



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:56 PM GMT

PDB ID : 1HZU  
Title : DOMAIN SWING UPON HIS TO ALA MUTATION IN NITRITE REDUCTASE OF PSEUDOMONAS AERUGINOSA  
Authors : Brown, K.; Cutruzzola, F.; Brunori, M.; Tegoni, M.; Cambillau, C.  
Deposited on : 2001-01-26  
Resolution : 2.70 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

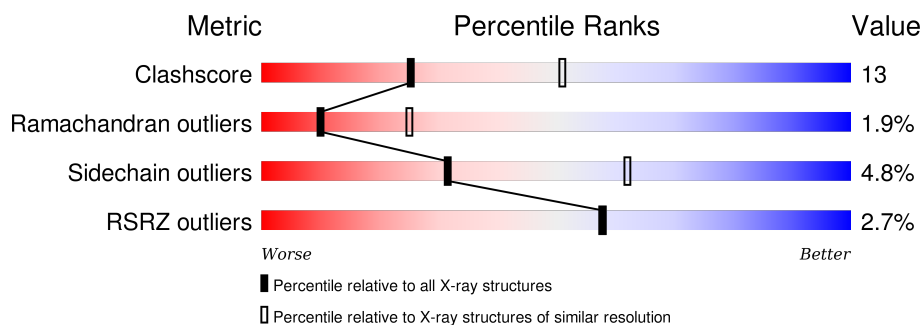
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.70 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	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (2.70-2.70)

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	543	

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	DHE	A	602	-	-	-	X

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4555 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 NITRITE REDUCTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	521	4081	2590	710	769	12	0	0	0

There is a discrepancy between the modelled and reference sequences:

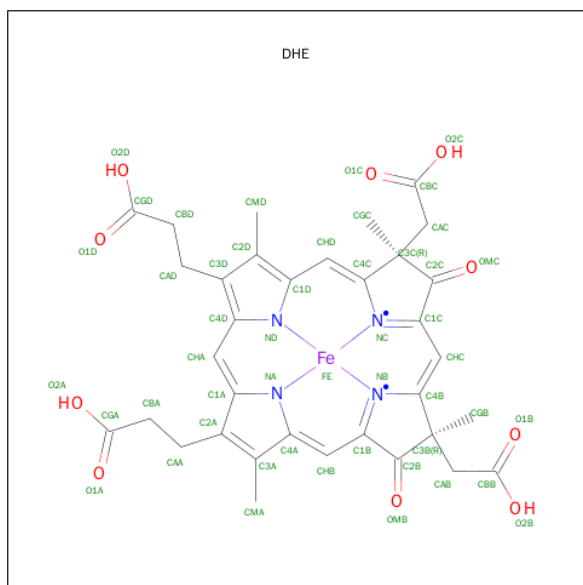
Chain	Residue	Modelled	Actual	Comment	Reference
A	327	ALA	HIS	ENGINEERED	UNP P24474

- Molecule 2 is HEME C (three-letter code: HEC) (formula:  $C_{34}H_{34}FeN_4O_4$ ).



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

- Molecule 3 is HEME D (three-letter code: DHE) (formula:  $C_{34}H_{32}FeN_4O_{10}$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	Fe	N	O	0	0
			49	34	1	4	10		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	382	Total O 382 382	0	0



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 43 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	70.52Å 70.52Å 281.23Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.86 – 2.70 26.48 – 2.70	Depositor EDS
% Data completeness (in resolution range)	94.8 (19.86-2.70) 94.7 (26.48-2.70)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	4.60	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.49 (at 2.72Å)	Xtriage
Refinement program	CNS 0.9	Depositor
R, $R_{free}$	0.210 , 0.248 0.202 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	71.7	Xtriage
Anisotropy	0.362	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 50.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 19418 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	4555	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.32% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: DHE, HEC

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	0.49	1/4182 (0.0%)	0.78	8/5679 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	33	GLU	CB-CG	-6.21	1.40	1.52

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	422	GLY	N-CA-C	7.55	131.99	113.10
1	A	33	GLU	N-CA-CB	-6.39	99.10	110.60
1	A	473	ASP	CA-C-N	5.97	130.34	117.20
1	A	473	ASP	C-N-CA	-5.96	106.81	121.70
1	A	499	ASN	N-CA-C	-5.83	95.25	111.00
1	A	26	ASN	N-CA-C	5.73	126.47	111.00
1	A	58	ALA	N-CA-C	-5.40	96.43	111.00
1	A	483	GLN	N-CA-C	5.23	125.13	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	472	ALA	Mainchain

## 5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4081	0	4051	99	0
2	A	43	0	30	5	0
3	A	49	0	24	7	0
4	A	382	0	0	28	1
All	All	4555	0	4105	104	1

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.

All (104) 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:71:ARG:HB3	1:A:75:TYR:HD1	1.34	0.93
1:A:25:THR:HB	1:A:111:HIS:HD2	1.34	0.90
1:A:25:THR:HB	1:A:111:HIS:CD2	2.09	0.88
1:A:474:LEU:HB3	4:A:811:HOH:O	1.80	0.82
1:A:28:ALA:HB1	1:A:29:PRO:HD2	1.62	0.79
1:A:79:LEU:HD13	2:A:601:HEC:HBA2	1.62	0.79
1:A:161:ILE:HB	4:A:767:HOH:O	1.81	0.79
1:A:155:LEU:HD12	1:A:160:GLN:HG3	1.64	0.79
1:A:417:GLN:HG2	4:A:799:HOH:O	1.81	0.78
1:A:71:ARG:HB3	1:A:75:TYR:CD1	2.20	0.76
3:A:602:DHE:HHA	4:A:711:HOH:O	1.89	0.71
1:A:229:SER:HB2	4:A:771:HOH:O	1.91	0.70
1:A:357:ARG:NH1	4:A:976:HOH:O	2.25	0.69
1:A:25:THR:HG21	1:A:111:HIS:HA	1.76	0.68
1:A:186:MET:SD	1:A:193:LEU:HD23	2.34	0.67
1:A:522:LYS:HA	4:A:817:HOH:O	1.93	0.67
1:A:58:ALA:O	1:A:60:GLY:N	2.28	0.67
1:A:423:SER:HB2	1:A:438:ASP:O	1.96	0.66
1:A:286:ILE:HG22	1:A:295:ILE:HB	1.77	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:164:VAL:HG22	1:A:171:ILE:HG12	1.78	0.65
1:A:426:ILE:HG22	1:A:437:VAL:HB	1.83	0.61
1:A:86:LEU:HB2	2:A:601:HEC:O1D	2.01	0.60
1:A:366:LYS:HB3	4:A:794:HOH:O	2.01	0.60
2:A:601:HEC:HMC1	2:A:601:HEC:HBC3	1.83	0.60
1:A:530:THR:HB	3:A:602:DHE:HAC1	1.84	0.59
1:A:258:LEU:HD23	4:A:865:HOH:O	2.04	0.58
1:A:261:LYS:HE2	4:A:779:HOH:O	2.02	0.58
1:A:53:VAL:HG21	1:A:246:TRP:CD1	2.38	0.58
1:A:476:GLU:HA	4:A:693:HOH:O	2.06	0.55
1:A:58:ALA:O	1:A:60:GLY:O	2.25	0.55
1:A:75:TYR:CZ	1:A:79:LEU:HD11	2.42	0.54
1:A:200:ALA:HB1	1:A:220:ILE:HG13	1.88	0.54
1:A:369:HIS:HB3	4:A:827:HOH:O	2.07	0.53
1:A:58:ALA:O	1:A:59:THR:C	2.46	0.53
1:A:369:HIS:HA	4:A:667:HOH:O	2.08	0.53
1:A:53:VAL:HG22	1:A:116:PRO:HB3	1.91	0.52
1:A:295:ILE:CD1	1:A:306:LEU:HD22	2.40	0.52
1:A:55:ARG:NH2	4:A:608:HOH:O	2.44	0.51
1:A:141:GLN:HG2	1:A:208:TRP:CD1	2.46	0.51
1:A:81:THR:HG22	1:A:102:ILE:HD13	1.92	0.51
1:A:312:ILE:HD11	4:A:779:HOH:O	2.11	0.51
1:A:293:GLU:HB3	1:A:308:ASN:HA	1.92	0.50
1:A:193:LEU:HB3	1:A:205:ILE:HB	1.92	0.50
3:A:602:DHE:CB	3:A:602:DHE:HMD1	2.40	0.50
1:A:201:ARG:HB2	4:A:880:HOH:O	2.11	0.50
1:A:306:LEU:HD12	1:A:356:ARG:HG2	1.92	0.50
1:A:126:GLU:HG2	4:A:674:HOH:O	2.12	0.50
1:A:524:PRO:HG2	4:A:770:HOH:O	2.12	0.50
1:A:82:TYR:O	1:A:90:ASN:HB3	2.11	0.50
1:A:182:HIS:HE1	3:A:602:DHE:C4A	2.25	0.49
1:A:90:ASN:HD22	1:A:93:SER:HB2	1.75	0.49
1:A:75:TYR:CE1	1:A:79:LEU:HD11	2.48	0.49
1:A:243:GLY:HA3	1:A:282:VAL:HG11	1.95	0.49
1:A:390:LEU:HD12	1:A:422:GLY:HA2	1.96	0.48
1:A:43:TYR:OH	1:A:55:ARG:HG2	2.14	0.48
1:A:141:GLN:NE2	1:A:143:ASN:H	2.12	0.48
1:A:180:ALA:HB1	1:A:198:ARG:HB2	1.97	0.47
1:A:201:ARG:NH1	4:A:906:HOH:O	2.48	0.47
1:A:372:ARG:NH2	3:A:602:DHE:CB	2.77	0.47
1:A:116:PRO:HB2	1:A:246:TRP:NE1	2.30	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:29:PRO:O	1:A:111:HIS:CE1	2.68	0.46
1:A:257:THR:O	1:A:258:LEU:HB2	2.16	0.46
1:A:220:ILE:HD13	1:A:242:ALA:HB3	1.97	0.46
1:A:480:ARG:HB2	1:A:498:TRP:HB3	1.98	0.46
1:A:160:GLN:HB2	1:A:175:ILE:O	2.16	0.45
1:A:334:SER:O	1:A:335:HIS:HB2	2.15	0.45
1:A:374:ALA:HB1	1:A:428:THR:HG22	1.97	0.45
1:A:96:GLU:HA	1:A:96:GLU:OE1	2.15	0.45
1:A:286:ILE:CG2	1:A:295:ILE:HB	2.44	0.45
1:A:389:HIS:HE1	1:A:396:SER:OG	1.99	0.45
1:A:291:HIS:HB3	1:A:292:PRO:HD2	1.98	0.45
1:A:84:THR:CG2	1:A:85:PRO:HD2	2.46	0.45
1:A:183:ILE:HD13	3:A:602:DHE:HBA1	1.98	0.45
1:A:295:ILE:HD12	1:A:306:LEU:HD22	1.98	0.44
1:A:43:TYR:CD1	1:A:47:CYS:HB2	2.52	0.44
1:A:386:SER:HB3	1:A:397:LEU:HD22	1.99	0.44
3:A:602:DHE:O1D	3:A:602:DHE:HMD1	2.18	0.44
1:A:500:GLY:HA2	4:A:903:HOH:O	2.17	0.44
1:A:444:ASP:OD2	1:A:446:ARG:HB3	2.18	0.43
1:A:511:ASP:HB3	1:A:514:THR:OG1	2.19	0.43
1:A:175:ILE:CD1	1:A:205:ILE:HD13	2.48	0.43
1:A:73:GLN:NE2	1:A:103:THR:HG23	2.33	0.43
1:A:81:THR:HG23	1:A:102:ILE:HG21	2.00	0.43
1:A:433:SER:HB3	4:A:809:HOH:O	2.19	0.43
1:A:128:TRP:CZ2	1:A:259:GLU:HB2	2.54	0.43
1:A:236:GLU:O	1:A:237:ASP:HB2	2.19	0.43
1:A:436:TYR:CE1	1:A:493:VAL:HG21	2.54	0.43
1:A:51:HIS:O	1:A:55:ARG:HA	2.18	0.43
1:A:170:LYS:HB2	4:A:941:HOH:O	2.19	0.43
1:A:267:ARG:NH1	4:A:782:HOH:O	2.50	0.42
2:A:601:HEC:HHA	2:A:601:HEC:HAA1	1.79	0.42
1:A:43:TYR:CZ	1:A:55:ARG:HG2	2.55	0.42
1:A:303:LYS:HD2	4:A:835:HOH:O	2.19	0.42
1:A:88:MET:HB2	2:A:601:HEC:C4D	2.50	0.41
1:A:146:ASP:HB3	4:A:854:HOH:O	2.20	0.41
1:A:239:TYR:CE1	1:A:312:ILE:HD13	2.56	0.41
1:A:185:ARG:NH2	1:A:372:ARG:HH12	2.19	0.41
1:A:32:SER:HB3	1:A:33:GLU:H	1.66	0.41
1:A:81:THR:HG22	1:A:102:ILE:CD1	2.51	0.41
1:A:526:LEU:HB3	4:A:816:HOH:O	2.20	0.41
1:A:388:SER:N	4:A:827:HOH:O	2.54	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:138:PRO:HB3	1:A:142:LEU:CD1	2.52	0.40
1:A:357:ARG:HD2	4:A:954:HOH:O	2.22	0.40
1:A:115:GLN:HA	1:A:116:PRO:HD2	1.93	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:783:HOH:O	4:A:984:HOH:O[5_556]	1.54	0.66

## 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	519/543 (96%)	471 (91%)	38 (7%)	10 (2%)	10	25

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	29	PRO
1	A	59	THR
1	A	473	ASP
1	A	499	ASN
1	A	501	LYS
1	A	421	GLY
1	A	26	ASN
1	A	483	GLN
1	A	27	GLY
1	A	365	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	442/457 (97%)	421 (95%)	21 (5%)	31 62

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	THR
1	A	29	PRO
1	A	33	GLU
1	A	34	SER
1	A	53	VAL
1	A	96	GLU
1	A	141	GLN
1	A	176	ASP
1	A	193	LEU
1	A	315	LEU
1	A	332	ASP
1	A	336	ARG
1	A	366	LYS
1	A	413	VAL
1	A	417	GLN
1	A	423	SER
1	A	437	VAL
1	A	439	THR
1	A	480	ARG
1	A	488	LYS
1	A	498	TRP

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

Mol	Chain	Res	Type
1	A	111	HIS
1	A	141	GLN
1	A	343	ASN
1	A	389	HIS

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Mol	Chain	Res	Type
1	A	539	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](#)

2 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	$\# Z  > 2$
2	HEC	A	601	1	24,50,50	1.24	2 (8%)	19,82,82	3.24	7 (36%)
3	DHE	A	602	1	29,56,56	2.20	8 (27%)	27,94,94	2.60	8 (29%)

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	HEC	A	601	1	-	0/6/54/54	0/0/8/8
3	DHE	A	602	1	-	0/12/108/108	0/0/8/8

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	602	DHE	CAC-C3C	-5.64	1.47	1.56
3	A	602	DHE	CAB-C3B	-4.91	1.48	1.56
3	A	602	DHE	CBD-CAD	-3.85	1.27	1.53
3	A	602	DHE	CAD-C3D	-3.38	1.46	1.52
2	A	601	HEC	C3B-C2B	-3.28	1.37	1.40
2	A	601	HEC	C3C-C2C	-3.09	1.37	1.40
3	A	602	DHE	FE-NB	2.53	2.05	1.95
3	A	602	DHE	CGC-C3C	3.31	1.61	1.54
3	A	602	DHE	CGB-C3B	3.35	1.61	1.54
3	A	602	DHE	FE-NC	3.41	2.15	1.97

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	HEC	CAA-C2A-C1A	-9.73	116.45	127.01
2	A	601	HEC	CBD-CAD-C3D	-4.22	104.96	112.53
2	A	601	HEC	CBB-CAB-C3B	-3.69	119.14	127.35
3	A	602	DHE	CMD-C2D-C1D	-2.69	123.91	128.36
2	A	601	HEC	CMB-C2B-C1B	-2.56	124.13	128.36
3	A	602	DHE	CMA-C3A-C4A	-2.52	124.19	128.36
2	A	601	HEC	CBC-CAC-C3C	-2.30	122.25	127.35
3	A	602	DHE	CAD-C3D-C4D	-2.06	124.77	127.01
2	A	601	HEC	CMA-C3A-C2A	2.27	129.99	125.24
3	A	602	DHE	CAA-CBA-CGA	2.60	117.52	112.75
3	A	602	DHE	CHB-C1B-NB	3.20	127.81	124.42
3	A	602	DHE	C1B-NB-C4B	4.57	109.38	106.90
2	A	601	HEC	CBA-CAA-C2A	6.30	123.82	112.53
3	A	602	DHE	CAD-CBD-CGD	6.87	125.34	112.75
3	A	602	DHE	CBD-CAD-C3D	7.47	125.92	112.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	HEC	5	0
3	A	602	DHE	7	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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	521/543 (95%)	-0.08	14 (2%) 58 58	41, 64, 104, 126	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	32	SER	3.4
1	A	94	SER	3.4
1	A	502	ASN	3.2
1	A	501	LYS	3.0
1	A	476	GLU	2.7
1	A	24	ARG	2.7
1	A	500	GLY	2.7
1	A	426	ILE	2.6
1	A	193	LEU	2.6
1	A	26	ASN	2.5
1	A	31	MET	2.4
1	A	265	SER	2.1
1	A	489	ARG	2.1
1	A	485	GLU	2.0

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

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

### 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.



## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 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
3	DHE	A	602	49/49	0.40	0.64	7.63	86,101,105,109	49
2	HEC	A	601	43/43	0.96	0.17	-0.56	54,62,72,75	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.