



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:40 AM GMT

PDB ID : 2NSI
Title : HUMAN INDUCIBLE NITRIC OXIDE SYNTHASE, ZN-FREE, SEITU COMPLEX
Authors : Li, H.; Raman, C.S.; Glaser, C.B.; Blasko, E.; Young, T.A.; Parkinson, J.F.; Whitlow, M.; Poulos, T.L.
Deposited on : 1999-01-11
Resolution : 3.00 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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

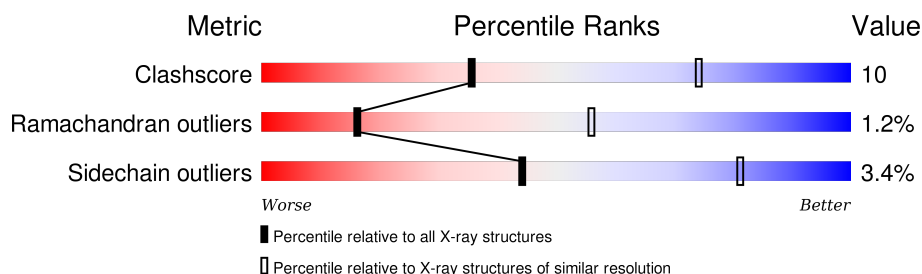
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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))
Clashscore	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (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 $\leq 5\%$. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	431	
1	B	431	
1	C	431	
1	D	431	

2 Entry composition [i](#)

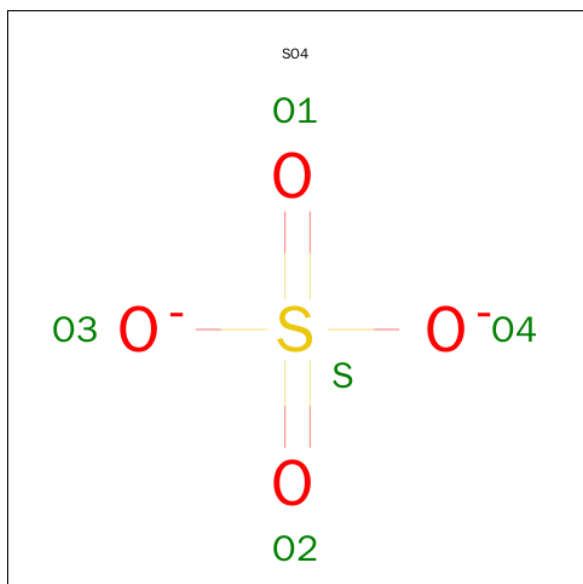
There are 5 unique types of molecules in this entry. The entry contains 13985 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PROTEIN (NITRIC OXIDE SYNTHASE).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	420	Total	C	N	O	S	0	0	0
			3419	2186	599	612	22			
1	B	420	Total	C	N	O	S	0	0	0
			3419	2186	599	612	22			
1	C	420	Total	C	N	O	S	0	0	0
			3419	2186	599	612	22			
1	D	420	Total	C	N	O	S	0	0	0
			3419	2186	599	612	22			

- Molecule 2 is SULFATE ION (three-letter code: SO₄) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		

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

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



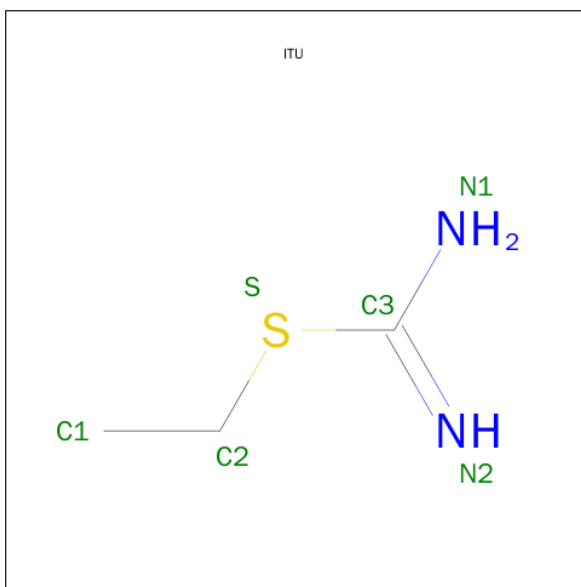
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
3	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
3	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
3	D	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 4 is 5,6,7,8-TETRAHYDROBIOPTERIN (three-letter code: H4B) (formula: $C_9H_{15}N_5O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			17	9	5	3		
4	B	1	Total	C	N	O	0	0
			17	9	5	3		
4	C	1	Total	C	N	O	0	0
			17	9	5	3		
4	D	1	Total	C	N	O	0	0
			17	9	5	3		

- Molecule 5 is ETHYLISOTHIOUREA (three-letter code: ITU) (formula: $C_3H_8N_2S$).



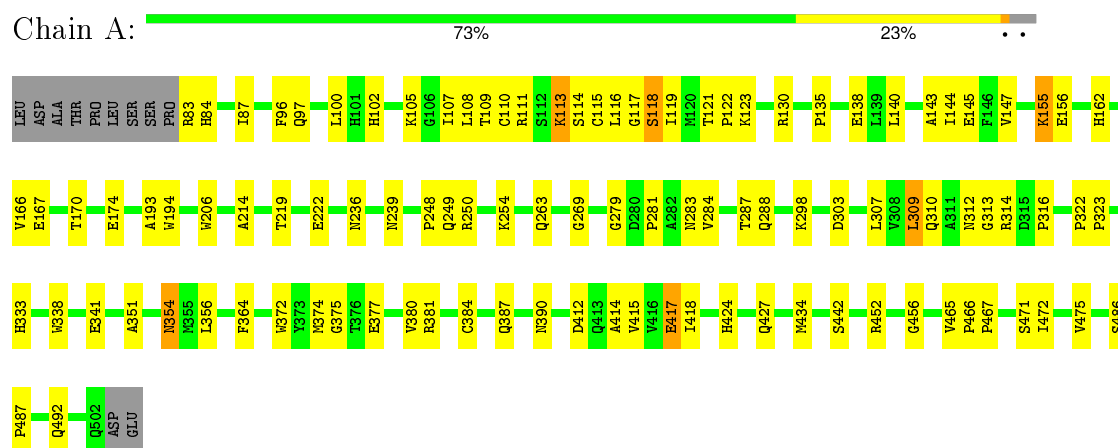
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	S	0	0
			6	3	2	1		
5	B	1	Total	C	N	S	0	0
			6	3	2	1		
5	C	1	Total	C	N	S	0	0
			6	3	2	1		
5	D	1	Total	C	N	S	0	0
			6	3	2	1		

3 Residue-property plots [i](#)

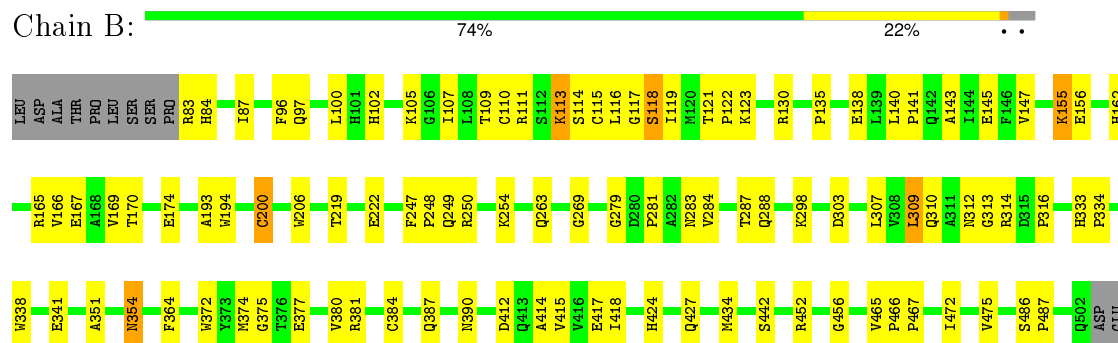
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

Note EDS was not executed.

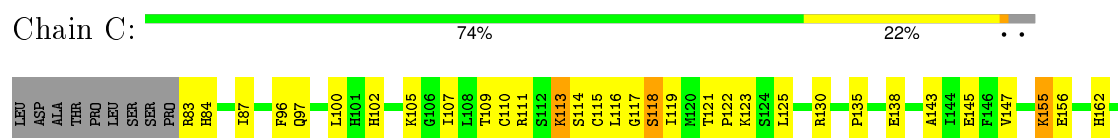
• Molecule 1: PROTEIN (NITRIC OXIDE SYNTHASE)

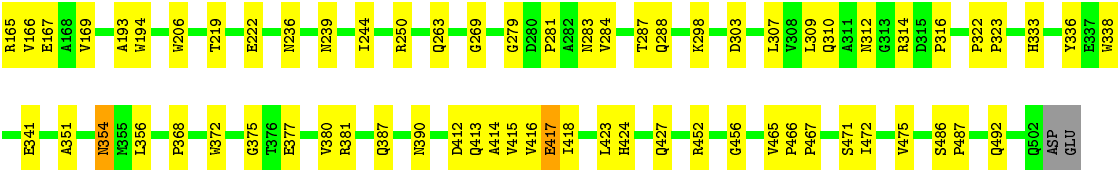


• Molecule 1: PROTEIN (NITRIC OXIDE SYNTHASE)



• Molecule 1: PROTEIN (NITRIC OXIDE SYNTHASE)





● Molecule 1: PROTEIN (NITRIC OXIDE SYNTHASE)



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	188.37Å 188.37Å 230.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 3.00	Depositor
% Data completeness (in resolution range)	80.8 (30.00-3.00)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
Refinement program	X-PLOR 3.8	Depositor
R, R_{free}	0.214 , 0.238	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	13985	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

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: HEM, ITU, H4B, SO4

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	0.37	0/3518	0.58	0/4774
1	B	0.37	0/3518	0.59	1/4774 (0.0%)
1	C	0.38	0/3518	0.59	0/4774
1	D	0.37	0/3518	0.58	0/4774
All	All	0.37	0/14072	0.58	1/19096 (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	B	200	CYS	CA-CB-SG	-5.68	103.77	114.00

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	3419	0	3327	71	0
1	B	3419	0	3327	70	0
1	C	3419	0	3327	66	0
1	D	3419	0	3327	73	0
2	A	15	0	0	1	0
2	B	10	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	10	0	0	0	0
2	D	10	0	0	0	0
3	A	43	0	30	1	0
3	B	43	0	30	2	0
3	C	43	0	30	1	0
3	D	43	0	30	1	0
4	A	17	0	15	0	0
4	B	17	0	15	0	0
4	C	17	0	15	0	0
4	D	17	0	15	0	0
5	A	6	0	7	0	0
5	B	6	0	7	0	0
5	C	6	0	7	0	0
5	D	6	0	7	0	0
All	All	13985	0	13516	267	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 10.

All (267) 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:A:115:CYS:HB2	1:B:110:CYS:SG	2.11	0.91
1:A:110:CYS:SG	1:B:115:CYS:HB2	2.17	0.85
1:B:83:ARG:HB3	1:B:102:HIS:HE1	1.45	0.80
1:D:83:ARG:HB3	1:D:102:HIS:HE1	1.45	0.80
1:C:83:ARG:HB3	1:C:102:HIS:HE1	1.46	0.79
1:A:83:ARG:HB3	1:A:102:HIS:HE1	1.47	0.78
1:B:269:GLY:H	1:B:287:THR:HG21	1.51	0.76
1:C:109:THR:HG22	1:C:116:LEU:HG	1.67	0.75
1:A:269:GLY:H	1:A:287:THR:HG21	1.52	0.73
1:D:109:THR:HG22	1:D:116:LEU:HG	1.70	0.73
1:D:269:GLY:H	1:D:287:THR:HG21	1.53	0.72
1:A:109:THR:HG22	1:A:116:LEU:HG	1.70	0.72
1:C:269:GLY:H	1:C:287:THR:HG21	1.54	0.72
1:C:115:CYS:HB2	1:D:110:CYS:SG	2.31	0.69
1:B:109:THR:HG22	1:B:116:LEU:HG	1.72	0.69
1:C:110:CYS:SG	1:D:115:CYS:HB2	2.33	0.68
1:D:110:CYS:HB3	1:D:116:LEU:H	1.57	0.68
1:C:372:TRP:H	3:C:550:HEM:HAB	1.59	0.67
1:B:83:ARG:HB3	1:B:102:HIS:CE1	2.29	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:83:ARG:HB3	1:D:102:HIS:CE1	2.28	0.67
1:A:372:TRP:H	3:A:550:HEM:HAB	1.59	0.66
1:D:372:TRP:H	3:D:550:HEM:HAB	1.59	0.66
1:B:110:CYS:HB3	1:B:116:LEU:H	1.61	0.66
1:A:110:CYS:HB3	1:A:116:LEU:H	1.62	0.65
1:C:83:ARG:HB3	1:C:102:HIS:CE1	2.31	0.65
1:A:83:ARG:HB3	1:A:102:HIS:CE1	2.31	0.65
1:B:372:TRP:H	3:B:550:HEM:HAB	1.62	0.65
1:B:155:LYS:H	1:B:155:LYS:HD3	1.63	0.62
1:B:354:ASN:HD22	1:B:354:ASN:H	1.48	0.61
1:D:155:LYS:H	1:D:155:LYS:HD3	1.66	0.61
1:A:354:ASN:HD22	1:A:354:ASN:H	1.49	0.61
1:C:354:ASN:HD22	1:C:354:ASN:H	1.47	0.61
1:C:414:ALA:O	1:C:418:ILE:HG13	2.01	0.61
1:C:110:CYS:HB3	1:C:116:LEU:H	1.65	0.60
1:D:354:ASN:H	1:D:354:ASN:HD22	1.49	0.59
1:C:110:CYS:HB2	1:C:115:CYS:HA	1.84	0.59
1:B:412:ASP:O	1:B:415:VAL:HG12	2.03	0.59
1:C:155:LYS:HD3	1:C:155:LYS:H	1.68	0.59
1:B:219:THR:HG23	1:B:222:GLU:H	1.68	0.59
1:C:219:THR:HG23	1:C:222:GLU:H	1.67	0.59
1:A:155:LYS:H	1:A:155:LYS:HD3	1.66	0.58
1:A:380:VAL:HG11	1:A:467:PRO:HB2	1.86	0.58
1:C:380:VAL:HG11	1:C:467:PRO:HB2	1.85	0.57
1:A:219:THR:HG23	1:A:222:GLU:H	1.69	0.57
1:C:354:ASN:HD22	1:C:354:ASN:N	2.03	0.57
1:C:110:CYS:HA	1:C:116:LEU:HD23	1.87	0.56
1:D:380:VAL:HG11	1:D:467:PRO:HB2	1.88	0.56
1:B:380:VAL:HG11	1:B:467:PRO:HB2	1.88	0.56
1:B:135:PRO:HB2	1:B:138:GLU:HG2	1.88	0.56
1:C:135:PRO:HB2	1:C:138:GLU:HG2	1.88	0.56
1:B:414:ALA:O	1:B:418:ILE:HG13	2.07	0.55
1:C:263:GLN:HA	1:C:351:ALA:O	2.07	0.55
1:C:412:ASP:O	1:C:415:VAL:HG12	2.07	0.55
1:A:414:ALA:O	1:A:418:ILE:HG13	2.07	0.55
1:A:412:ASP:O	1:A:415:VAL:HG12	2.06	0.54
1:B:263:GLN:HA	1:B:351:ALA:O	2.07	0.54
1:A:110:CYS:HB2	1:A:115:CYS:HA	1.89	0.54
1:D:219:THR:HG23	1:D:222:GLU:H	1.72	0.54
1:C:279:GLY:O	1:C:281:PRO:HD3	2.08	0.54
1:B:283:ASN:O	1:B:287:THR:HG23	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:354:ASN:HD22	1:B:354:ASN:N	2.04	0.53
1:B:110:CYS:HB2	1:B:115:CYS:HA	1.89	0.53
1:A:354:ASN:HD22	1:A:354:ASN:N	2.06	0.53
1:D:110:CYS:HB3	1:D:116:LEU:N	2.22	0.53
1:C:110:CYS:CB	1:C:116:LEU:H	2.21	0.53
1:C:283:ASN:O	1:C:287:THR:HG23	2.09	0.53
1:D:354:ASN:N	1:D:354:ASN:HD22	2.07	0.52
1:A:135:PRO:HB2	1:A:138:GLU:HG2	1.91	0.52
1:D:143:ALA:O	1:D:147:VAL:HG23	2.08	0.52
1:A:263:GLN:HA	1:A:351:ALA:O	2.09	0.52
1:B:111:ARG:HG3	1:B:113:LYS:HG2	1.92	0.52
1:D:263:GLN:HA	1:D:351:ALA:O	2.10	0.52
1:D:412:ASP:O	1:D:415:VAL:HG12	2.09	0.52
1:D:110:CYS:CB	1:D:116:LEU:H	2.22	0.52
1:B:110:CYS:HA	1:B:116:LEU:HD23	1.92	0.52
1:D:110:CYS:HB2	1:D:115:CYS:HA	1.91	0.52
1:D:465:VAL:HG22	1:D:475:VAL:HG23	1.92	0.51
1:A:110:CYS:HA	1:A:116:LEU:HD23	1.92	0.51
1:D:117:GLY:O	1:D:118:SER:CB	2.59	0.51
1:C:193:ALA:HB2	1:C:487:PRO:HB2	1.92	0.51
1:C:119:ILE:HB	1:C:122:PRO:HG3	1.92	0.51
1:A:110:CYS:CB	1:A:116:LEU:H	2.23	0.51
1:A:111:ARG:HG3	1:A:113:LYS:HG2	1.93	0.51
1:B:110:CYS:CB	1:B:116:LEU:H	2.23	0.50
1:B:281:PRO:HB2	1:B:387:GLN:O	2.11	0.50
1:A:143:ALA:O	1:A:147:VAL:HG23	2.11	0.50
1:A:283:ASN:O	1:A:287:THR:HG23	2.11	0.50
1:D:135:PRO:HB2	1:D:138:GLU:HG2	1.93	0.50
1:A:117:GLY:O	1:A:118:SER:CB	2.59	0.50
1:B:110:CYS:HB3	1:B:116:LEU:N	2.26	0.49
1:B:452:ARG:NH1	1:B:452:ARG:HG2	2.27	0.49
1:D:194:TRP:CE3	1:D:206:TRP:HA	2.47	0.49
1:B:117:GLY:O	1:B:118:SER:CB	2.60	0.49
1:A:110:CYS:HB3	1:A:116:LEU:N	2.27	0.49
1:C:117:GLY:O	1:C:118:SER:CB	2.61	0.49
1:A:193:ALA:HB2	1:A:487:PRO:HB2	1.94	0.48
1:C:143:ALA:O	1:C:147:VAL:HG23	2.12	0.48
1:D:193:ALA:HB2	1:D:487:PRO:HB2	1.95	0.48
1:A:119:ILE:HB	1:A:122:PRO:HG3	1.96	0.48
1:C:194:TRP:CE3	1:C:206:TRP:HA	2.48	0.48
1:D:162:HIS:O	1:D:166:VAL:HG23	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:105:LYS:HG3	1:C:107:ILE:HG12	1.96	0.48
1:D:110:CYS:HA	1:D:116:LEU:HD23	1.96	0.48
1:C:377:GLU:O	1:C:381:ARG:HB2	2.14	0.48
1:D:111:ARG:HG3	1:D:113:LYS:HG2	1.95	0.48
1:C:111:ARG:HG3	1:C:113:LYS:HG2	1.94	0.47
1:A:466:PRO:HA	1:A:467:PRO:HD3	1.78	0.47
1:B:298:LYS:HB2	1:B:303:ASP:OD2	2.14	0.47
1:D:140:LEU:O	1:D:144:ILE:HG12	2.14	0.47
1:A:471:SER:HB2	1:B:467:PRO:HB3	1.95	0.47
1:C:452:ARG:HG2	1:C:452:ARG:NH1	2.30	0.47
1:D:283:ASN:O	1:D:287:THR:HG23	2.15	0.47
1:B:194:TRP:CE3	1:B:206:TRP:HA	2.50	0.47
1:A:105:LYS:HG3	1:A:107:ILE:HG12	1.96	0.47
1:D:298:LYS:HB2	1:D:303:ASP:OD2	2.14	0.47
1:A:281:PRO:HB2	1:A:387:GLN:O	2.14	0.47
1:D:105:LYS:HG3	1:D:107:ILE:HG12	1.95	0.47
1:A:452:ARG:NH1	1:A:452:ARG:HG2	2.30	0.46
1:A:162:HIS:O	1:A:166:VAL:HG23	2.15	0.46
1:D:219:THR:HG22	1:D:222:GLU:HG3	1.97	0.46
1:B:193:ALA:HB2	1:B:487:PRO:HB2	1.97	0.46
1:B:84:HIS:HB3	1:B:97:GLN:NE2	2.31	0.46
1:B:87:ILE:HG13	1:B:96:PHE:HB2	1.98	0.46
1:D:414:ALA:O	1:D:418:ILE:HG13	2.15	0.46
1:B:380:VAL:O	1:B:384:CYS:HB2	2.16	0.46
1:C:486:SER:HA	1:C:487:PRO:C	2.36	0.46
1:D:130:ARG:HG2	1:D:250:ARG:HD3	1.98	0.46
1:B:424:HIS:HA	1:B:427:GLN:HE21	1.80	0.46
1:C:110:CYS:HB3	1:C:116:LEU:N	2.30	0.46
1:A:219:THR:HG22	1:A:222:GLU:HG3	1.97	0.46
1:A:281:PRO:O	1:A:284:VAL:HG23	2.15	0.46
1:B:143:ALA:O	1:B:147:VAL:HG23	2.15	0.46
1:D:424:HIS:HA	1:D:427:GLN:HE21	1.81	0.46
1:A:465:VAL:HG22	1:A:475:VAL:HG23	1.98	0.46
1:A:84:HIS:HB3	1:A:97:GLN:NE2	2.31	0.46
1:C:413:GLN:HG3	1:D:416:VAL:HG11	1.98	0.46
1:A:194:TRP:CE3	1:A:206:TRP:HA	2.50	0.46
1:C:336:TYR:CD1	1:C:336:TYR:N	2.84	0.45
1:A:140:LEU:O	1:A:144:ILE:HG12	2.16	0.45
1:A:111:ARG:C	1:B:111:ARG:HA	2.36	0.45
1:C:244:ILE:HG13	1:C:368:PRO:O	2.16	0.45
1:B:130:ARG:HG2	1:B:250:ARG:HD3	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:119:ILE:HB	1:D:122:PRO:HG3	1.97	0.45
1:B:338:TRP:O	1:B:341:GLU:HB2	2.16	0.45
1:B:281:PRO:O	1:B:284:VAL:HG23	2.17	0.45
1:A:377:GLU:O	1:A:381:ARG:HB2	2.16	0.45
1:A:322:PRO:HA	1:A:323:PRO:HD2	1.80	0.45
1:D:84:HIS:HB3	1:D:97:GLN:NE2	2.32	0.45
1:D:309:LEU:O	1:D:316:PRO:HA	2.16	0.45
1:D:279:GLY:O	1:D:281:PRO:HD3	2.16	0.45
1:B:249:GLN:HB3	1:B:364:PHE:CE2	2.52	0.45
1:A:380:VAL:O	1:A:384:CYS:HB2	2.17	0.44
1:D:380:VAL:O	1:D:384:CYS:HB2	2.18	0.44
1:C:471:SER:HB2	1:D:467:PRO:HB3	2.00	0.44
1:B:100:LEU:CB	1:B:456:GLY:HA3	2.47	0.44
1:B:119:ILE:HB	1:B:122:PRO:HG3	1.98	0.44
1:D:356:LEU:HB2	1:D:492:GLN:CD	2.37	0.44
1:A:87:ILE:HG13	1:A:96:PHE:HB2	1.98	0.44
1:C:130:ARG:HG2	1:C:250:ARG:HD3	1.99	0.44
1:B:107:ILE:HG13	1:B:107:ILE:O	2.17	0.44
1:D:423:LEU:O	1:D:427:GLN:HG3	2.18	0.44
1:D:87:ILE:HG13	1:D:96:PHE:HB2	1.98	0.44
1:A:130:ARG:HG2	1:A:250:ARG:HD3	1.99	0.44
1:C:84:HIS:HB3	1:C:97:GLN:NE2	2.32	0.44
1:A:108:LEU:HD13	1:B:113:LYS:HA	2.00	0.44
1:B:472:ILE:HG22	1:B:472:ILE:O	2.18	0.44
1:B:100:LEU:HB3	1:B:456:GLY:HA3	1.98	0.44
1:B:465:VAL:HG22	1:B:475:VAL:HG23	2.00	0.44
1:B:310:GLN:HG3	1:B:314:ARG:O	2.18	0.44
1:B:140:LEU:HB3	1:B:141:PRO:HD3	2.00	0.44
1:D:254:LYS:O	1:D:313:GLY:HA3	2.18	0.44
1:D:377:GLU:O	1:D:381:ARG:HB2	2.18	0.44
1:A:279:GLY:O	1:A:281:PRO:HD3	2.17	0.43
1:D:281:PRO:O	1:D:284:VAL:HG23	2.18	0.43
1:D:452:ARG:HG2	1:D:452:ARG:NH1	2.34	0.43
1:A:333:HIS:NE2	1:A:417:GLU:HG3	2.33	0.43
1:B:219:THR:HG22	1:B:222:GLU:HG3	2.00	0.43
1:C:281:PRO:HB2	1:C:387:GLN:O	2.18	0.43
1:A:111:ARG:HA	1:B:111:ARG:C	2.37	0.43
1:B:452:ARG:HG2	1:B:452:ARG:HH11	1.83	0.43
1:D:249:GLN:HB3	1:D:364:PHE:CE2	2.52	0.43
1:A:310:GLN:HG3	1:A:314:ARG:O	2.18	0.43
1:A:298:LYS:HB2	1:A:303:ASP:OD2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:338:TRP:O	1:D:341:GLU:HB2	2.18	0.43
1:D:247:PHE:HB3	1:D:248:PRO:HD2	2.01	0.43
1:C:322:PRO:HA	1:C:323:PRO:HD2	1.83	0.43
1:C:219:THR:HG22	1:C:222:GLU:HG3	2.00	0.43
1:A:111:ARG:HA	1:B:111:ARG:HA	2.00	0.43
1:B:170:THR:O	1:B:174:GLU:HG2	2.18	0.43
1:B:105:LYS:HG3	1:B:107:ILE:HG12	1.99	0.43
1:B:162:HIS:O	1:B:166:VAL:HG23	2.18	0.43
1:C:281:PRO:O	1:C:284:VAL:HG23	2.19	0.43
1:B:279:GLY:O	1:B:281:PRO:HD3	2.18	0.43
1:C:107:ILE:HG13	1:C:107:ILE:O	2.19	0.43
1:D:307:LEU:HG	1:D:309:LEU:HD11	2.01	0.43
1:A:100:LEU:HB3	1:A:456:GLY:HA3	2.00	0.43
1:A:338:TRP:O	1:A:341:GLU:HB2	2.19	0.43
1:D:135:PRO:HA	1:D:136:PRO:HD3	1.92	0.42
1:C:333:HIS:NE2	1:C:417:GLU:HG3	2.34	0.42
1:A:333:HIS:HE2	1:A:417:GLU:HG3	1.83	0.42
1:A:236:ASN:ND2	2:A:930:SO4:O2	2.51	0.42
1:C:310:GLN:HG3	1:C:314:ARG:O	2.19	0.42
1:A:100:LEU:CB	1:A:456:GLY:HA3	2.49	0.42
1:D:100:LEU:CB	1:D:456:GLY:HA3	2.49	0.42
1:D:465:VAL:HA	1:D:466:PRO:HD3	1.88	0.42
1:B:309:LEU:O	1:B:316:PRO:HA	2.19	0.42
1:B:333:HIS:CG	1:B:334:PRO:HD2	2.54	0.42
1:B:155:LYS:H	1:B:155:LYS:CD	2.25	0.42
1:A:486:SER:HA	1:A:487:PRO:C	2.40	0.42
1:B:486:SER:HA	1:B:487:PRO:C	2.39	0.42
1:D:281:PRO:HB2	1:D:387:GLN:O	2.20	0.42
1:D:459:ALA:HB3	1:D:480:MET:HB2	2.01	0.42
1:A:249:GLN:HB3	1:A:364:PHE:CE2	2.55	0.42
1:C:333:HIS:HE2	1:C:417:GLU:HG3	1.84	0.42
1:B:333:HIS:NE2	1:B:417:GLU:HG3	2.35	0.42
1:C:162:HIS:O	1:C:166:VAL:HG23	2.20	0.42
1:C:309:LEU:O	1:C:316:PRO:HA	2.20	0.42
1:B:254:LYS:O	1:B:313:GLY:HA3	2.19	0.42
1:D:333:HIS:NE2	1:D:417:GLU:HG3	2.35	0.42
1:C:356:LEU:HB2	1:C:492:GLN:CD	2.39	0.42
1:A:424:HIS:HA	1:A:427:GLN:HE21	1.84	0.42
1:A:214:ALA:O	1:A:248:PRO:HD3	2.20	0.42
1:C:87:ILE:HG13	1:C:96:PHE:HB2	2.00	0.42
1:A:465:VAL:HA	1:A:466:PRO:HD3	1.87	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:423:LEU:O	1:C:427:GLN:HG3	2.20	0.42
1:C:165:ARG:O	1:C:169:VAL:HG23	2.19	0.42
1:A:107:ILE:HG13	1:A:107:ILE:O	2.20	0.42
1:D:171:LYS:O	1:D:175:THR:HG23	2.20	0.42
1:A:254:LYS:O	1:A:313:GLY:HA3	2.20	0.42
1:C:236:ASN:HB3	1:C:239:ASN:O	2.20	0.42
1:D:310:GLN:HG3	1:D:314:ARG:O	2.19	0.42
1:A:356:LEU:HB2	1:A:492:GLN:CD	2.41	0.41
1:C:472:ILE:O	1:C:472:ILE:HG22	2.20	0.41
1:D:472:ILE:HG22	1:D:472:ILE:O	2.19	0.41
1:C:416:VAL:HG11	1:D:413:GLN:HG3	2.00	0.41
1:D:100:LEU:HB3	1:D:456:GLY:HA3	2.01	0.41
1:D:155:LYS:CD	1:D:155:LYS:H	2.29	0.41
1:B:377:GLU:O	1:B:381:ARG:HB2	2.21	0.41
1:A:472:ILE:O	1:A:472:ILE:HG22	2.19	0.41
1:B:200:CYS:HB2	3:B:550:HEM:ND	2.36	0.41
1:A:236:ASN:HB3	1:A:239:ASN:O	2.20	0.41
1:B:247:PHE:HB3	1:B:248:PRO:HD2	2.02	0.41
1:D:115:CYS:C	1:D:117:GLY:H	2.25	0.41
1:B:466:PRO:HA	1:B:467:PRO:HD3	1.76	0.41
1:C:109:THR:O	1:C:116:LEU:HG	2.21	0.41
1:C:466:PRO:HA	1:C:467:PRO:HD3	1.79	0.41
1:A:471:SER:CB	1:B:467:PRO:HB3	2.51	0.41
1:D:214:ALA:O	1:D:248:PRO:HD3	2.20	0.41
1:C:424:HIS:HA	1:C:427:GLN:HE21	1.86	0.41
1:C:465:VAL:HG22	1:C:475:VAL:HG23	2.02	0.41
1:C:338:TRP:O	1:C:341:GLU:HB2	2.21	0.41
1:C:100:LEU:HB3	1:C:456:GLY:HA3	2.03	0.41
1:A:374:MET:HA	1:A:434:MET:O	2.21	0.41
1:A:170:THR:O	1:A:174:GLU:HG2	2.20	0.41
1:B:374:MET:HA	1:B:434:MET:O	2.20	0.41
1:D:296:LYS:HA	1:D:297:PRO:HD2	1.91	0.41
1:D:333:HIS:HE2	1:D:417:GLU:HG3	1.86	0.40
1:D:244:ILE:HG13	1:D:368:PRO:O	2.21	0.40
1:A:309:LEU:O	1:A:316:PRO:HA	2.20	0.40
1:B:165:ARG:O	1:B:169:VAL:HG23	2.21	0.40
1:D:466:PRO:HA	1:D:467:PRO:HD3	1.77	0.40
1:C:336:TYR:CE2	1:C:417:GLU:HG2	2.56	0.40
1:D:433:ILE:HG13	1:D:434:MET:N	2.36	0.40
1:C:298:LYS:HB2	1:C:303:ASP:OD2	2.21	0.40
1:C:115:CYS:C	1:C:117:GLY:H	2.25	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:122:PRO:HD2	1:C:125:LEU:HD12	2.03	0.40
1:D:107:ILE:O	1:D:107:ILE:HG13	2.22	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	418/431 (97%)	392 (94%)	21 (5%)	5 (1%)	16	56
1	B	418/431 (97%)	387 (93%)	26 (6%)	5 (1%)	16	56
1	C	418/431 (97%)	389 (93%)	24 (6%)	5 (1%)	16	56
1	D	418/431 (97%)	390 (93%)	23 (6%)	5 (1%)	16	56
All	All	1672/1724 (97%)	1558 (93%)	94 (6%)	20 (1%)	16	56

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	118	SER
1	B	118	SER
1	C	118	SER
1	D	118	SER
1	A	114	SER
1	A	390	ASN
1	B	114	SER
1	B	390	ASN
1	C	114	SER
1	D	114	SER
1	D	390	ASN
1	A	113	LYS
1	B	113	LYS

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Mol	Chain	Res	Type
1	C	113	LYS
1	C	390	ASN
1	D	113	LYS
1	C	375	GLY
1	A	375	GLY
1	D	375	GLY
1	B	375	GLY

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	366/376 (97%)	353 (96%)	13 (4%)	42	79
1	B	366/376 (97%)	354 (97%)	12 (3%)	45	82
1	C	366/376 (97%)	355 (97%)	11 (3%)	48	83
1	D	366/376 (97%)	352 (96%)	14 (4%)	40	78
All	All	1464/1504 (97%)	1414 (97%)	50 (3%)	44	81

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	121	THR
1	A	123	LYS
1	A	145	GLU
1	A	155	LYS
1	A	156	GLU
1	A	167	GLU
1	A	288	GLN
1	A	307	LEU
1	A	309	LEU
1	A	312	ASN
1	A	354	ASN
1	A	417	GLU
1	A	442	SER
1	B	121	THR

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Mol	Chain	Res	Type
1	B	123	LYS
1	B	145	GLU
1	B	155	LYS
1	B	156	GLU
1	B	167	GLU
1	B	288	GLN
1	B	307	LEU
1	B	309	LEU
1	B	312	ASN
1	B	354	ASN
1	B	442	SER
1	C	121	THR
1	C	123	LYS
1	C	145	GLU
1	C	155	LYS
1	C	156	GLU
1	C	167	GLU
1	C	288	GLN
1	C	307	LEU
1	C	312	ASN
1	C	354	ASN
1	C	417	GLU
1	D	121	THR
1	D	123	LYS
1	D	145	GLU
1	D	155	LYS
1	D	156	GLU
1	D	167	GLU
1	D	288	GLN
1	D	307	LEU
1	D	309	LEU
1	D	312	ASN
1	D	354	ASN
1	D	395	VAL
1	D	417	GLU
1	D	442	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (56) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	97	GLN
1	A	102	HIS

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Mol	Chain	Res	Type
1	A	142	GLN
1	A	192	GLN
1	A	205	GLN
1	A	210	GLN
1	A	249	GLN
1	A	283	ASN
1	A	354	ASN
1	A	390	ASN
1	A	424	HIS
1	A	427	GLN
1	A	429	GLN
1	A	482	ASN
1	A	502	GLN
1	B	97	GLN
1	B	102	HIS
1	B	142	GLN
1	B	192	GLN
1	B	205	GLN
1	B	249	GLN
1	B	283	ASN
1	B	354	ASN
1	B	390	ASN
1	B	424	HIS
1	B	427	GLN
1	B	429	GLN
1	B	482	ASN
1	B	502	GLN
1	C	97	GLN
1	C	102	HIS
1	C	142	GLN
1	C	192	GLN
1	C	205	GLN
1	C	210	GLN
1	C	249	GLN
1	C	354	ASN
1	C	390	ASN
1	C	424	HIS
1	C	427	GLN
1	C	429	GLN
1	C	482	ASN
1	C	502	GLN
1	D	97	GLN

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Mol	Chain	Res	Type
1	D	102	HIS
1	D	142	GLN
1	D	205	GLN
1	D	249	GLN
1	D	283	ASN
1	D	354	ASN
1	D	390	ASN
1	D	424	HIS
1	D	427	GLN
1	D	429	GLN
1	D	482	ASN
1	D	502	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

21 ligands 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	$\# Z > 2$
3	HEM	A	550	1	30,50,50	2.74	8 (26%)	24,82,82	2.20	8 (33%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	H4B	A	600	-	13,18,18	1.89	3 (23%)	11,26,26	4.10	6 (54%)
5	ITU	A	800	-	4,5,5	0.98	0	4,5,5	2.68	1 (25%)
2	SO4	A	910	-	4,4,4	1.15	0	6,6,6	0.50	0
2	SO4	A	920	-	4,4,4	1.01	0	6,6,6	0.43	0
2	SO4	A	930	-	4,4,4	1.02	0	6,6,6	0.28	0
3	HEM	B	550	1	30,50,50	2.86	8 (26%)	24,82,82	2.25	9 (37%)
4	H4B	B	601	-	13,18,18	1.98	4 (30%)	11,26,26	4.08	6 (54%)
5	ITU	B	801	-	4,5,5	0.94	0	4,5,5	2.31	1 (25%)
2	SO4	B	911	-	4,4,4	1.12	0	6,6,6	0.39	0
2	SO4	B	921	-	4,4,4	0.99	0	6,6,6	0.38	0
3	HEM	C	550	1	30,50,50	2.86	8 (26%)	24,82,82	2.20	8 (33%)
4	H4B	C	602	-	13,18,18	1.86	3 (23%)	11,26,26	4.12	6 (54%)
5	ITU	C	802	-	4,5,5	0.99	0	4,5,5	2.52	1 (25%)
2	SO4	C	912	-	4,4,4	1.03	0	6,6,6	0.40	0
2	SO4	C	922	-	4,4,4	0.98	0	6,6,6	0.38	0
3	HEM	D	550	1	30,50,50	2.90	8 (26%)	24,82,82	2.21	9 (37%)
4	H4B	D	603	-	13,18,18	1.95	4 (30%)	11,26,26	4.11	6 (54%)
5	ITU	D	803	-	4,5,5	0.97	0	4,5,5	2.43	1 (25%)
2	SO4	D	913	-	4,4,4	1.11	0	6,6,6	0.35	0
2	SO4	D	923	-	4,4,4	1.06	0	6,6,6	0.34	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
3	HEM	A	550	1	-	0/10/54/54	0/0/8/8
4	H4B	A	600	-	-	0/8/17/17	0/2/2/2
5	ITU	A	800	-	-	0/3/3/3	0/0/0/0
2	SO4	A	910	-	-	0/0/0/0	0/0/0/0
2	SO4	A	920	-	-	0/0/0/0	0/0/0/0
2	SO4	A	930	-	-	0/0/0/0	0/0/0/0
3	HEM	B	550	1	-	0/10/54/54	0/0/8/8
4	H4B	B	601	-	-	0/8/17/17	0/2/2/2
5	ITU	B	801	-	-	0/3/3/3	0/0/0/0
2	SO4	B	911	-	-	0/0/0/0	0/0/0/0
2	SO4	B	921	-	-	0/0/0/0	0/0/0/0
3	HEM	C	550	1	-	0/10/54/54	0/0/8/8
4	H4B	C	602	-	-	0/8/17/17	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	ITU	C	802	-	-	0/3/3/3	0/0/0/0
2	SO4	C	912	-	-	0/0/0/0	0/0/0/0
2	SO4	C	922	-	-	0/0/0/0	0/0/0/0
3	HEM	D	550	1	-	0/10/54/54	0/0/8/8
4	H4B	D	603	-	-	0/8/17/17	0/2/2/2
5	ITU	D	803	-	-	0/3/3/3	0/0/0/0
2	SO4	D	913	-	-	0/0/0/0	0/0/0/0
2	SO4	D	923	-	-	0/0/0/0	0/0/0/0

All (46) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	550	HEM	C3B-C4B	-8.03	1.44	1.51
3	C	550	HEM	C3B-C4B	-7.92	1.44	1.51
3	B	550	HEM	C3B-C4B	-7.19	1.45	1.51
3	B	550	HEM	C2D-C3D	-7.10	1.33	1.54
3	C	550	HEM	C2D-C3D	-7.08	1.33	1.54
3	A	550	HEM	C3B-C4B	-6.86	1.45	1.51
3	D	550	HEM	C3B-CAB	-6.74	1.38	1.51
3	A	550	HEM	C2D-C3D	-6.70	1.34	1.54
3	B	550	HEM	C3B-CAB	-6.63	1.38	1.51
3	C	550	HEM	C3B-CAB	-6.55	1.39	1.51
3	D	550	HEM	C2D-C3D	-6.52	1.34	1.54
3	A	550	HEM	C3B-CAB	-6.26	1.39	1.51
3	A	550	HEM	C3C-CAC	-6.22	1.39	1.51
3	B	550	HEM	C3C-CAC	-6.20	1.39	1.51
3	C	550	HEM	C3C-CAC	-5.69	1.40	1.51
3	D	550	HEM	C3C-CAC	-5.61	1.40	1.51
3	D	550	HEM	C3D-C4D	-5.38	1.44	1.51
3	B	550	HEM	C3D-C4D	-4.90	1.45	1.51
3	A	550	HEM	C3D-C4D	-4.25	1.46	1.51
3	C	550	HEM	C3D-C4D	-4.21	1.46	1.51
3	D	550	HEM	C2C-C1C	-3.83	1.45	1.52
3	B	550	HEM	C2C-C1C	-3.80	1.45	1.52
3	C	550	HEM	C2C-C1C	-3.78	1.45	1.52
3	A	550	HEM	C2C-C1C	-3.40	1.46	1.52
3	C	550	HEM	C2B-C1B	-2.41	1.43	1.51
3	D	550	HEM	C2B-C1B	-2.40	1.44	1.51
3	A	550	HEM	C2B-C1B	-2.38	1.44	1.51
3	A	550	HEM	C2D-C1D	-2.31	1.44	1.51
3	B	550	HEM	C2D-C1D	-2.30	1.44	1.51
3	C	550	HEM	C2D-C1D	-2.29	1.44	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	550	HEM	C2B-C1B	-2.17	1.44	1.51
3	D	550	HEM	C2D-C1D	-2.16	1.44	1.51
4	D	603	H4B	C8A-N1	2.15	1.38	1.34
4	B	601	H4B	C8A-N1	2.23	1.38	1.34
4	B	601	H4B	C4A-N5	3.33	1.45	1.38
4	D	603	H4B	C6-N5	3.37	1.52	1.45
4	C	602	H4B	C4A-N5	3.43	1.46	1.38
4	A	600	H4B	C4A-N5	3.50	1.46	1.38
4	C	602	H4B	C4-N3	3.53	1.39	1.33
4	C	602	H4B	C6-N5	3.53	1.52	1.45
4	D	603	H4B	C4A-N5	3.54	1.46	1.38
4	A	600	H4B	C6-N5	3.66	1.53	1.45
4	A	600	H4B	C4-N3	3.67	1.39	1.33
4	B	601	H4B	C6-N5	3.81	1.53	1.45
4	B	601	H4B	C4-N3	3.90	1.40	1.33
4	D	603	H4B	C4-N3	4.05	1.40	1.33

All (62) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	603	H4B	N3-C2-N1	-4.50	118.16	125.53
4	C	602	H4B	N3-C2-N1	-4.42	118.29	125.53
4	A	600	H4B	N3-C2-N1	-4.31	118.47	125.53
4	B	601	H4B	N3-C2-N1	-4.28	118.51	125.53
3	C	550	HEM	CBD-CAD-C3D	-2.79	105.42	113.55
3	D	550	HEM	CBD-CAD-C3D	-2.66	105.80	113.55
3	A	550	HEM	CBD-CAD-C3D	-2.64	105.87	113.55
3	B	550	HEM	CBD-CAD-C3D	-2.44	106.47	113.55
3	B	550	HEM	C3B-CAB-CBB	-2.41	120.76	124.46
3	D	550	HEM	C3B-CAB-CBB	-2.29	120.95	124.46
3	B	550	HEM	CMD-C2D-C3D	2.44	125.14	114.35
3	D	550	HEM	C2D-C3D-C4D	2.62	105.94	101.50
3	A	550	HEM	CMD-C2D-C3D	2.64	126.04	114.35
3	D	550	HEM	CMD-C2D-C3D	2.77	126.60	114.35
3	B	550	HEM	C2D-C3D-C4D	2.82	106.28	101.50
3	C	550	HEM	CMD-C2D-C3D	2.84	126.89	114.35
3	A	550	HEM	C2D-C3D-C4D	2.90	106.41	101.50
3	C	550	HEM	C2D-C3D-C4D	3.11	106.77	101.50
3	C	550	HEM	C3C-CAC-CBC	3.21	129.38	124.46
4	B	601	H4B	N2-C2-N1	3.30	122.66	117.20
4	B	601	H4B	C2-N1-C8A	3.31	121.99	114.54
4	A	600	H4B	C2-N1-C8A	3.32	122.01	114.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	603	H4B	C2-N1-C8A	3.33	122.02	114.54
3	D	550	HEM	C3C-CAC-CBC	3.33	129.57	124.46
3	A	550	HEM	C3C-CAC-CBC	3.35	129.59	124.46
4	C	602	H4B	N2-C2-N1	3.42	122.86	117.20
4	A	600	H4B	N2-C2-N1	3.42	122.87	117.20
4	C	602	H4B	C2-N1-C8A	3.46	122.31	114.54
4	D	603	H4B	N2-C2-N1	3.51	123.00	117.20
3	B	550	HEM	C3C-CAC-CBC	3.61	130.00	124.46
3	D	550	HEM	CMB-C2B-C3B	3.62	125.57	116.53
3	A	550	HEM	CMB-C2B-C3B	3.69	125.74	116.53
3	C	550	HEM	CAD-C3D-C2D	3.79	124.12	113.22
3	B	550	HEM	CMB-C2B-C3B	3.80	126.01	116.53
4	D	603	H4B	C4A-C8A-N8	3.83	122.94	118.43
3	C	550	HEM	CMB-C2B-C3B	3.85	126.13	116.53
4	B	601	H4B	C4A-C8A-N8	4.01	123.15	118.43
4	A	600	H4B	C4A-C8A-N8	4.17	123.34	118.43
3	D	550	HEM	CAD-C3D-C4D	4.20	127.30	112.47
3	A	550	HEM	CAD-C3D-C2D	4.25	125.42	113.22
4	C	602	H4B	C4A-C8A-N8	4.25	123.44	118.43
3	A	550	HEM	CAD-C3D-C4D	4.27	127.54	112.47
3	B	550	HEM	CAD-C3D-C2D	4.29	125.55	113.22
3	B	550	HEM	CAD-C3D-C4D	4.29	127.61	112.47
3	D	550	HEM	CAD-C3D-C2D	4.45	126.02	113.22
5	B	801	ITU	C2-S-C3	4.50	108.13	101.53
3	C	550	HEM	CMC-C2C-C3C	4.55	127.88	116.53
3	C	550	HEM	CAD-C3D-C4D	4.55	128.51	112.47
3	B	550	HEM	CMC-C2C-C3C	4.64	128.11	116.53
3	A	550	HEM	CMC-C2C-C3C	4.64	128.11	116.53
5	D	803	ITU	C2-S-C3	4.74	108.48	101.53
3	D	550	HEM	CMC-C2C-C3C	4.81	128.55	116.53
5	C	802	ITU	C2-S-C3	4.91	108.74	101.53
5	A	800	ITU	C2-S-C3	5.23	109.21	101.53
4	B	601	H4B	C4-N3-C2	6.71	125.25	115.94
4	A	600	H4B	C4-N3-C2	6.82	125.40	115.94
4	D	603	H4B	C4-N3-C2	6.87	125.48	115.94
4	C	602	H4B	C4-N3-C2	6.93	125.56	115.94
4	C	602	H4B	C4-C4A-C8A	8.64	122.38	114.56
4	A	600	H4B	C4-C4A-C8A	8.74	122.47	114.56
4	D	603	H4B	C4-C4A-C8A	8.74	122.47	114.56
4	B	601	H4B	C4-C4A-C8A	8.88	122.60	114.56

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	550	HEM	1	0
2	A	930	SO4	1	0
3	B	550	HEM	2	0
3	C	550	HEM	1	0
3	D	550	HEM	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 ⓘ

EDS was not executed - this section will therefore be empty.

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.