ	1:C:392:LYS:HD2	2.56	0.41
1:E:9:GLY:O	1:E:14:GLY:HA3	2.21	0.41
1:B:79:GLU:HG3	1:B:91:GLN:HA	2.03	0.41
1:F:277:VAL:HB	1:F:281:PHE:HA	2.03	0.41
1:F:372:THR:O	1:F:372:THR:HG23	2.21	0.40
1:A:279:ARG:NH1	1:A:430:CYS:HA	2.37	0.40
1:E:149:GLY:HA3	1:E:185:ILE:O	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:150:ALA:HB3	1:C:186:THR:HG22	2.03	0.40
1:A:124[A]:CYS:HB2	8:A:510:HOH:O	2.22	0.40
1:D:151:ILE:HB	1:D:152:PRO:CD	2.52	0.40
1:A:173:LYS:HD3	1:A:173:LYS:HA	1.92	0.40
1:F:364:ILE:HA	1:F:365:PRO:HD3	1.95	0.40
3:C:600:LMT:O2B	3:C:600:LMT:H6E	2.22	0.40
1:A:235:ASP:OD2	1:A:241:HIS:HE1	2.04	0.40
1:F:63:ALA:HB3	1:F:64:PRO:HD3	2.03	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	433/430 (101%)	427 (99%)	6 (1%)	0	100	100
1	B	432/430 (100%)	426 (99%)	6 (1%)	0	100	100
1	C	431/430 (100%)	423 (98%)	8 (2%)	0	100	100
1	D	431/430 (100%)	421 (98%)	9 (2%)	1 (0%)	52	64
1	E	429/430 (100%)	417 (97%)	12 (3%)	0	100	100
1	F	429/430 (100%)	422 (98%)	7 (2%)	0	100	100
All	All	2585/2580 (100%)	2536 (98%)	48 (2%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	334	ASN

### 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	362/356 (102%)	350 (97%)	12 (3%)	45	61
1	B	361/356 (101%)	350 (97%)	11 (3%)	48	65
1	C	360/356 (101%)	343 (95%)	17 (5%)	32	43
1	D	360/356 (101%)	349 (97%)	11 (3%)	47	64
1	E	358/356 (101%)	345 (96%)	13 (4%)	42	57
1	F	358/356 (101%)	348 (97%)	10 (3%)	51	68
All	All	2159/2136 (101%)	2085 (97%)	74 (3%)	44	59

All (74) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	LEU
1	A	42	THR
1	A	147	VAL
1	A	165	LEU
1	A	177	ARG
1	A	191	LEU
1	A	214	ASN
1	A	227	GLU
1	A	285	THR
1	A	390	LEU
1	A	408	GLU
1	A	409	ILE
1	B	24	LEU
1	B	28	LEU
1	B	64	PRO
1	B	81	ILE
1	B	177	ARG
1	B	191	LEU
1	B	308	THR
1	B	334	ASN
1	B	343	LEU
1	B	407	LEU

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Mol	Chain	Res	Type
1	B	412	LYS
1	C	24	LEU
1	C	171	LEU
1	C	172	LYS
1	C	176	ILE
1	C	177	ARG
1	C	186	THR
1	C	191	LEU
1	C	240	THR
1	C	243	VAL
1	C	285	THR
1	C	338	LYS
1	C	343	LEU
1	C	347[A]	CYS
1	C	347[B]	CYS
1	C	390	LEU
1	C	407	LEU
1	C	420	LYS
1	D	28	LEU
1	D	42	THR
1	D	48	LEU
1	D	177	ARG
1	D	277	VAL
1	D	288	ASN
1	D	308	THR
1	D	369	ARG
1	D	407	LEU
1	D	409	ILE
1	D	413	VAL
1	E	5	VAL
1	E	24	LEU
1	E	28	LEU
1	E	147	VAL
1	E	165	LEU
1	E	171	LEU
1	E	177	ARG
1	E	191	LEU
1	E	308	THR
1	E	343	LEU
1	E	381	PHE
1	E	390	LEU
1	E	407	LEU

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Mol	Chain	Res	Type
1	F	24	LEU
1	F	142	ASN
1	F	165	LEU
1	F	177	ARG
1	F	226	ILE
1	F	343	LEU
1	F	369	ARG
1	F	372	THR
1	F	381	PHE
1	F	417	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (43) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	ASN
1	A	241	HIS
1	A	283	ASN
1	A	327	ASN
1	A	379	HIS
1	B	128	HIS
1	B	214	ASN
1	B	241	HIS
1	B	283	ASN
1	B	327	ASN
1	B	334	ASN
1	B	379	HIS
1	C	23	ASN
1	C	128	HIS
1	C	133	GLN
1	C	137	GLN
1	C	214	ASN
1	C	237	ASN
1	C	241	HIS
1	C	283	ASN
1	C	327	ASN
1	C	379	HIS
1	C	395	ASN
1	D	70	ASN
1	D	214	ASN
1	D	283	ASN
1	D	288	ASN
1	D	327	ASN

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Mol	Chain	Res	Type
1	D	379	HIS
1	E	128	HIS
1	E	214	ASN
1	E	237	ASN
1	E	241	HIS
1	E	283	ASN
1	E	327	ASN
1	E	379	HIS
1	E	395	ASN
1	F	128	HIS
1	F	133	GLN
1	F	214	ASN
1	F	241	HIS
1	F	283	ASN
1	F	327	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

12 non-standard protein/DNA/RNA residues are 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
1	CSS	A	156[A]	-	4,5,7	1.15	0	3,5,8	1.62	1 (33%)
1	CSS	A	156[B]	5	4,6,7	0.78	0	3,6,8	1.64	1 (33%)
1	CSS	B	156[A]	-	4,5,7	1.14	0	3,5,8	1.51	1 (33%)
1	CSS	B	156[B]	5	4,6,7	0.88	0	3,6,8	2.26	2 (66%)
1	CSS	C	156[A]	-	4,5,7	1.20	0	3,5,8	1.44	0
1	CSS	C	156[B]	5	4,6,7	0.96	0	3,6,8	1.27	0
1	CSS	D	156[A]	-	4,5,7	1.38	1 (25%)	3,5,8	2.20	2 (66%)



Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
1	CSS	D	156[B]	5	4,6,7	0.91	0	3,6,8	1.60	1 (33%)
1	CSS	E	156[A]	-	4,5,7	1.33	0	3,5,8	1.48	1 (33%)
1	CSS	E	156[B]	5	4,6,7	0.91	0	3,6,8	1.53	1 (33%)
1	CSS	F	156[A]	-	4,5,7	1.30	0	3,5,8	1.47	1 (33%)
1	CSS	F	156[B]	5	4,6,7	0.96	0	3,6,8	1.91	1 (33%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	CSS	A	156[A]	-	-	0/1/4/7	0/0/0/0
1	CSS	A	156[B]	5	-	0/1/5/7	0/0/0/0
1	CSS	B	156[A]	-	-	0/1/4/7	0/0/0/0
1	CSS	B	156[B]	5	-	0/1/5/7	0/0/0/0
1	CSS	C	156[A]	-	-	0/1/4/7	0/0/0/0
1	CSS	C	156[B]	5	-	0/1/5/7	0/0/0/0
1	CSS	D	156[A]	-	-	0/1/4/7	0/0/0/0
1	CSS	D	156[B]	5	-	0/1/5/7	0/0/0/0
1	CSS	E	156[A]	-	-	0/1/4/7	0/0/0/0
1	CSS	E	156[B]	5	-	0/1/5/7	0/0/0/0
1	CSS	F	156[A]	-	-	0/1/4/7	0/0/0/0
1	CSS	F	156[B]	5	-	0/1/5/7	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	156[A]	CSS	CB-SG	-2.42	1.76	1.81

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	156[A]	CSS	CA-CB-SG	-2.82	108.02	114.48
1	A	156[A]	CSS	O-C-CA	-2.78	118.26	125.49
1	A	156[B]	CSS	O-C-CA	-2.78	118.26	125.49
1	D	156[B]	CSS	O-C-CA	-2.52	118.93	125.49
1	D	156[A]	CSS	O-C-CA	-2.52	118.93	125.49
1	F	156[B]	CSS	O-C-CA	-2.49	119.00	125.49
1	F	156[A]	CSS	O-C-CA	-2.49	119.00	125.49
1	B	156[B]	CSS	O-C-CA	-2.46	119.08	125.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	156[A]	CSS	O-C-CA	-2.46	119.08	125.49
1	E	156[A]	CSS	O-C-CA	-2.45	119.10	125.49
1	E	156[B]	CSS	O-C-CA	-2.45	119.10	125.49
1	B	156[B]	CSS	CB-SG-SD	3.04	109.91	103.94

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	C	156[A]	CSS	1	0

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 55 ligands modelled in this entry, 7 are modelled with single atom - leaving 48 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
7	SO4	A	431	-	4,4,4	0.26	0	6,6,6	0.08	0
7	SO4	A	432	-	4,4,4	0.24	0	6,6,6	0.15	0
7	SO4	A	433	-	4,4,4	0.25	0	6,6,6	0.27	0
7	SO4	A	434	-	4,4,4	0.17	0	6,6,6	0.13	0
2	FAD	A	441	4	48,58,58	1.04	3 (6%)	54,89,89	1.94	6 (11%)
3	LMT	A	600	-	36,36,36	0.52	0	47,47,47	0.93	2 (4%)
5	PS9	A	800	1	8,8,8	0.68	0	8,8,8	1.16	0
6	MES	A	900	-	11,12,12	0.61	0	14,16,16	2.62	7 (50%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
7	SO4	B	431	-	4,4,4	0.26	0	6,6,6	0.20	0
7	SO4	B	432	-	4,4,4	0.19	0	6,6,6	0.18	0
7	SO4	B	433	-	4,4,4	0.14	0	6,6,6	0.29	0
7	SO4	B	434	-	4,4,4	0.21	0	6,6,6	0.10	0
7	SO4	B	435	-	4,4,4	0.29	0	6,6,6	0.09	0
2	FAD	B	441	4	48,58,58	1.00	2 (4%)	54,89,89	1.92	7 (12%)
3	LMT	B	600	-	36,36,36	0.59	1 (2%)	47,47,47	1.46	4 (8%)
5	PS9	B	800	1	3,5,8	1.59	0	2,4,8	0.71	0
6	MES	B	900	-	11,12,12	0.63	0	14,16,16	2.64	6 (42%)
7	SO4	C	431	-	4,4,4	0.26	0	6,6,6	0.19	0
7	SO4	C	432	-	4,4,4	0.28	0	6,6,6	0.20	0
7	SO4	C	433	-	4,4,4	0.07	0	6,6,6	0.36	0
7	SO4	C	434	-	4,4,4	0.17	0	6,6,6	0.20	0
2	FAD	C	441	4	48,58,58	1.06	2 (4%)	54,89,89	2.01	8 (14%)
3	LMT	C	600	-	36,36,36	0.50	0	47,47,47	1.04	4 (8%)
5	PS9	C	800	1	3,5,8	1.60	0	2,4,8	1.03	0
6	MES	C	900	-	11,12,12	0.62	0	14,16,16	2.78	5 (35%)
7	SO4	D	432	-	4,4,4	0.23	0	6,6,6	0.18	0
7	SO4	D	433	-	4,4,4	0.23	0	6,6,6	0.09	0
7	SO4	D	434	-	4,4,4	0.16	0	6,6,6	0.21	0
2	FAD	D	441	4	48,58,58	1.07	3 (6%)	54,89,89	2.01	8 (14%)
3	LMT	D	600	-	36,36,36	0.47	0	47,47,47	0.89	1 (2%)
5	PS9	D	802	-	2,4,8	1.59	0	1,3,8	1.31	0
6	MES	D	900	-	11,12,12	0.58	0	14,16,16	2.21	3 (21%)
7	SO4	E	431	-	4,4,4	0.17	0	6,6,6	0.11	0
7	SO4	E	432	-	4,4,4	0.22	0	6,6,6	0.13	0
7	SO4	E	433	-	4,4,4	0.20	0	6,6,6	0.11	0
7	SO4	E	434	-	4,4,4	0.20	0	6,6,6	0.16	0
7	SO4	E	435	-	4,4,4	0.25	0	6,6,6	0.08	0
2	FAD	E	441	4	48,58,58	1.12	3 (6%)	54,89,89	2.14	8 (14%)
3	LMT	E	600	-	36,36,36	0.51	0	47,47,47	0.99	3 (6%)
5	PS9	E	800	1	8,8,8	0.64	0	8,8,8	0.99	0
6	MES	E	900	-	11,12,12	0.60	0	14,16,16	2.56	6 (42%)
7	SO4	F	431	-	4,4,4	0.23	0	6,6,6	0.07	0
7	SO4	F	433	-	4,4,4	0.12	0	6,6,6	0.21	0
7	SO4	F	434	-	4,4,4	0.17	0	6,6,6	0.12	0
2	FAD	F	441	4	48,58,58	1.05	2 (4%)	54,89,89	1.97	7 (12%)
3	LMT	F	600	-	36,36,36	0.48	0	47,47,47	0.99	1 (2%)
5	PS9	F	800	1	8,8,8	0.67	0	8,8,8	1.00	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
6	MES	F	900	-	11,12,12	0.61	0	14,16,16	2.78	7 (50%)

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
7	SO4	A	431	-	-	0/0/0/0	0/0/0/0
7	SO4	A	432	-	-	0/0/0/0	0/0/0/0
7	SO4	A	433	-	-	0/0/0/0	0/0/0/0
7	SO4	A	434	-	-	0/0/0/0	0/0/0/0
2	FAD	A	441	4	-	0/30/50/50	0/6/6/6
3	LMT	A	600	-	-	0/21/61/61	0/2/2/2
5	PS9	A	800	1	-	0/0/8/8	0/1/1/1
6	MES	A	900	-	-	0/6/14/14	0/1/1/1
7	SO4	B	431	-	-	0/0/0/0	0/0/0/0
7	SO4	B	432	-	-	0/0/0/0	0/0/0/0
7	SO4	B	433	-	-	0/0/0/0	0/0/0/0
7	SO4	B	434	-	-	0/0/0/0	0/0/0/0
7	SO4	B	435	-	-	0/0/0/0	0/0/0/0
2	FAD	B	441	4	-	0/30/50/50	0/6/6/6
3	LMT	B	600	-	-	0/21/61/61	0/2/2/2
5	PS9	B	800	1	-	0/3/3/8	0/0/0/1
6	MES	B	900	-	-	0/6/14/14	0/1/1/1
7	SO4	C	431	-	-	0/0/0/0	0/0/0/0
7	SO4	C	432	-	-	0/0/0/0	0/0/0/0
7	SO4	C	433	-	-	0/0/0/0	0/0/0/0
7	SO4	C	434	-	-	0/0/0/0	0/0/0/0
2	FAD	C	441	4	-	0/30/50/50	0/6/6/6
3	LMT	C	600	-	-	0/21/61/61	0/2/2/2
5	PS9	C	800	1	-	0/3/3/8	0/0/0/1
6	MES	C	900	-	-	0/6/14/14	0/1/1/1
7	SO4	D	432	-	-	0/0/0/0	0/0/0/0
7	SO4	D	433	-	-	0/0/0/0	0/0/0/0
7	SO4	D	434	-	-	0/0/0/0	0/0/0/0
2	FAD	D	441	4	-	0/30/50/50	0/6/6/6
3	LMT	D	600	-	-	0/21/61/61	0/2/2/2
5	PS9	D	802	-	-	0/2/2/8	0/0/0/1
6	MES	D	900	-	-	0/6/14/14	0/1/1/1
7	SO4	E	431	-	-	0/0/0/0	0/0/0/0
7	SO4	E	432	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	SO4	E	433	-	-	0/0/0/0	0/0/0/0
7	SO4	E	434	-	-	0/0/0/0	0/0/0/0
7	SO4	E	435	-	-	0/0/0/0	0/0/0/0
2	FAD	E	441	4	-	0/30/50/50	0/6/6/6
3	LMT	E	600	-	-	0/21/61/61	0/2/2/2
5	PS9	E	800	1	-	0/0/8/8	0/1/1/1
6	MES	E	900	-	-	0/6/14/14	0/1/1/1
7	SO4	F	431	-	-	0/0/0/0	0/0/0/0
7	SO4	F	433	-	-	0/0/0/0	0/0/0/0
7	SO4	F	434	-	-	0/0/0/0	0/0/0/0
2	FAD	F	441	4	-	0/30/50/50	0/6/6/6
3	LMT	F	600	-	-	0/21/61/61	0/2/2/2
5	PS9	F	800	1	-	0/0/8/8	0/1/1/1
6	MES	F	900	-	-	0/6/14/14	0/1/1/1

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	441	FAD	C10-N10	-2.78	1.35	1.39
2	F	441	FAD	C10-N10	-2.76	1.35	1.39
2	E	441	FAD	C10-N10	-2.54	1.36	1.39
2	A	441	FAD	C10-N10	-2.48	1.36	1.39
2	D	441	FAD	C10-N10	-2.47	1.36	1.39
2	B	441	FAD	C10-N10	-2.09	1.36	1.39
2	A	441	FAD	O4B-C1B	2.20	1.44	1.41
2	E	441	FAD	O4B-C1B	2.38	1.44	1.41
2	D	441	FAD	O4B-C1B	2.56	1.44	1.41
3	B	600	LMT	O1'-C1'	2.64	1.44	1.40
2	B	441	FAD	O4-C4	4.35	1.35	1.24
2	F	441	FAD	O4-C4	4.58	1.35	1.24
2	D	441	FAD	O4-C4	4.64	1.35	1.24
2	A	441	FAD	O4-C4	4.70	1.36	1.24
2	C	441	FAD	O4-C4	4.82	1.36	1.24
2	E	441	FAD	O4-C4	4.86	1.36	1.24

All (93) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	441	FAD	N3A-C2A-N1A	-11.18	120.33	128.89
2	F	441	FAD	N3A-C2A-N1A	-10.47	120.88	128.89
2	A	441	FAD	N3A-C2A-N1A	-10.39	120.94	128.89
2	C	441	FAD	N3A-C2A-N1A	-10.31	121.00	128.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	441	FAD	N3A-C2A-N1A	-10.17	121.11	128.89
2	B	441	FAD	N3A-C2A-N1A	-9.94	121.28	128.89
6	E	900	MES	C6-C5-N4	-3.34	105.07	110.12
6	E	900	MES	C2-C3-N4	-3.21	105.27	110.12
6	F	900	MES	C6-C5-N4	-2.84	105.82	110.12
2	B	441	FAD	C4A-C5A-N7A	-2.75	106.95	109.48
2	B	441	FAD	P-O3P-PA	-2.70	125.14	132.73
2	C	441	FAD	C4X-C4-N3	-2.68	119.93	123.59
6	B	900	MES	C2-C3-N4	-2.67	106.07	110.12
2	E	441	FAD	C4X-C4-N3	-2.61	120.02	123.59
6	A	900	MES	C6-C5-N4	-2.60	106.18	110.12
6	F	900	MES	C2-C3-N4	-2.60	106.18	110.12
2	D	441	FAD	P-O3P-PA	-2.55	125.57	132.73
2	E	441	FAD	C4A-C5A-N7A	-2.51	107.17	109.48
3	A	600	LMT	C1B-O1B-C4'	-2.50	111.48	118.01
6	A	900	MES	C2-C3-N4	-2.49	106.36	110.12
2	D	441	FAD	C4A-C5A-N7A	-2.43	107.24	109.48
2	B	441	FAD	C1B-N9A-C4A	-2.42	123.28	126.94
2	A	441	FAD	C4X-C4-N3	-2.42	120.28	123.59
6	B	900	MES	C6-C5-N4	-2.36	106.55	110.12
2	E	441	FAD	C1B-N9A-C4A	-2.34	123.41	126.94
2	D	441	FAD	C4X-C4-N3	-2.30	120.45	123.59
2	A	441	FAD	C4A-C5A-N7A	-2.17	107.48	109.48
2	C	441	FAD	C4A-C5A-N7A	-2.17	107.49	109.48
2	F	441	FAD	C4X-C4-N3	-2.15	120.65	123.59
2	C	441	FAD	C1B-N9A-C4A	-2.10	123.78	126.94
2	F	441	FAD	C1B-N9A-C4A	-2.09	123.79	126.94
3	C	600	LMT	C3'-C4'-C5'	-2.03	106.24	110.84
2	F	441	FAD	C4A-C5A-N7A	-2.00	107.64	109.48
3	E	600	LMT	C1B-O5B-C5B	2.19	117.99	113.75
3	B	600	LMT	O5B-C5B-C6B	2.21	111.95	106.36
2	B	441	FAD	C4X-N5-C5X	2.37	119.49	116.76
6	D	900	MES	C7-N4-C5	2.43	117.50	111.27
2	E	441	FAD	C2B-C1B-N9A	2.47	118.07	114.29
6	C	900	MES	C7-N4-C5	2.54	117.77	111.27
2	D	441	FAD	C4X-N5-C5X	2.55	119.69	116.76
3	C	600	LMT	O1B-C4'-C5'	2.56	116.06	109.32
3	E	600	LMT	O1B-C4'-C5'	2.59	116.13	109.32
3	D	600	LMT	C1B-O5B-C5B	2.63	118.85	113.75
3	A	600	LMT	C1B-C2B-C3B	2.65	115.20	109.97
2	D	441	FAD	C1'-N10-C9A	2.66	121.84	118.86
6	A	900	MES	O1S-S-C8	2.70	109.20	106.91

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	441	FAD	C4X-N5-C5X	2.72	119.89	116.76
2	A	441	FAD	C4X-N5-C5X	2.74	119.92	116.76
2	F	441	FAD	C4X-N5-C5X	2.98	120.19	116.76
3	C	600	LMT	C1B-O5B-C5B	2.98	119.53	113.75
3	B	600	LMT	C1-O1'-C1'	2.98	119.16	113.94
2	E	441	FAD	C4X-N5-C5X	2.99	120.20	116.76
3	E	600	LMT	O1B-C4'-C3'	3.03	115.00	107.17
2	C	441	FAD	C1'-N10-C9A	3.06	122.30	118.86
6	C	900	MES	C7-N4-C3	3.14	119.31	111.27
6	F	900	MES	C7-N4-C5	3.15	119.34	111.27
6	A	900	MES	C7-N4-C5	3.16	119.38	111.27
6	D	900	MES	C7-N4-C3	3.24	119.58	111.27
6	B	900	MES	C7-N4-C3	3.30	119.72	111.27
6	E	900	MES	C7-N4-C5	3.34	119.82	111.27
3	B	600	LMT	C1B-O5B-C5B	3.38	120.30	113.75
3	C	600	LMT	O1B-C4'-C3'	3.48	116.16	107.17
6	A	900	MES	C7-N4-C3	3.49	120.22	111.27
2	C	441	FAD	C5X-C9A-N10	3.50	120.28	117.62
6	F	900	MES	C7-N4-C3	3.51	120.26	111.27
6	B	900	MES	C7-N4-C5	3.52	120.28	111.27
6	E	900	MES	C7-N4-C3	3.60	120.50	111.27
2	D	441	FAD	C5X-C9A-N10	3.63	120.38	117.62
2	F	441	FAD	C5X-C9A-N10	3.65	120.39	117.62
6	F	900	MES	O2S-S-C8	3.72	110.08	106.91
2	A	441	FAD	C5X-C9A-N10	3.73	120.45	117.62
6	E	900	MES	O1S-S-C8	3.74	110.09	106.91
2	E	441	FAD	C5X-C9A-N10	3.89	120.58	117.62
3	F	600	LMT	O1B-C1B-C2B	3.90	117.59	108.10
2	B	441	FAD	C5X-C9A-N10	3.93	120.60	117.62
6	C	900	MES	O2S-S-C8	4.19	110.48	106.91
6	A	900	MES	O2S-S-C8	4.38	110.64	106.91
2	A	441	FAD	C4-N3-C2	4.58	119.20	115.25
2	B	441	FAD	C4-N3-C2	4.69	119.30	115.25
6	C	900	MES	O1S-S-C8	4.70	110.91	106.91
6	F	900	MES	O1S-S-C8	4.88	111.07	106.91
6	E	900	MES	C5-N4-C3	4.97	119.67	108.90
2	F	441	FAD	C4-N3-C2	5.01	119.58	115.25
6	B	900	MES	C5-N4-C3	5.13	120.00	108.90
6	F	900	MES	C5-N4-C3	5.31	120.40	108.90
6	A	900	MES	C5-N4-C3	5.31	120.41	108.90
6	B	900	MES	O2S-S-C8	5.45	111.55	106.91
2	C	441	FAD	C4-N3-C2	5.65	120.13	115.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	441	FAD	C4-N3-C2	5.69	120.17	115.25
2	E	441	FAD	C4-N3-C2	6.04	120.47	115.25
6	C	900	MES	C5-N4-C3	6.47	122.92	108.90
6	D	900	MES	C5-N4-C3	6.47	122.92	108.90
3	B	600	LMT	O1'-C1'-C2'	7.17	117.09	108.04

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

17 monomers are involved in 29 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	441	FAD	1	0
5	A	800	PS9	2	0
7	B	435	SO4	1	0
5	B	800	PS9	2	0
7	C	432	SO4	2	0
3	C	600	LMT	2	0
5	C	800	PS9	1	0
6	C	900	MES	1	0
7	D	432	SO4	1	0
2	D	441	FAD	1	0
5	D	802	PS9	1	0
2	E	441	FAD	1	0
3	E	600	LMT	8	0
7	F	431	SO4	1	0
2	F	441	FAD	1	0
3	F	600	LMT	9	0
5	F	800	PS9	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	428/430 (99%)	0.19	12 (2%)	56	66	31, 38, 45, 51	0
1	B	428/430 (99%)	0.18	12 (2%)	56	66	31, 38, 45, 54	0
1	C	428/430 (99%)	0.16	11 (2%)	59	68	30, 38, 44, 49	0
1	D	428/430 (99%)	0.42	43 (10%)	9	14	30, 38, 44, 47	0
1	E	428/430 (99%)	0.16	16 (3%)	45	54	31, 38, 45, 47	0
1	F	428/430 (99%)	0.15	12 (2%)	56	66	32, 38, 45, 54	0
All	All	2568/2580 (99%)	0.21	106 (4%)	41	50	30, 38, 44, 54	0

All (106) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	430	CYS	6.8
1	A	2	ALA	6.7
1	C	430	CYS	6.5
1	A	430	CYS	6.4
1	B	430	CYS	5.1
1	E	430	CYS	4.6
1	B	237	ASN	4.1
1	D	2	ALA	4.1
1	D	430	CYS	4.0
1	D	237	ASN	3.9
1	F	2	ALA	3.7
1	F	237	ASN	3.6
1	D	236	LEU	3.6
1	E	2	ALA	3.6
1	A	423	GLU	3.5
1	E	92	SER	3.4
1	D	96	ILE	3.4
1	D	89	THR	3.3
1	D	97	GLU	3.3

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Mol	Chain	Res	Type	RSRZ
1	D	92	SER	3.3
1	D	265	ASP	3.3
1	B	91	GLN	3.2
1	B	420	LYS	3.2
1	C	2	ALA	3.2
1	D	239	ASN	3.2
1	C	27	ASP	3.1
1	E	141	ALA	3.1
1	D	29	LYS	3.1
1	F	303	LYS	3.1
1	D	119	ASN	3.0
1	E	91	GLN	3.0
1	D	423	GLU	3.0
1	D	303	LYS	3.0
1	D	93	GLY	3.0
1	D	271	ALA	2.9
1	D	118	GLU	2.9
1	D	91	GLN	2.9
1	D	228	PRO	2.9
1	D	79	GLU	2.9
1	F	271	ALA	2.8
1	D	27	ASP	2.8
1	C	303	LYS	2.8
1	B	236	LEU	2.8
1	D	95	LYS	2.8
1	D	261	ALA	2.8
1	B	167	LEU	2.8
1	D	90	THR	2.8
1	B	239	ASN	2.7
1	D	117	GLU	2.7
1	D	141	ALA	2.7
1	E	93	GLY	2.7
1	F	423	GLU	2.7
1	F	56	GLU	2.6
1	D	334	ASN	2.6
1	A	237	ASN	2.6
1	D	421	ASP	2.6
1	F	81	ILE	2.6
1	E	27	ASP	2.5
1	D	266	LYS	2.5
1	E	408	GLU	2.5
1	E	142	ASN	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	27	ASP	2.5
1	D	227	GLU	2.5
1	F	420	LYS	2.5
1	D	142	ASN	2.5
1	C	164	ALA	2.4
1	D	240	THR	2.4
1	A	8	ILE	2.4
1	D	138	GLU	2.4
1	D	74	ILE	2.4
1	B	2	ALA	2.4
1	B	423	GLU	2.4
1	E	265	ASP	2.4
1	A	250	PHE	2.4
1	C	142	ASN	2.3
1	A	91	GLN	2.3
1	E	336	PRO	2.3
1	C	91	GLN	2.3
1	F	27	ASP	2.3
1	E	137	GLN	2.3
1	B	303	LYS	2.3
1	F	97	GLU	2.2
1	D	263	ALA	2.2
1	E	303	LYS	2.2
1	D	270	PRO	2.2
1	D	82	ASP	2.2
1	A	7	VAL	2.2
1	E	94	LYS	2.2
1	B	164	ALA	2.2
1	D	35	ASP	2.1
1	D	4	HIS	2.1
1	B	27	ASP	2.1
1	F	142	ASN	2.1
1	A	420	LYS	2.1
1	D	336	PRO	2.1
1	E	95	LYS	2.1
1	A	167	LEU	2.1
1	C	185	ILE	2.1
1	D	241	HIS	2.1
1	C	149	GLY	2.1
1	A	164	ALA	2.1
1	D	242	GLU	2.0
1	D	285	THR	2.0

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Mol	Chain	Res	Type	RSRZ
1	E	89	THR	2.0
1	C	343	LEU	2.0
1	C	423	GLU	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	CSS	D	156[B]	7/8	0.97	0.09	-	34,35,35,35	2
1	CSS	C	156[B]	7/8	0.93	0.12	-	38,39,40,40	2
1	CSS	C	156[A]	6/8	0.93	0.12	-	39,39,40,43	1
1	CSS	A	156[A]	6/8	0.95	0.18	-	37,38,39,42	1
1	CSS	E	156[A]	6/8	0.95	0.14	-	37,38,38,40	1
1	CSS	B	156[B]	7/8	0.93	0.14	-	35,36,36,36	2
1	CSS	A	156[B]	7/8	0.95	0.18	-	35,38,38,39	2
1	CSS	F	156[B]	7/8	0.95	0.15	-	35,37,37,37	2
1	CSS	E	156[B]	7/8	0.95	0.14	-	37,38,38,38	2
1	CSS	B	156[A]	6/8	0.93	0.14	-	36,36,36,36	1
1	CSS	F	156[A]	6/8	0.95	0.15	-	37,37,37,38	1
1	CSS	D	156[A]	6/8	0.97	0.09	-	35,35,35,37	1

## 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, 95<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
5	PS9	D	802	5/8	0.48	0.96	32.40	48,48,48,49	5
4	H2S	F	700	1/1	0.95	0.29	29.53	30,30,30,30	1
5	PS9	C	800	6/8	0.47	0.43	24.50	37,38,38,39	6
5	PS9	E	800	8/8	0.62	0.46	17.12	39,39,40,40	8
5	PS9	B	800	6/8	0.66	0.51	16.92	36,37,37,37	6
5	PS9	F	800	8/8	0.61	0.47	15.76	35,35,36,37	8
4	H2S	A	700	1/1	0.97	0.35	11.60	31,31,31,31	1
5	PS9	A	800	8/8	0.66	0.45	11.39	34,35,35,35	8
3	LMT	F	600	35/35	0.41	0.45	9.88	36,45,52,53	35
4	H2S	E	700	1/1	0.95	0.26	9.73	32,32,32,32	1
4	H2S	C	700	1/1	0.98	0.33	9.64	34,34,34,34	1
3	LMT	C	600	35/35	0.55	0.40	8.96	34,46,51,52	35
3	LMT	A	600	35/35	0.30	0.48	8.69	40,44,49,50	35
4	H2S	B	700	1/1	0.94	0.21	7.29	32,32,32,32	1
3	LMT	D	600	35/35	0.58	0.37	6.78	43,49,50,51	35
3	LMT	B	600	35/35	0.50	0.38	6.57	41,47,51,52	35
3	LMT	E	600	35/35	0.61	0.40	6.54	40,43,47,47	35
4	H2S	D	700	1/1	0.93	0.16	2.28	37,37,37,37	1
6	MES	D	900	12/12	0.91	0.22	1.58	48,53,57,58	0
6	MES	A	900	12/12	0.92	0.20	0.78	56,60,63,63	0
6	MES	C	900	12/12	0.92	0.19	0.63	50,57,61,62	0
7	SO4	B	433	5/5	0.86	0.16	0.06	65,66,66,67	0
6	MES	E	900	12/12	0.96	0.14	-0.42	55,59,63,63	0
6	MES	F	900	12/12	0.96	0.15	-0.50	47,51,55,55	0
2	FAD	D	441	53/53	0.95	0.11	-0.66	31,36,38,40	0
7	SO4	F	433	5/5	0.93	0.15	-0.67	62,63,63,64	0
2	FAD	E	441	53/53	0.95	0.10	-0.95	26,29,42,43	0
7	SO4	C	433	5/5	0.94	0.11	-1.19	60,61,61,62	0
7	SO4	E	433	5/5	0.95	0.09	-1.25	72,72,73,73	0
2	FAD	B	441	53/53	0.97	0.09	-1.61	23,30,30,32	0
2	FAD	A	441	53/53	0.96	0.10	-1.61	23,28,30,32	0
7	SO4	A	433	5/5	0.98	0.07	-1.71	42,42,43,43	0
6	MES	B	900	12/12	0.97	0.10	-1.74	47,52,55,55	0
2	FAD	F	441	53/53	0.97	0.09	-1.76	25,30,34,35	0
7	SO4	D	433	5/5	0.91	0.12	-1.77	101,101,101,101	0
2	FAD	C	441	53/53	0.97	0.08	-1.78	22,30,32,32	0
7	SO4	B	435	5/5	0.95	0.15	-	70,70,70,71	0
7	SO4	C	434	5/5	0.92	0.21	-	62,62,63,63	0
7	SO4	E	431	5/5	0.86	0.22	-	69,69,70,70	0
7	SO4	F	431	5/5	0.93	0.16	-	72,72,73,73	0
7	SO4	B	431	5/5	0.96	0.12	-	77,77,78,78	0
7	SO4	C	431	5/5	0.95	0.20	-	71,71,71,71	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
7	SO4	E	435	5/5	0.94	0.14	-	75,75,76,76	0
7	SO4	B	434	5/5	0.95	0.33	-	64,64,64,65	0
7	SO4	B	432	5/5	0.96	0.17	-	54,54,55,55	0
7	SO4	A	434	5/5	0.96	0.19	-	61,62,62,63	0
5	PS9	D	800	1/8	0.71	0.51	-	33,33,33,33	1
7	SO4	D	434	5/5	0.90	0.28	-	69,69,70,70	0
7	SO4	E	432	5/5	0.94	0.32	-	66,66,67,67	0
7	SO4	F	434	5/5	0.95	0.26	-	63,64,64,65	0
7	SO4	A	432	5/5	0.90	0.21	-	61,61,62,62	0
7	SO4	E	434	5/5	0.91	0.20	-	68,68,68,69	0
7	SO4	A	431	5/5	0.89	0.24	-	75,76,76,76	0
7	SO4	C	432	5/5	0.93	0.16	-	66,66,66,66	0
7	SO4	D	432	5/5	0.88	0.17	-	71,72,72,72	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.