wwPDB X-ray Structure Validation Summary Report (i)

Feb 1, 2016 – 11:50 AM GMT

PDB ID	:	3Q09
Title	:	Crystal Structure of Chlorite Dismutase from D. Aromatica at pH 9.0
Authors	:	Goblirsch, B.R.; Wilmot, C.M.
Deposited on	:	2010-12-15
Resolution	:	3.00 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

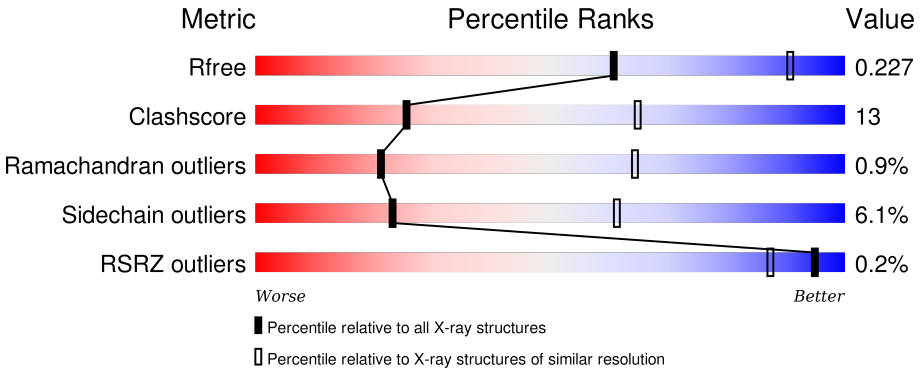
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	1578 (3.00-3.00)
Clashscore	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)
RSRZ outliers	91569	1592 (3.00-3.00)

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 <=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	248	70% 25%
1	B	248	72% 22%
1	C	248	% 73% 20%
1	D	248	73% 22%
1	E	248	71% 23%

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Mol	Chain	Length	Quality of chain
1	F	248	71% 25%
1	G	248	70% 24%
1	H	248	76% 19%
1	I	248	75% 19%
1	J	248	73% 20%
1	K	248	75% 21%
1	L	248	74% 20%
1	M	248	74% 21%
1	N	248	70% 24%
1	O	248	76% 19%
1	P	248	70% 24%
1	Q	248	67% 25% 5%
1	R	248	70% 24%
1	S	248	2% 67% 28%
1	T	248	62% 31% 5%

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	HEM	G	1000	-	-	-	X
3	NO2	A	2000	-	-	-	X
3	NO2	B	2000	-	-	-	X
3	NO2	C	2000	-	-	-	X
3	NO2	I	2000	-	-	-	X
3	NO2	K	2000	-	-	-	X
3	NO2	N	2000	-	-	-	X
3	NO2	R	2000	-	-	-	X

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 39520 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 Chlorite dismutase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	241	Total	C	N	O	S	0	0	0
			1928	1239	325	359	5			
1	B	241	Total	C	N	O	S	0	0	0
			1928	1239	325	359	5			
1	C	241	Total	C	N	O	S	0	0	0
			1928	1239	325	359	5			
1	D	241	Total	C	N	O	S	0	0	0
			1928	1239	325	359	5			
1	E	241	Total	C	N	O	S	0	0	0
			1928	1239	325	359	5			
1	F	241	Total	C	N	O	S	0	0	0
			1928	1239	325	359	5			
1	G	241	Total	C	N	O	S	0	0	0
			1928	1239	325	359	5			
1	H	241	Total	C	N	O	S	0	0	0
			1928	1239	325	359	5			
1	I	241	Total	C	N	O	S	0	0	0
			1928	1239	325	359	5			
1	J	241	Total	C	N	O	S	0	0	0
			1928	1239	325	359	5			
1	K	241	Total	C	N	O	S	0	0	0
			1928	1239	325	359	5			
1	L	241	Total	C	N	O	S	0	0	0
			1928	1239	325	359	5			
1	M	241	Total	C	N	O	S	0	0	0
			1928	1239	325	359	5			
1	N	241	Total	C	N	O	S	0	0	0
			1928	1239	325	359	5			
1	O	241	Total	C	N	O	S	0	0	0
			1928	1239	325	359	5			
1	P	241	Total	C	N	O	S	0	0	0
			1928	1239	325	359	5			

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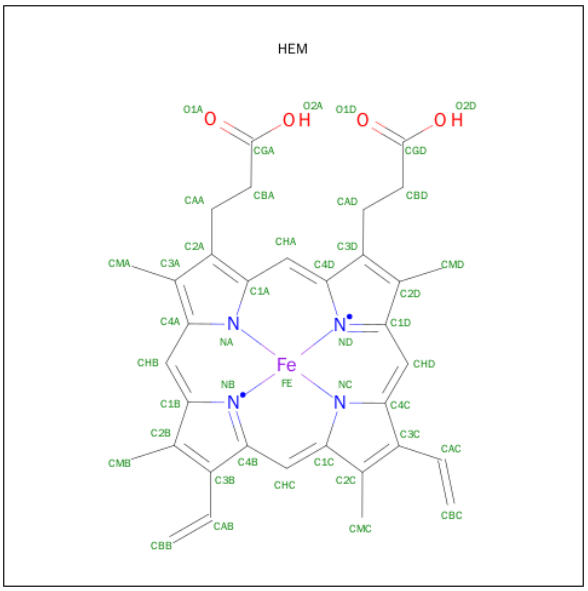
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Q	241	Total	C	N	O	S	0	0	0
			1928	1239	325	359	5			
1	R	241	Total	C	N	O	S	0	0	0
			1928	1239	325	359	5			
1	S	241	Total	C	N	O	S	0	0	0
			1928	1239	325	359	5			
1	T	241	Total	C	N	O	S	0	0	0
			1928	1239	325	359	5			

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	C	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	D	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	E	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	F	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	G	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

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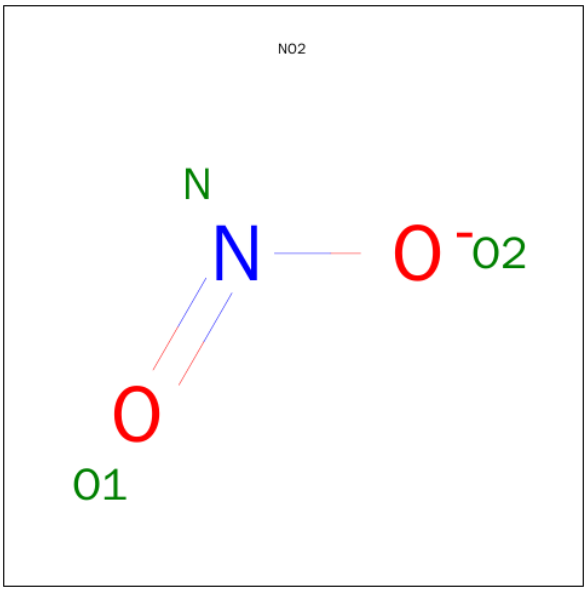
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	H	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	I	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	J	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	K	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	L	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	M	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	N	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	O	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	P	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	Q	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	R	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	S	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	T	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 3 is NITRITE ION (three-letter code: NO2) (formula: NO_2).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total N O 3 1 2	0	0
3	B	1	Total N O 3 1 2	0	0
3	C	1	Total N O 3 1 2	0	0
3	D	1	Total N O 3 1 2	0	0
3	E	1	Total N O 3 1 2	0	0
3	F	1	Total N O 3 1 2	0	0
3	G	1	Total N O 3 1 2	0	0
3	H	1	Total N O 3 1 2	0	0
3	I	1	Total N O 3 1 2	0	0
3	J	1	Total N O 3 1 2	0	0
3	K	1	Total N O 3 1 2	0	0
3	L	1	Total N O 3 1 2	0	0
3	M	1	Total N O 3 1 2	0	0
3	N	1	Total N O 3 1 2	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	O	1	Total N O 3 1 2	0	0
3	P	1	Total N O 3 1 2	0	0
3	Q	1	Total N O 3 1 2	0	0
3	R	1	Total N O 3 1 2	0	0
3	S	1	Total N O 3 1 2	0	0
3	T	1	Total N O 3 1 2	0	0

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	P	1	Total Ca 1 1	0	0
4	G	1	Total Ca 1 1	0	0
4	J	1	Total Ca 1 1	0	0
4	Q	1	Total Ca 1 1	0	0
4	D	1	Total Ca 1 1	0	0
4	K	1	Total Ca 1 1	0	0
4	E	1	Total Ca 1 1	0	0
4	H	1	Total Ca 1 1	0	0
4	B	1	Total Ca 1 1	0	0
4	I	1	Total Ca 1 1	0	0
4	C	1	Total Ca 1 1	0	0
4	A	1	Total Ca 1 1	0	0
4	T	1	Total Ca 1 1	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	N	1	Total Ca 1 1	0	0
4	O	1	Total Ca 1 1	0	0
4	R	1	Total Ca 1 1	0	0
4	L	1	Total Ca 1 1	0	0
4	S	1	Total Ca 1 1	0	0
4	F	1	Total Ca 1 1	0	0
4	M	1	Total Ca 1 1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total O 1 1	0	0
5	B	1	Total O 1 1	0	0
5	C	1	Total O 1 1	0	0
5	D	1	Total O 1 1	0	0
5	E	1	Total O 1 1	0	0
5	F	1	Total O 1 1	0	0
5	G	1	Total O 1 1	0	0
5	H	1	Total O 1 1	0	0
5	I	1	Total O 1 1	0	0
5	J	1	Total O 1 1	0	0
5	K	1	Total O 1 1	0	0
5	L	1	Total O 1 1	0	0

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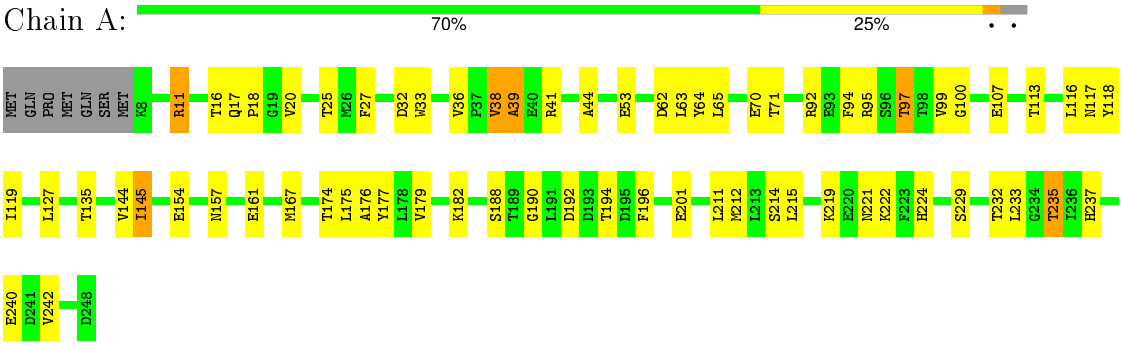
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	M	1	Total O 1 1	0	0
5	N	1	Total O 1 1	0	0
5	O	1	Total O 1 1	0	0
5	P	1	Total O 1 1	0	0
5	Q	1	Total O 1 1	0	0
5	R	1	Total O 1 1	0	0
5	S	1	Total O 1 1	0	0
5	T	1	Total O 1 1	0	0

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.

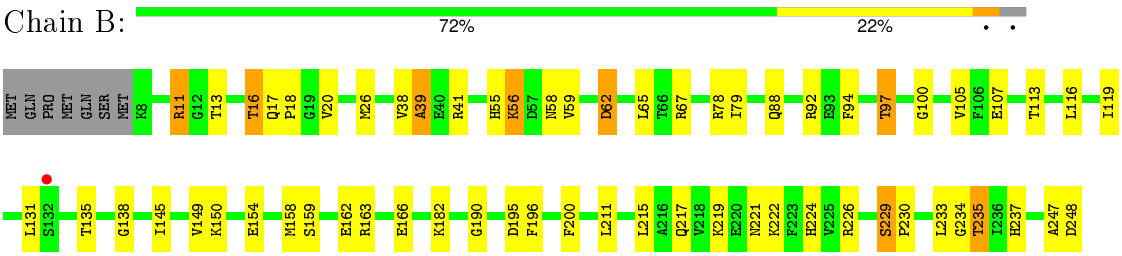
- Molecule 1: Chlorite dismutase

Chain A:



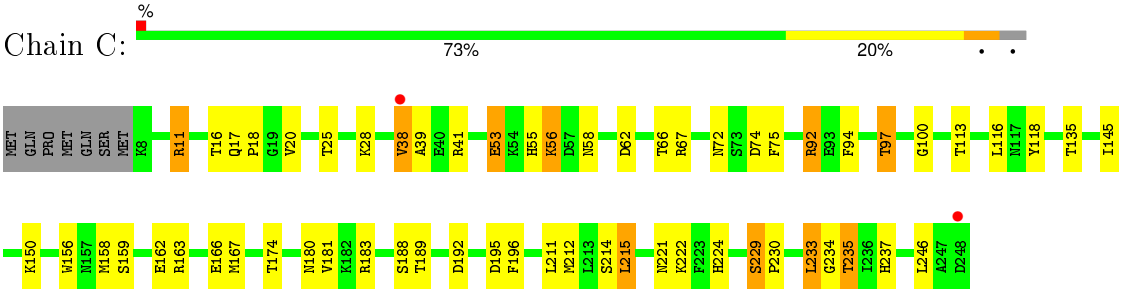
- Molecule 1: Chlorite dismutase

Chain B:



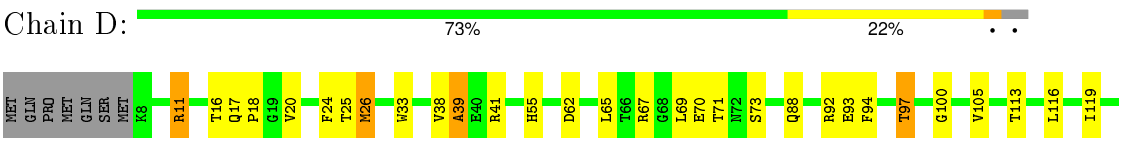
- Molecule 1: Chlorite dismutase

Chain C:



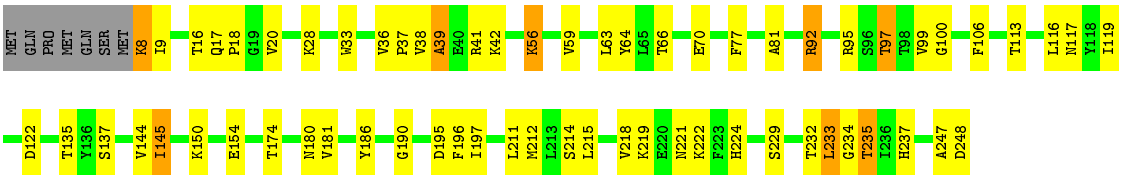
- Molecule 1: Chlorite dismutase

Chain D:



- Molecule 1: Chlorite dismutase

Chain E:



- Molecule 1: Chlorite dismutase

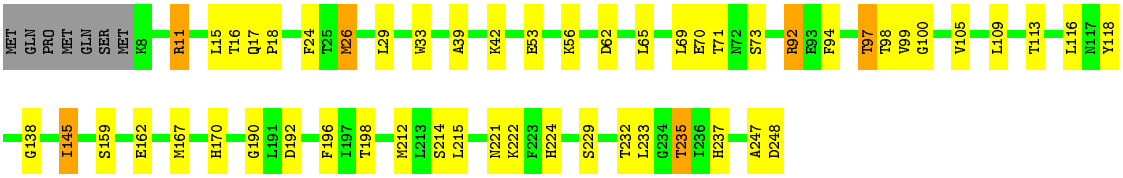
Chain F:

- Molecule 1: Chlorite dismutase

Chain G:

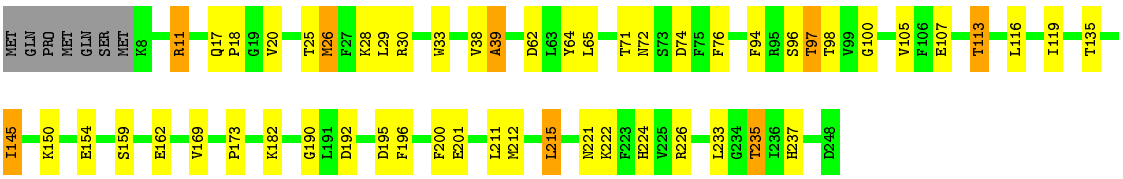
- Molecule 1: Chlorite dismutase

Chain H:



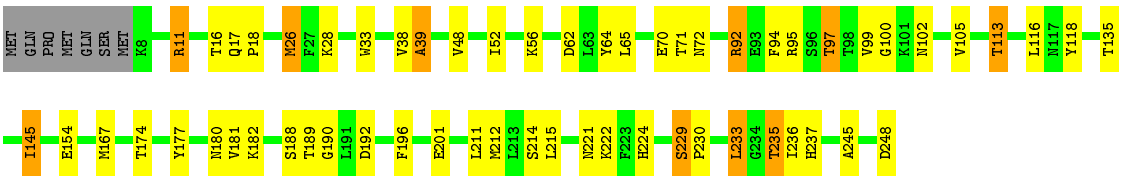
- Molecule 1: Chlorite dismutase

Chain I:



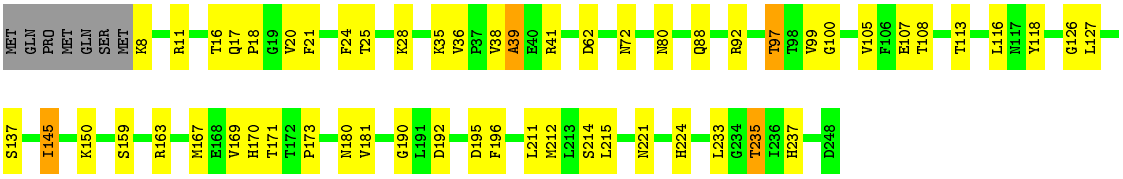
• Molecule 1: Chlorite dismutase

Chain J:



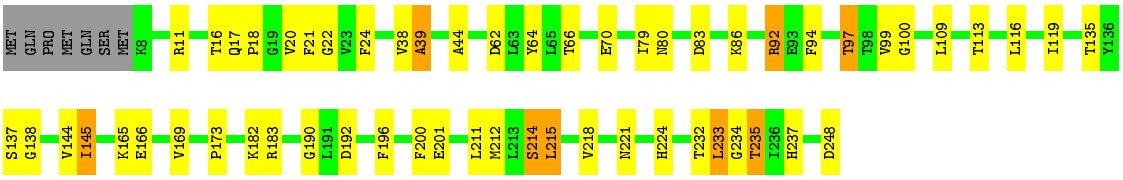
• Molecule 1: Chlorite dismutase

Chain K:



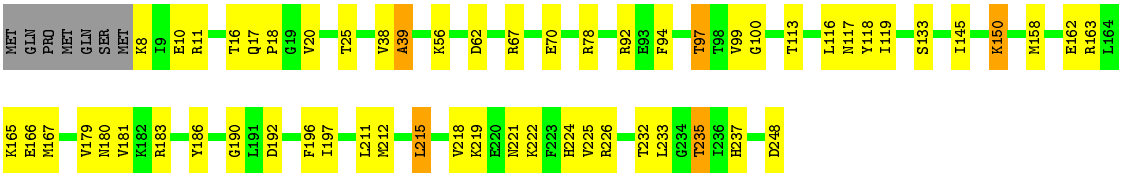
• Molecule 1: Chlorite dismutase

Chain L:



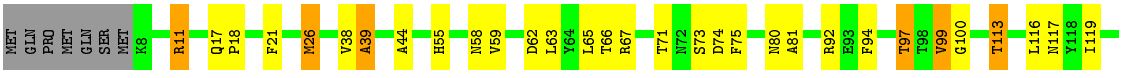
• Molecule 1: Chlorite dismutase

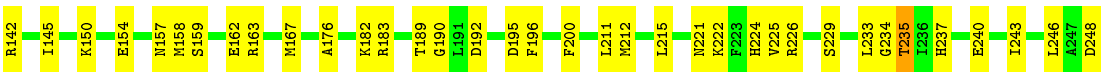
Chain M:



• Molecule 1: Chlorite dismutase

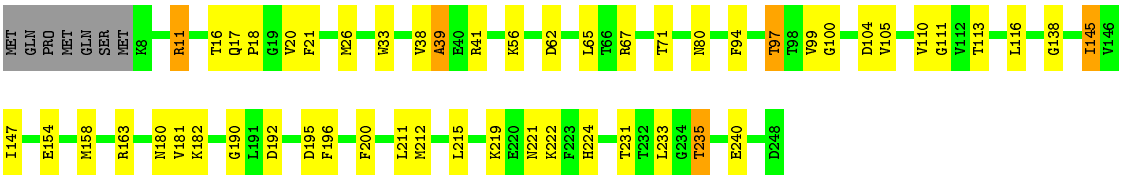
Chain N:





• Molecule 1: Chlorite dismutase

Chain O:

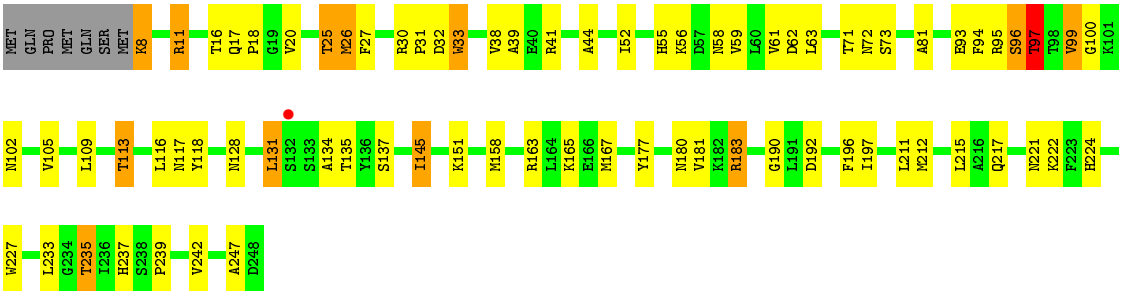


• Molecule 1: Chlorite dismutase

Chain P:

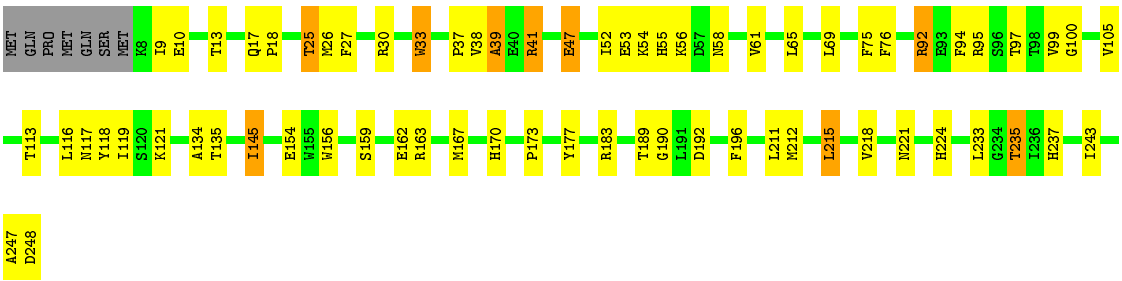
• Molecule 1: Chlorite dismutase

Chain Q:



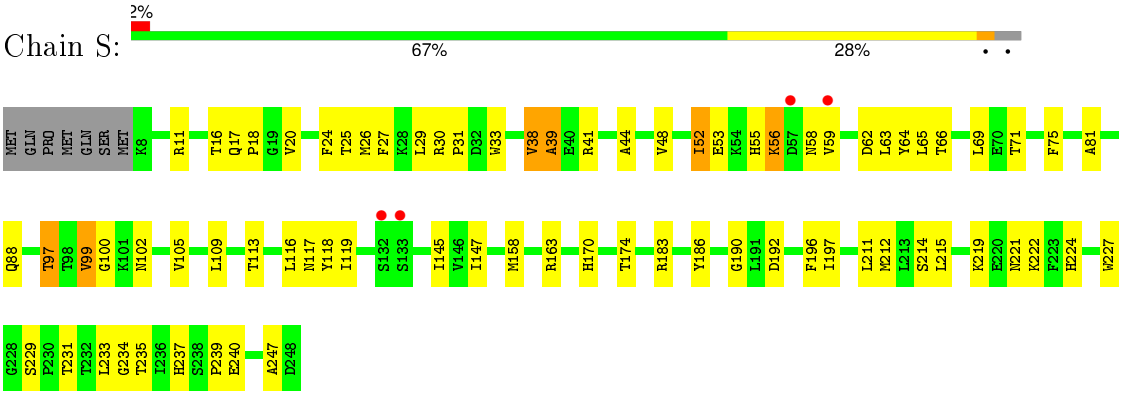
• Molecule 1: Chlorite dismutase

Chain R:



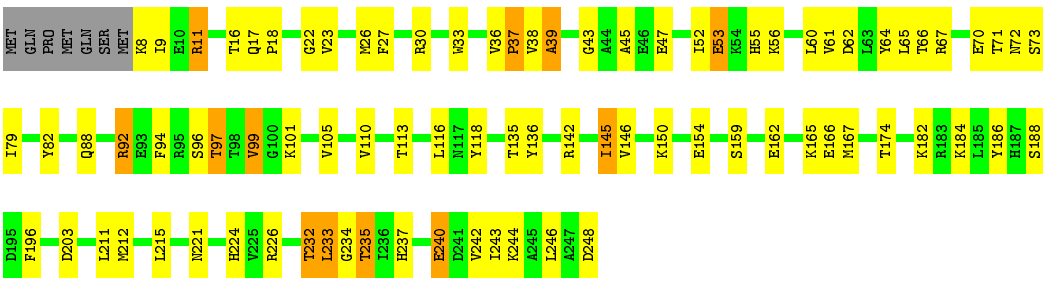
• Molecule 1: Chlorite dismutase

Chain S:



• Molecule 1: Chlorite dismutase

Chain T:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	122.70Å 202.85Å 247.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.31 – 3.00 48.31 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.7 (48.31-3.00) 99.7 (48.31-3.00)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	5.59 (at 3.01Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.185 , 0.234 0.183 , 0.227	Depositor DCC
R_{free} test set	6174 reflections (5.28%)	DCC
Wilson B-factor (Å^2)	39.8	Xtriage
Anisotropy	0.038	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.28 , 17.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.34$	Xtriage
Outliers	2 of 123079 reflections (0.002%)	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	39520	wwPDB-VP
Average B, all atoms (Å^2)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.33% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: HEM, CA, NO2

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	0.83	2/1973 (0.1%)	0.86	2/2676 (0.1%)
1	B	0.87	0/1973	0.89	2/2676 (0.1%)
1	C	0.85	1/1973 (0.1%)	0.85	1/2676 (0.0%)
1	D	0.80	0/1973	0.83	0/2676
1	E	0.84	0/1973	0.86	2/2676 (0.1%)
1	F	0.75	1/1973 (0.1%)	0.81	0/2676
1	G	0.77	0/1973	0.81	0/2676
1	H	0.76	0/1973	0.83	0/2676
1	I	0.80	0/1973	0.84	0/2676
1	J	0.75	0/1973	0.81	0/2676
1	K	0.76	0/1973	0.83	0/2676
1	L	0.76	0/1973	0.79	0/2676
1	M	0.80	1/1973 (0.1%)	0.85	1/2676 (0.0%)
1	N	0.75	1/1973 (0.1%)	0.79	0/2676
1	O	0.71	0/1973	0.77	1/2676 (0.0%)
1	P	0.82	6/1973 (0.3%)	0.83	5/2676 (0.2%)
1	Q	0.99	6/1973 (0.3%)	0.88	6/2676 (0.2%)
1	R	0.78	4/1973 (0.2%)	0.77	1/2676 (0.0%)
1	S	0.76	1/1973 (0.1%)	0.73	0/2676
1	T	0.88	7/1973 (0.4%)	0.83	1/2676 (0.0%)
All	All	0.80	30/39460 (0.1%)	0.82	22/53520 (0.0%)

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	H	0	1
1	L	0	1
1	M	0	2

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Mol	Chain	#Chirality outliers	#Planarity outliers
1	O	0	1
All	All	0	5

The worst 5 of 30 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	Q	93	GLU	CD-OE1	16.32	1.43	1.25
1	Q	33	TRP	CG-CD1	11.41	1.52	1.36
1	Q	33	TRP	CB-CG	10.37	1.69	1.50
1	R	33	TRP	CB-CG	8.30	1.65	1.50
1	Q	96	SER	CB-OG	7.63	1.52	1.42

The worst 5 of 22 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	Q	33	TRP	CD1-CG-CD2	-8.99	99.11	106.30
1	P	75	PHE	CB-CG-CD1	8.80	126.96	120.80
1	Q	33	TRP	CG-CD2-CE3	-8.24	126.48	133.90
1	Q	33	TRP	CB-CG-CD2	8.17	137.22	126.60
1	Q	131	LEU	CA-CB-CG	7.86	133.39	115.30

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	H	138	GLY	Peptide
1	L	138	GLY	Peptide
1	M	133	SER	Peptide
1	M	179	VAL	Peptide
1	O	138	GLY	Peptide

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1928	0	1938	57	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1928	0	1938	51	0
1	C	1928	0	1938	50	0
1	D	1928	0	1938	57	0
1	E	1928	0	1938	54	0
1	F	1928	0	1938	54	0
1	G	1928	0	1938	65	0
1	H	1928	0	1938	43	0
1	I	1928	0	1938	49	0
1	J	1928	0	1938	56	0
1	K	1928	0	1938	41	0
1	L	1928	0	1938	47	0
1	M	1928	0	1938	41	0
1	N	1928	0	1938	59	0
1	O	1928	0	1938	41	0
1	P	1928	0	1938	58	0
1	Q	1928	0	1938	60	0
1	R	1928	0	1938	60	0
1	S	1928	0	1938	62	0
1	T	1928	0	1938	61	0
2	A	43	0	30	8	0
2	B	43	0	30	6	0
2	C	43	0	30	7	0
2	D	43	0	30	9	0
2	E	43	0	30	8	0
2	F	43	0	30	11	0
2	G	43	0	30	15	0
2	H	43	0	30	9	0
2	I	43	0	30	7	0
2	J	43	0	30	10	0
2	K	43	0	30	5	0
2	L	43	0	30	7	0
2	M	43	0	30	6	0
2	N	43	0	30	10	0
2	O	43	0	30	5	0
2	P	43	0	30	3	0
2	Q	43	0	30	10	0
2	R	43	0	30	14	0
2	S	43	0	30	9	0
2	T	43	0	30	6	0
3	A	3	0	0	0	0
3	B	3	0	0	0	0
3	C	3	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	3	0	0	0	0
3	E	3	0	0	0	0
3	F	3	0	0	0	0
3	G	3	0	0	1	0
3	H	3	0	0	0	0
3	I	3	0	0	0	0
3	J	3	0	0	0	0
3	K	3	0	0	0	0
3	L	3	0	0	1	0
3	M	3	0	0	0	0
3	N	3	0	0	0	0
3	O	3	0	0	0	0
3	P	3	0	0	0	0
3	Q	3	0	0	0	0
3	R	3	0	0	0	0
3	S	3	0	0	0	0
3	T	3	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	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	0	0
4	G	1	0	0	0	0
4	H	1	0	0	0	0
4	I	1	0	0	0	0
4	J	1	0	0	0	0
4	K	1	0	0	0	0
4	L	1	0	0	0	0
4	M	1	0	0	0	0
4	N	1	0	0	0	0
4	O	1	0	0	0	0
4	P	1	0	0	0	0
4	Q	1	0	0	0	0
4	R	1	0	0	0	0
4	S	1	0	0	0	0
4	T	1	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	2	0
5	C	1	0	0	1	0
5	D	1	0	0	0	0
5	E	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	F	1	0	0	1	0
5	G	1	0	0	0	0
5	H	1	0	0	2	0
5	I	1	0	0	0	0
5	J	1	0	0	0	0
5	K	1	0	0	0	0
5	L	1	0	0	2	0
5	M	1	0	0	0	0
5	N	1	0	0	0	0
5	O	1	0	0	0	0
5	P	1	0	0	0	0
5	Q	1	0	0	0	0
5	R	1	0	0	0	0
5	S	1	0	0	0	0
5	T	1	0	0	0	0
All	All	39520	0	39360	988	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 13.

The worst 5 of 988 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:O:97:THR:HG22	1:O:100:GLY:H	1.10	1.15
1:C:18:PRO:HA	1:C:113:THR:HG22	1.23	1.15
1:J:18:PRO:HA	1:J:113:THR:HG22	1.26	1.12
1:K:18:PRO:HA	1:K:113:THR:HG22	1.33	1.10
1:N:18:PRO:HA	1:N:113:THR:HG22	1.28	1.10

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	239/248 (96%)	230 (96%)	8 (3%)	1 (0%)	39	80
1	B	239/248 (96%)	228 (95%)	8 (3%)	3 (1%)	15	53
1	C	239/248 (96%)	231 (97%)	5 (2%)	3 (1%)	15	53
1	D	239/248 (96%)	232 (97%)	6 (2%)	1 (0%)	39	80
1	E	239/248 (96%)	228 (95%)	8 (3%)	3 (1%)	15	53
1	F	239/248 (96%)	226 (95%)	11 (5%)	2 (1%)	24	66
1	G	239/248 (96%)	228 (95%)	9 (4%)	2 (1%)	24	66
1	H	239/248 (96%)	227 (95%)	11 (5%)	1 (0%)	39	80
1	I	239/248 (96%)	226 (95%)	12 (5%)	1 (0%)	39	80
1	J	239/248 (96%)	229 (96%)	9 (4%)	1 (0%)	39	80
1	K	239/248 (96%)	228 (95%)	10 (4%)	1 (0%)	39	80
1	L	239/248 (96%)	231 (97%)	7 (3%)	1 (0%)	39	80
1	M	239/248 (96%)	228 (95%)	9 (4%)	2 (1%)	24	66
1	N	239/248 (96%)	225 (94%)	11 (5%)	3 (1%)	15	53
1	O	239/248 (96%)	227 (95%)	11 (5%)	1 (0%)	39	80
1	P	239/248 (96%)	223 (93%)	13 (5%)	3 (1%)	15	53
1	Q	239/248 (96%)	223 (93%)	12 (5%)	4 (2%)	11	46
1	R	239/248 (96%)	225 (94%)	11 (5%)	3 (1%)	15	53
1	S	239/248 (96%)	226 (95%)	9 (4%)	4 (2%)	11	46
1	T	239/248 (96%)	225 (94%)	12 (5%)	2 (1%)	24	66
All	All	4780/4960 (96%)	4546 (95%)	192 (4%)	42 (1%)	21	64

5 of 42 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	39	ALA
1	C	39	ALA
1	E	39	ALA
1	F	38	VAL
1	F	39	ALA

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	212/219 (97%)	197 (93%)	15 (7%)	18	54
1	B	212/219 (97%)	200 (94%)	12 (6%)	25	64
1	C	212/219 (97%)	200 (94%)	12 (6%)	25	64
1	D	212/219 (97%)	199 (94%)	13 (6%)	23	61
1	E	212/219 (97%)	199 (94%)	13 (6%)	23	61
1	F	212/219 (97%)	197 (93%)	15 (7%)	18	54
1	G	212/219 (97%)	200 (94%)	12 (6%)	25	64
1	H	212/219 (97%)	199 (94%)	13 (6%)	23	61
1	I	212/219 (97%)	201 (95%)	11 (5%)	29	68
1	J	212/219 (97%)	196 (92%)	16 (8%)	17	51
1	K	212/219 (97%)	200 (94%)	12 (6%)	25	64
1	L	212/219 (97%)	199 (94%)	13 (6%)	23	61
1	M	212/219 (97%)	200 (94%)	12 (6%)	25	64
1	N	212/219 (97%)	198 (93%)	14 (7%)	21	57
1	O	212/219 (97%)	201 (95%)	11 (5%)	29	68
1	P	212/219 (97%)	203 (96%)	9 (4%)	36	76
1	Q	212/219 (97%)	193 (91%)	19 (9%)	12	41
1	R	212/219 (97%)	202 (95%)	10 (5%)	32	72
1	S	212/219 (97%)	201 (95%)	11 (5%)	29	68
1	T	212/219 (97%)	195 (92%)	17 (8%)	15	47
All	All	4240/4380 (97%)	3980 (94%)	260 (6%)	23	61

5 of 260 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	J	16	THR
1	L	92	ARG
1	S	227	TRP
1	J	97	THR
1	K	8	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 100 such sidechains are listed below:

Mol	Chain	Res	Type
1	H	224	HIS
1	J	224	HIS
1	S	157	ASN
1	I	58	ASN
1	I	224	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 60 ligands modelled in this entry, 20 are monoatomic - leaving 40 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	HEM	A	1000	1,3	30,50,50	2.46	7 (23%)	24,82,82	2.35	10 (41%)
3	NO2	A	2000	2	2,2,2	1.41	0	1,1,1	0.81	0
2	HEM	B	1000	1,3	30,50,50	2.42	5 (16%)	24,82,82	3.22	14 (58%)
3	NO2	B	2000	2	2,2,2	1.59	0	1,1,1	0.97	0
2	HEM	C	1000	1,3	30,50,50	2.28	5 (16%)	24,82,82	2.59	12 (50%)
3	NO2	C	2000	2	2,2,2	1.60	0	1,1,1	0.50	0
2	HEM	D	1000	1,3	30,50,50	2.48	9 (30%)	24,82,82	2.34	12 (50%)
3	NO2	D	2000	2	2,2,2	1.53	0	1,1,1	0.49	0
2	HEM	E	1000	1,3	30,50,50	1.90	4 (13%)	24,82,82	2.89	13 (54%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
3	NO2	E	2000	2	2,2,2	1.44	0	1,1,1	0.86	0
2	HEM	F	1000	1,3	30,50,50	2.18	7 (23%)	24,82,82	2.44	12 (50%)
3	NO2	F	2000	2	2,2,2	1.57	0	1,1,1	0.30	0
2	HEM	G	1000	1,3	30,50,50	2.19	9 (30%)	24,82,82	3.07	13 (54%)
3	NO2	G	2000	2	2,2,2	1.60	0	1,1,1	0.57	0
2	HEM	H	1000	1,3	30,50,50	2.09	6 (20%)	24,82,82	2.93	11 (45%)
3	NO2	H	2000	2	2,2,2	1.53	0	1,1,1	0.76	0
2	HEM	I	1000	1,3	30,50,50	2.30	9 (30%)	24,82,82	2.61	8 (33%)
3	NO2	I	2000	2	2,2,2	1.37	0	1,1,1	0.94	0
2	HEM	J	1000	1,3	30,50,50	2.47	7 (23%)	24,82,82	2.70	11 (45%)
3	NO2	J	2000	2	2,2,2	1.51	0	1,1,1	0.58	0
2	HEM	K	1000	1,3	30,50,50	1.95	4 (13%)	24,82,82	2.72	9 (37%)
3	NO2	K	2000	2	2,2,2	1.61	0	1,1,1	0.85	0
2	HEM	L	1000	1,3	30,50,50	2.10	7 (23%)	24,82,82	2.73	11 (45%)
3	NO2	L	2000	2	2,2,2	1.57	0	1,1,1	0.50	0
2	HEM	M	1000	1,3	30,50,50	1.91	6 (20%)	24,82,82	3.20	16 (66%)
3	NO2	M	2000	2	2,2,2	1.82	1 (50%)	1,1,1	0.87	0
2	HEM	N	1000	1,3	30,50,50	1.76	4 (13%)	24,82,82	2.55	12 (50%)
3	NO2	N	2000	2	2,2,2	1.56	0	1,1,1	0.57	0
2	HEM	O	1000	1,3	30,50,50	1.98	8 (26%)	24,82,82	2.55	8 (33%)
3	NO2	O	2000	2	2,2,2	1.47	0	1,1,1	0.55	0
2	HEM	P	1000	1,3	30,50,50	1.80	5 (16%)	24,82,82	2.80	11 (45%)
3	NO2	P	2000	2	2,2,2	1.61	0	1,1,1	0.24	0
2	HEM	Q	1000	1,3	30,50,50	2.08	8 (26%)	24,82,82	3.04	9 (37%)
3	NO2	Q	2000	2	2,2,2	1.51	0	1,1,1	1.02	0
2	HEM	R	1000	1,3	30,50,50	2.07	9 (30%)	24,82,82	2.50	12 (50%)
3	NO2	R	2000	2	2,2,2	1.66	0	1,1,1	0.31	0
2	HEM	S	1000	1,3	30,50,50	2.12	9 (30%)	24,82,82	2.58	11 (45%)
3	NO2	S	2000	2	2,2,2	1.57	0	1,1,1	0.37	0
2	HEM	T	1000	1,3	30,50,50	2.15	6 (20%)	24,82,82	2.56	10 (41%)
3	NO2	T	2000	2	2,2,2	1.49	0	1,1,1	0.66	0

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	HEM	A	1000	1,3	-	0/10/54/54	0/0/8/8

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NO2	A	2000	2	-	0/0/0/0	0/0/0/0
2	HEM	B	1000	1,3	-	0/10/54/54	0/0/8/8
3	NO2	B	2000	2	-	0/0/0/0	0/0/0/0
2	HEM	C	1000	1,3	-	0/10/54/54	0/0/8/8
3	NO2	C	2000	2	-	0/0/0/0	0/0/0/0
2	HEM	D	1000	1,3	-	0/10/54/54	0/0/8/8
3	NO2	D	2000	2	-	0/0/0/0	0/0/0/0
2	HEM	E	1000	1,3	-	0/10/54/54	0/0/8/8
3	NO2	E	2000	2	-	0/0/0/0	0/0/0/0
2	HEM	F	1000	1,3	-	0/10/54/54	0/0/8/8
3	NO2	F	2000	2	-	0/0/0/0	0/0/0/0
2	HEM	G	1000	1,3	-	0/10/54/54	0/0/8/8
3	NO2	G	2000	2	-	0/0/0/0	0/0/0/0
2	HEM	H	1000	1,3	-	0/10/54/54	0/0/8/8
3	NO2	H	2000	2	-	0/0/0/0	0/0/0/0
2	HEM	I	1000	1,3	-	0/10/54/54	0/0/8/8
3	NO2	I	2000	2	-	0/0/0/0	0/0/0/0
2	HEM	J	1000	1,3	-	0/10/54/54	0/0/8/8
3	NO2	J	2000	2	-	0/0/0/0	0/0/0/0
2	HEM	K	1000	1,3	-	0/10/54/54	0/0/8/8
3	NO2	K	2000	2	-	0/0/0/0	0/0/0/0
2	HEM	L	1000	1,3	-	0/10/54/54	0/0/8/8
3	NO2	L	2000	2	-	0/0/0/0	0/0/0/0
2	HEM	M	1000	1,3	-	0/10/54/54	0/0/8/8
3	NO2	M	2000	2	-	0/0/0/0	0/0/0/0
2	HEM	N	1000	1,3	-	0/10/54/54	0/0/8/8
3	NO2	N	2000	2	-	0/0/0/0	0/0/0/0
2	HEM	O	1000	1,3	-	0/10/54/54	0/0/8/8
3	NO2	O	2000	2	-	0/0/0/0	0/0/0/0
2	HEM	P	1000	1,3	-	0/10/54/54	0/0/8/8
3	NO2	P	2000	2	-	0/0/0/0	0/0/0/0
2	HEM	Q	1000	1,3	-	0/10/54/54	0/0/8/8
3	NO2	Q	2000	2	-	0/0/0/0	0/0/0/0
2	HEM	R	1000	1,3	-	0/10/54/54	0/0/8/8
3	NO2	R	2000	2	-	0/0/0/0	0/0/0/0
2	HEM	S	1000	1,3	-	0/10/54/54	0/0/8/8
3	NO2	S	2000	2	-	0/0/0/0	0/0/0/0
2	HEM	T	1000	1,3	-	0/10/54/54	0/0/8/8
3	NO2	T	2000	2	-	0/0/0/0	0/0/0/0

The worst 5 of 135 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1000	HEM	C3B-C4B	-9.26	1.43	1.51
2	B	1000	HEM	C3B-C4B	-8.15	1.44	1.51
2	A	1000	HEM	C3B-C4B	-7.78	1.44	1.51
2	J	1000	HEM	C3D-C4D	-7.15	1.42	1.51
2	T	1000	HEM	C3B-C4B	-6.98	1.45	1.51

The worst 5 of 225 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	M	1000	HEM	C3B-CAB-CBB	-9.41	110.03	124.46
2	Q	1000	HEM	C3B-CAB-CBB	-9.27	110.23	124.46
2	B	1000	HEM	C3B-CAB-CBB	-8.59	111.28	124.46
2	G	1000	HEM	C3C-CAC-CBC	-6.98	113.75	124.46
2	E	1000	HEM	C3C-CAC-CBC	-6.53	114.44	124.46

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

23 monomers are involved in 168 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1000	HEM	8	0
2	B	1000	HEM	6	0
2	C	1000	HEM	7	0
3	C	2000	NO2	1	0
2	D	1000	HEM	9	0
2	E	1000	HEM	8	0
2	F	1000	HEM	11	0
2	G	1000	HEM	15	0
3	G	2000	NO2	1	0
2	H	1000	HEM	9	0
2	I	1000	HEM	7	0
2	J	1000	HEM	10	0
2	K	1000	HEM	5	0
2	L	1000	HEM	7	0
3	L	2000	NO2	1	0
2	M	1000	HEM	6	0
2	N	1000	HEM	10	0
2	O	1000	HEM	5	0
2	P	1000	HEM	3	0
2	Q	1000	HEM	10	0
2	R	1000	HEM	14	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	S	1000	HEM	9	0
2	T	1000	HEM	6	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^{th} percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å^2)	Q<0.9
1	A	241/248 (97%)	-0.78	0 100 100	11, 24, 52, 62	0
1	B	241/248 (97%)	-0.74	1 (0%) 93 80	13, 24, 48, 70	0
1	C	241/248 (97%)	-0.68	2 (0%) 87 67	11, 23, 50, 68	0
1	D	241/248 (97%)	-0.78	0 100 100	14, 27, 54, 77	0
1	E	241/248 (97%)	-0.82	0 100 100	9, 21, 47, 64	0
1	F	241/248 (97%)	-0.51	0 100 100	17, 40, 75, 96	0
1	G	241/248 (97%)	-0.68	0 100 100	18, 34, 61, 90	0
1	H	241/248 (97%)	-0.77	0 100 100	14, 27, 55, 74	0
1	I	241/248 (97%)	-0.70	0 100 100	13, 28, 62, 80	0
1	J	241/248 (97%)	-0.67	0 100 100	15, 32, 69, 88	0
1	K	241/248 (97%)	-0.65	0 100 100	19, 34, 64, 87	0
1	L	241/248 (97%)	-0.62	0 100 100	17, 31, 60, 81	0
1	M	241/248 (97%)	-0.68	0 100 100	16, 29, 56, 79	0
1	N	241/248 (97%)	-0.59	0 100 100	19, 36, 71, 82	0
1	O	241/248 (97%)	-0.52	0 100 100	22, 41, 71, 96	0
1	P	241/248 (97%)	-0.30	0 100 100	38, 62, 108, 134	0
1	Q	241/248 (97%)	-0.49	1 (0%) 93 80	23, 52, 99, 120	0
1	R	241/248 (97%)	-0.30	0 100 100	27, 56, 101, 124	0
1	S	241/248 (97%)	-0.12	4 (1%) 73 45	39, 68, 110, 129	0
1	T	241/248 (97%)	-0.23	0 100 100	32, 63, 111, 135	0
All	All	4820/4960 (97%)	-0.58	8 (0%) 95 87	9, 36, 86, 135	0

The worst 5 of 8 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	132	SER	3.7
1	S	132	SER	2.6
1	Q	132	SER	2.4
1	C	38	VAL	2.2
1	S	57	ASP	2.2

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, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å^2)	Q<0.9
3	NO2	C	2000	3/3	0.98	0.18	3.91	25,25,28,31	0
3	NO2	A	2000	3/3	0.99	0.22	3.70	29,29,31,34	0
3	NO2	R	2000	3/3	0.97	0.33	3.30	48,48,50,52	0
3	NO2	I	2000	3/3	0.98	0.20	3.22	27,27,29,32	0
3	NO2	B	2000	3/3	0.95	0.23	2.64	32,32,36,43	0
2	HEM	G	1000	43/43	0.96	0.20	2.42	21,26,36,39	0
3	NO2	N	2000	3/3	0.97	0.20	2.32	41,41,45,47	0
3	NO2	K	2000	3/3	0.98	0.17	2.03	34,34,36,39	0
4	CA	K	1001	1/1	0.94	0.18	1.70	26,26,26,26	0
2	HEM	D	1000	43/43	0.97	0.18	1.34	15,22,31,35	0
4	CA	M	1001	1/1	0.97	0.16	1.32	26,26,26,26	0
3	NO2	J	2000	3/3	0.99	0.16	1.26	29,29,30,31	0
3	NO2	G	2000	3/3	0.99	0.17	1.20	36,36,37,39	0
3	NO2	D	2000	3/3	0.99	0.16	1.19	30,30,32,34	0
3	NO2	M	2000	3/3	0.96	0.18	1.07	31,31,34,41	0
2	HEM	Q	1000	43/43	0.96	0.18	1.05	29,31,39,45	0
2	HEM	E	1000	43/43	0.97	0.16	1.04	11,16,23,30	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å^2)	Q<0.9
2	HEM	I	1000	43/43	0.97	0.18	1.03	13,19,27,32	0
2	HEM	A	1000	43/43	0.98	0.17	0.90	13,19,28,36	0
3	NO2	H	2000	3/3	0.99	0.17	0.89	28,28,30,33	0
2	HEM	M	1000	43/43	0.97	0.18	0.83	17,22,29,34	0
3	NO2	O	2000	3/3	0.99	0.14	0.81	40,40,42,43	0
2	HEM	K	1000	43/43	0.97	0.17	0.78	21,25,34,37	0
2	HEM	H	1000	43/43	0.97	0.17	0.77	12,19,29,33	0
2	HEM	B	1000	43/43	0.96	0.19	0.63	15,18,27,33	0
2	HEM	F	1000	43/43	0.96	0.20	0.56	24,33,42,47	0
4	CA	C	1001	1/1	0.96	0.14	0.55	20,20,20,20	0
2	HEM	C	1000	43/43	0.97	0.16	0.51	13,17,24,31	0
2	HEM	L	1000	43/43	0.97	0.17	0.50	18,22,30,35	0
3	NO2	Q	2000	3/3	0.98	0.16	0.44	45,45,45,49	0
2	HEM	J	1000	43/43	0.97	0.17	0.44	15,20,28,35	0
2	HEM	N	1000	43/43	0.97	0.17	0.41	22,31,40,48	0
3	NO2	S	2000	3/3	0.97	0.18	0.39	63,63,64,65	0
2	HEM	O	1000	43/43	0.97	0.17	0.32	27,34,42,46	0
4	CA	G	1001	1/1	0.92	0.14	0.29	24,24,24,24	0
3	NO2	F	2000	3/3	0.98	0.17	0.23	40,40,42,44	0
3	NO2	L	2000	3/3	0.99	0.13	0.11	31,31,32,35	0
2	HEM	R	1000	43/43	0.96	0.19	0.09	30,36,43,53	0
4	CA	O	1001	1/1	0.96	0.15	0.02	29,29,29,29	0
4	CA	J	1001	1/1	0.98	0.17	0.01	29,29,29,29	0
2	HEM	S	1000	43/43	0.94	0.20	-0.05	45,55,64,71	0
3	NO2	P	2000	3/3	0.98	0.15	-0.05	55,55,55,55	0
2	HEM	P	1000	43/43	0.97	0.17	-0.13	41,45,52,55	0
2	HEM	T	1000	43/43	0.96	0.20	-0.16	39,46,52,59	0
4	CA	L	1001	1/1	0.95	0.15	-0.55	25,25,25,25	0
4	CA	A	1001	1/1	0.97	0.12	-0.58	21,21,21,21	0
4	CA	H	1001	1/1	0.97	0.13	-0.63	22,22,22,22	0
4	CA	T	1001	1/1	0.94	0.16	-0.65	52,52,52,52	0
4	CA	N	1001	1/1	0.96	0.14	-0.78	32,32,32,32	0
4	CA	R	1001	1/1	0.94	0.14	-0.85	50,50,50,50	0
4	CA	D	1001	1/1	0.97	0.12	-0.88	24,24,24,24	0
3	NO2	T	2000	3/3	0.99	0.17	-1.13	57,57,57,58	0
4	CA	E	1001	1/1	0.98	0.12	-1.14	13,13,13,13	0
4	CA	P	1001	1/1	0.97	0.12	-1.19	54,54,54,54	0
4	CA	I	1001	1/1	0.94	0.11	-1.34	23,23,23,23	0
4	CA	S	1001	1/1	0.84	0.14	-1.61	60,60,60,60	0
4	CA	B	1001	1/1	0.98	0.12	-1.64	17,17,17,17	0
4	CA	F	1001	1/1	0.99	0.10	-1.80	30,30,30,30	0
4	CA	Q	1001	1/1	0.96	0.12	-1.99	48,48,48,48	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å^2)	Q<0.9
3	NO2	E	2000	3/3	1.00	0.10	-2.82	22,22,24,28	0

6.5 Other polymers (i)

There are no such residues in this entry.