



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:43 PM GMT

PDB ID : 4MK4
Title : S197C variant of human ferrochelatase.
Authors : Lanzilotta, W.N.; Medlock, A.E.; Dailey, T.E.; Dailey, H.A.
Deposited on : 2013-09-04
Resolution : 2.50 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) : 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

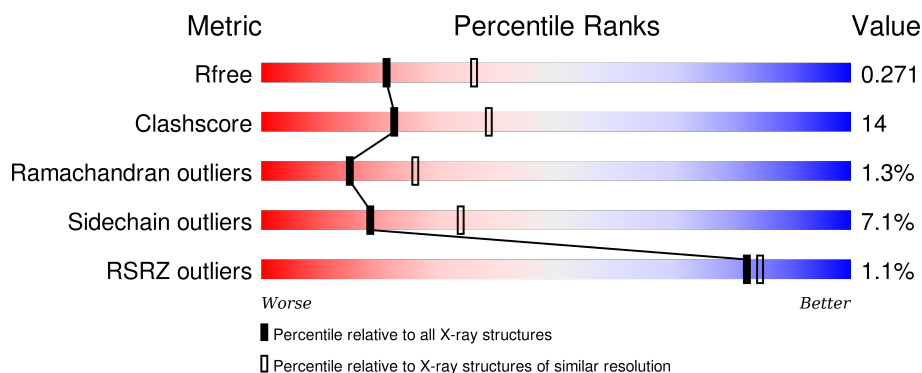
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	359	<div> <div></div> <div>72%</div> <div>23%</div> <div>.</div> </div>
1	B	359	<div> <div></div> <div>71%</div> <div>25%</div> <div>.</div> </div>

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	GOL	A	502	-	-	-	X
4	CHD	A	503	-	-	-	X
4	CHD	A	504	-	-	-	X
4	CHD	B	503	-	-	-	X

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6092 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.

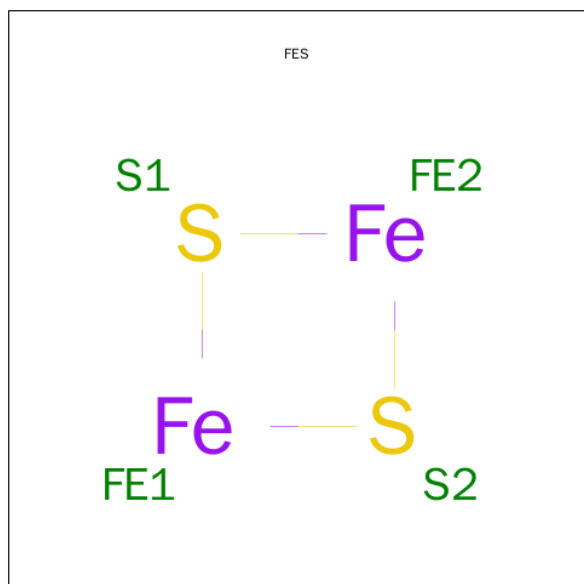
- Molecule 1 is a protein called Ferrochelatase, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	359	Total	C	N	O	S	0	0	0
			2891	1841	503	528	19			
1	B	359	Total	C	N	O	S	0	1	0
			2900	1846	505	530	19			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	197	CYS	SER	ENGINEERED MUTATION	UNP P22830
B	197	CYS	SER	ENGINEERED MUTATION	UNP P22830

- Molecule 2 is FE2/S2 (INORGANIC) CLUSTER (three-letter code: FES) (formula: Fe₂S₂).



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

Continued on next page...

Continued from previous page...

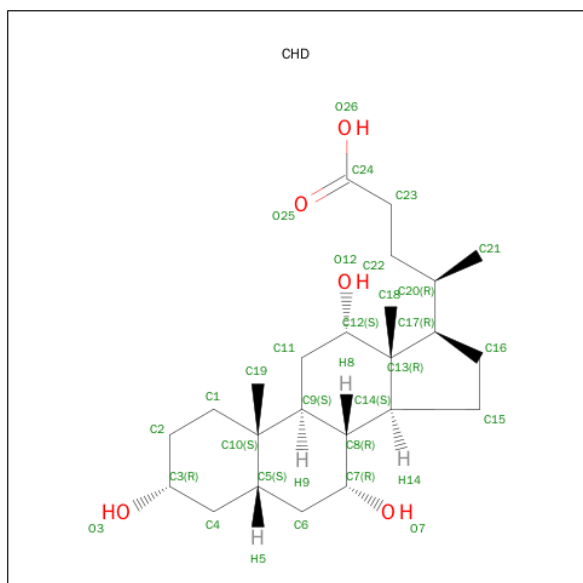
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	Fe	S		
2	B	1	4	2	2	0	0

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
3	A	1	6	3	3	0	0

- Molecule 4 is CHOLIC ACID (three-letter code: CHD) (formula: $C_{24}H_{40}O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			29	24	5		
4	A	1	Total	C	O	0	0
			29	24	5		
4	B	1	Total	C	O	0	0
			29	24	5		
4	B	1	Total	C	O	0	0
			29	24	5		
4	B	1	Total	C	O	0	0
			29	24	5		

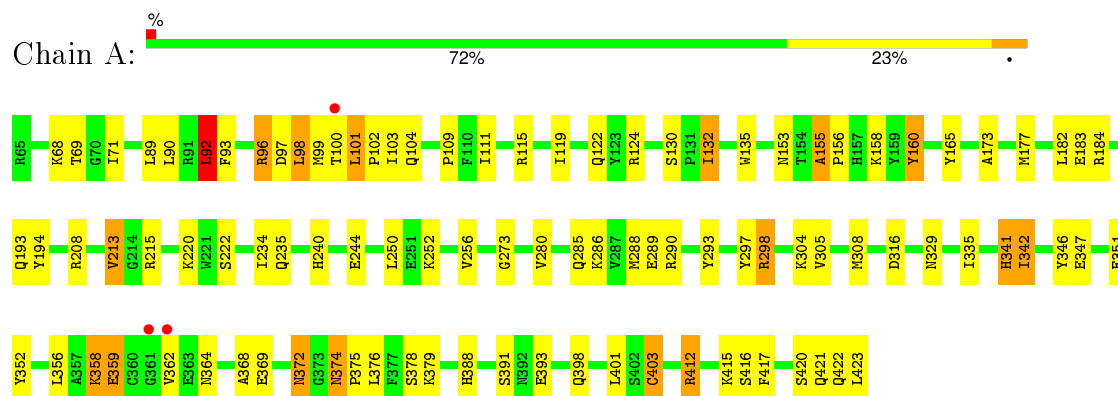
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	79	Total	O	0	0
			79	79		
5	B	63	Total	O	0	0
			63	63		

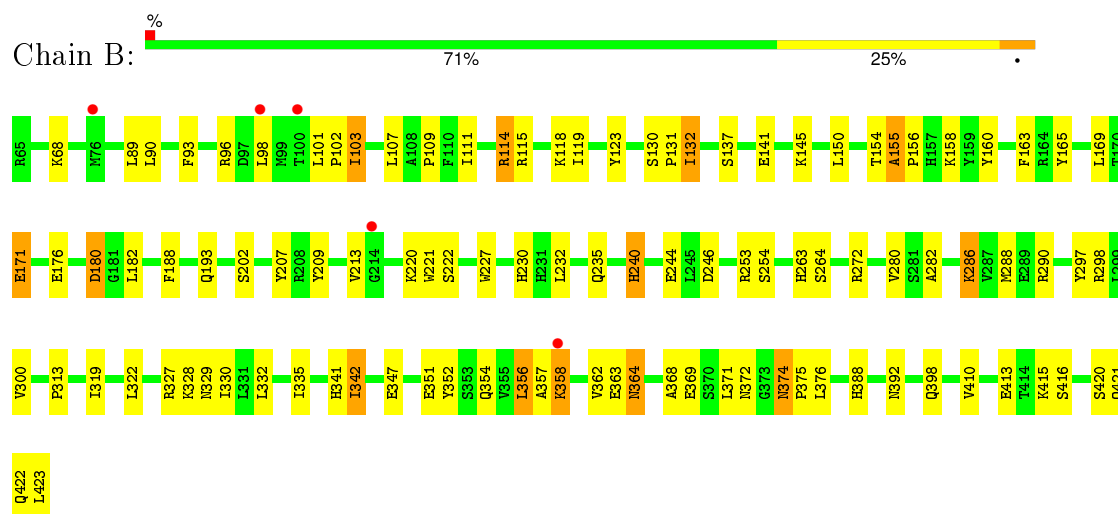
3 Residue-property plots [i](#)

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: Ferrochelatase, mitochondrial



• Molecule 1: Ferrochelatase, mitochondrial



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	88.17Å 93.43Å 111.74Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	60.00 – 2.50 64.13 – 2.50	Depositor EDS
% Data completeness (in resolution range)	94.2 (60.00-2.50) 94.2 (64.13-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.29 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.221 , 0.277 0.217 , 0.271	Depositor DCC
R_{free} test set	1563 reflections (5.38%)	DCC
Wilson B-factor (Å ²)	35.2	Xtriage
Anisotropy	0.522	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 36.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 30640 reflections (0.003%)	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6092	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 67.14 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 5.3960e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, CHD, FES

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	0.79	0/2961	0.84	1/4011 (0.0%)
1	B	0.75	0/2970	0.80	1/4023 (0.0%)
All	All	0.77	0/5931	0.82	2/8034 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	92	LEU	CA-CB-CG	9.17	136.40	115.30
1	B	356	LEU	CA-CB-CG	6.61	130.50	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2891	0	2899	78	0
1	B	2900	0	2906	79	0
2	A	4	0	0	0	0
2	B	4	0	0	1	0
3	A	6	0	8	0	0
4	A	58	0	78	6	0
4	B	87	0	117	13	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	79	0	0	4	0
5	B	63	0	0	4	0
All	All	6092	0	6008	164	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 14.

All (164) 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:304:LYS:HE2	5:A:649:HOH:O	1.44	1.13
1:B:300:VAL:HG12	1:B:313:PRO:HG2	1.37	1.05
1:A:155:ALA:HB1	1:A:156:PRO:HD3	1.32	1.03
1:B:235:GLN:HG3	1:B:290:ARG:NH1	1.79	0.98
1:B:155:ALA:HB1	1:B:156:PRO:CD	1.93	0.97
1:A:155:ALA:HB1	1:A:156:PRO:CD	1.96	0.96
1:B:155:ALA:HB1	1:B:156:PRO:HD3	1.49	0.92
1:A:374:ASN:ND2	1:A:376:LEU:H	1.67	0.92
1:B:240:HIS:HD2	1:B:369:GLU:O	1.55	0.88
1:B:244:GLU:HG3	1:B:368:ALA:HA	1.56	0.85
1:A:222:SER:OG	1:A:388:HIS:HE1	1.58	0.84
1:A:398:GLN:HE22	1:B:297:TYR:H	1.27	0.81
1:A:240:HIS:HD2	1:A:369:GLU:O	1.62	0.81
1:A:115:ARG:HG2	1:A:115:ARG:HH11	1.48	0.79
1:A:297:TYR:H	1:B:398:GLN:HE22	1.28	0.77
1:B:222:SER:OG	1:B:388:HIS:HE1	1.68	0.77
1:A:374:ASN:HD22	1:A:376:LEU:H	1.32	0.77
1:A:244:GLU:HG3	1:A:368:ALA:HA	1.67	0.76
1:A:155:ALA:CB	1:A:156:PRO:CD	2.64	0.76
1:A:288:MET:HG3	1:A:297:TYR:CD2	2.21	0.75
1:B:68:LYS:CB	1:B:155:ALA:HB3	2.18	0.73
1:B:68:LYS:HB3	1:B:155:ALA:HB3	1.69	0.72
1:A:222:SER:OG	1:A:388:HIS:CE1	2.44	0.71
1:A:374:ASN:HD22	1:A:374:ASN:C	1.94	0.70
1:B:155:ALA:CB	1:B:156:PRO:CD	2.69	0.69
1:B:123:TYR:CZ	1:B:342:ILE:HG13	2.27	0.69
1:B:319:ILE:HD11	1:B:332:LEU:HD21	1.76	0.68
1:B:240:HIS:CD2	1:B:369:GLU:O	2.44	0.67
1:B:244:GLU:HG2	5:B:655:HOH:O	1.95	0.66
1:B:328:LYS:O	1:B:363:GLU:O	2.13	0.66
4:A:504:CHD:H12	4:A:504:CHD:H212	1.79	0.65

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:114:ARG:HH11	1:B:114:ARG:HG2	1.60	0.64
1:B:68:LYS:HD3	1:B:155:ALA:CB	2.27	0.64
1:A:329:ASN:HD22	1:A:364:ASN:HB2	1.63	0.64
1:A:102:PRO:O	1:A:103:ILE:HG12	1.98	0.63
1:A:250:LEU:H	1:A:250:LEU:HD12	1.64	0.62
1:B:330:ILE:N	1:B:330:ILE:HD12	2.14	0.62
1:A:234:ILE:HG13	1:A:286:LYS:HG2	1.81	0.62
1:A:173:ALA:O	1:A:177:MET:HG3	2.00	0.62
1:B:222:SER:OG	1:B:388:HIS:CE1	2.51	0.62
1:A:89:LEU:HA	1:A:92:LEU:HD23	1.82	0.62
1:A:89:LEU:HA	1:A:92:LEU:CD2	2.29	0.62
1:A:358:LYS:O	1:A:359:GLU:CB	2.48	0.61
1:B:115:ARG:HD2	4:B:502:CHD:H221	1.83	0.61
1:B:235:GLN:HG3	1:B:290:ARG:HH11	1.63	0.60
1:A:69:THR:HG23	1:A:184:ARG:HD2	1.82	0.60
1:A:412:ARG:HB2	1:A:412:ARG:HH21	1.66	0.60
1:A:417:PHE:O	1:A:421:GLN:HG2	2.03	0.59
1:A:98:LEU:HD11	1:A:165:TYR:OH	2.03	0.58
1:B:68:LYS:HD3	1:B:155:ALA:HB2	1.84	0.58
1:A:374:ASN:HD22	1:A:376:LEU:N	2.00	0.58
1:B:171:GLU:H	1:B:171:GLU:CD	2.07	0.57
1:A:374:ASN:HD22	1:A:375:PRO:N	2.01	0.57
1:B:416:SER:O	1:B:420:SER:HB2	2.04	0.57
1:B:422:GLN:O	1:B:423:LEU:HB2	2.04	0.57
1:B:374:ASN:HD22	1:B:376:LEU:H	1.53	0.57
1:B:288:MET:HG3	1:B:297:TYR:CD2	2.40	0.56
1:A:93:PHE:CE1	1:A:111:ILE:HG22	2.40	0.56
1:A:374:ASN:ND2	1:A:376:LEU:N	2.46	0.56
1:B:111:ILE:HG23	4:B:502:CHD:H213	1.89	0.55
1:B:374:ASN:ND2	1:B:376:LEU:H	2.04	0.55
1:A:401:LEU:O	1:B:298:ARG:NH2	2.40	0.54
1:B:272:ARG:NH1	2:B:501:FES:S2	2.81	0.54
1:B:388:HIS:HD2	5:B:656:HOH:O	1.91	0.54
4:A:503:CHD:H222	5:A:613:HOH:O	2.07	0.54
1:B:354[B]:GLN:O	1:B:358:LYS:HE3	2.07	0.54
1:B:93:PHE:CE1	1:B:111:ILE:HG22	2.42	0.54
4:A:503:CHD:C22	5:A:613:HOH:O	2.54	0.54
1:A:115:ARG:HG2	1:A:115:ARG:NH1	2.17	0.53
1:A:97:ASP:OD1	1:A:208:ARG:NH2	2.41	0.53
1:A:288:MET:HG3	1:A:297:TYR:CE2	2.42	0.53
1:A:297:TYR:CD1	1:A:297:TYR:C	2.82	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:90:LEU:HD13	1:B:109:PRO:HA	1.89	0.53
1:B:89:LEU:HD21	1:B:119:ILE:HD13	1.89	0.52
1:A:194:TYR:HA	5:A:636:HOH:O	2.09	0.51
1:A:71:ILE:N	1:A:71:ILE:HD12	2.25	0.51
1:A:235:GLN:HG3	1:A:290:ARG:NH1	2.26	0.51
1:B:263:HIS:HD2	1:B:264:SER:O	1.93	0.50
1:A:101:LEU:HB2	1:A:104:GLN:HG3	1.92	0.50
1:B:114:ARG:HG2	1:B:114:ARG:NH1	2.26	0.50
1:A:391:SER:OG	1:A:393:GLU:HB2	2.11	0.50
1:B:288:MET:HG3	1:B:297:TYR:CE2	2.47	0.49
1:A:122:GLN:HG2	1:A:346:TYR:CE2	2.47	0.49
1:A:358:LYS:O	1:A:359:GLU:HB3	2.12	0.49
1:A:403:CYS:HB3	1:B:298:ARG:NH1	2.27	0.49
1:A:68:LYS:CB	1:A:155:ALA:HB3	2.42	0.49
1:B:68:LYS:HD3	1:B:155:ALA:HB3	1.94	0.49
4:A:504:CHD:H212	4:A:504:CHD:C12	2.42	0.49
1:A:135:TRP:CZ3	1:A:372:ASN:HB3	2.48	0.49
1:A:68:LYS:HB2	1:A:155:ALA:HB3	1.96	0.48
1:A:416:SER:O	1:A:420:SER:HB2	2.14	0.48
1:B:352:TYR:HA	1:B:356:LEU:HB3	1.94	0.48
1:A:90:LEU:HD13	1:A:109:PRO:HA	1.95	0.48
1:B:329:ASN:HD22	1:B:364:ASN:HB2	1.77	0.48
1:A:273:GLY:O	1:B:298:ARG:NH1	2.43	0.48
1:B:111:ILE:CG2	4:B:502:CHD:H213	2.44	0.47
1:B:163:PHE:O	1:B:169:LEU:HA	2.15	0.47
1:A:96:ARG:HH12	1:A:104:GLN:HB3	1.78	0.47
1:A:285:GLN:HG2	1:B:282:ALA:HB1	1.96	0.47
4:B:502:CHD:H193	4:B:504:CHD:H162	1.96	0.47
1:A:374:ASN:C	1:A:374:ASN:ND2	2.67	0.47
1:A:316:ASP:HB3	1:A:352:TYR:CE1	2.50	0.46
1:A:215:ARG:HB2	1:A:215:ARG:HE	1.54	0.46
1:B:98:LEU:HD11	1:B:165:TYR:OH	2.16	0.46
1:B:68:LYS:HB2	1:B:155:ALA:HB3	1.97	0.46
1:B:356:LEU:O	1:B:357:ALA:HB3	2.16	0.46
1:B:193:GLN:HG2	1:B:280:VAL:HA	1.97	0.46
1:B:130:SER:OG	1:B:132:ILE:HG22	2.16	0.45
1:A:89:LEU:HD21	1:A:119:ILE:CD1	2.46	0.45
1:B:347:GLU:O	1:B:351:GLU:HB2	2.16	0.45
1:B:354[A]:GLN:O	1:B:358:LYS:HE3	2.16	0.45
4:B:504:CHD:H222	4:B:504:CHD:H162	1.58	0.45
1:A:193:GLN:HG2	1:A:280:VAL:HA	1.98	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:412:ARG:CB	1:A:412:ARG:HH21	2.30	0.45
1:A:342:ILE:HD13	1:A:346:TYR:HB3	1.98	0.45
1:A:289:GLU:HG3	1:A:293:TYR:OH	2.17	0.45
4:B:502:CHD:H232	4:B:502:CHD:H211	1.73	0.44
1:B:374:ASN:HD22	1:B:375:PRO:N	2.15	0.44
1:B:221:TRP:HB2	1:B:421:GLN:NE2	2.33	0.44
1:B:137:SER:O	1:B:141:GLU:HG3	2.18	0.44
1:B:371:LEU:O	1:B:372:ASN:C	2.56	0.44
1:A:256:VAL:HA	1:A:329:ASN:O	2.18	0.44
1:A:347:GLU:O	1:A:351:GLU:HB2	2.17	0.44
1:A:240:HIS:CD2	1:A:369:GLU:O	2.54	0.44
1:A:97:ASP:CG	1:A:208:ARG:HH22	2.21	0.44
1:B:322:LEU:HB3	1:B:327:ARG:HB2	2.00	0.44
1:A:335:ILE:O	1:A:335:ILE:HG13	2.18	0.44
4:A:503:CHD:H12	4:A:503:CHD:H212	2.00	0.43
1:B:102:PRO:C	1:B:103:ILE:HG13	2.37	0.43
1:A:252:LYS:HD3	1:A:329:ASN:OD1	2.19	0.43
1:B:68:LYS:HE2	1:B:154:THR:HA	2.00	0.43
1:B:176:GLU:O	1:B:180:ASP:OD1	2.36	0.43
1:A:297:TYR:HD1	1:A:298:ARG:N	2.17	0.42
1:B:374:ASN:HD22	1:B:374:ASN:C	2.23	0.42
1:B:111:ILE:HG23	4:B:502:CHD:C21	2.50	0.42
1:A:130:SER:OG	1:A:132:ILE:HG22	2.19	0.42
4:B:503:CHD:H162	4:B:503:CHD:H222	1.72	0.42
1:B:253:ARG:NH2	5:B:660:HOH:O	2.44	0.42
1:A:160:TYR:CD1	1:A:160:TYR:N	2.87	0.42
1:A:341:HIS:N	1:A:341:HIS:CD2	2.86	0.42
1:A:96:ARG:HA	1:A:99:MET:O	2.19	0.42
1:B:230:HIS:CE1	1:B:232:LEU:HB2	2.55	0.42
1:B:286:LYS:HD3	5:B:659:HOH:O	2.20	0.42
1:A:422:GLN:HG2	1:A:423:LEU:HG	2.02	0.42
1:B:398:GLN:NE2	1:B:398:GLN:HA	2.35	0.42
1:B:114:ARG:NH1	4:B:502:CHD:O26	2.53	0.41
4:B:502:CHD:H162	4:B:502:CHD:H222	1.91	0.41
1:B:107:LEU:HA	1:B:107:LEU:HD23	1.85	0.41
1:A:352:TYR:HA	1:A:356:LEU:HB3	2.02	0.41
1:B:131:PRO:O	1:B:132:ILE:C	2.59	0.41
1:A:305:VAL:HG23	4:A:504:CHD:H7	2.02	0.41
4:B:502:CHD:H20	4:B:502:CHD:H183	1.83	0.41
1:A:158:LYS:HD3	1:A:160:TYR:OH	2.21	0.41
1:B:158:LYS:HD3	1:B:160:TYR:CZ	2.55	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:304:LYS:NZ	1:A:308:MET:O	2.54	0.40
1:A:101:LEU:HB2	1:A:104:GLN:CG	2.51	0.40
1:B:171:GLU:CD	1:B:171:GLU:N	2.74	0.40
1:B:160:TYR:OH	1:B:182:LEU:HD11	2.22	0.40
1:B:207:TYR:CZ	1:B:413:GLU:HB3	2.57	0.40
1:B:335:ILE:O	1:B:335:ILE:HG13	2.21	0.40
4:B:502:CHD:H183	4:B:504:CHD:H151	2.04	0.40
4:B:504:CHD:H9	4:B:504:CHD:H21	1.81	0.40
1:A:160:TYR:OH	1:A:182:LEU:HD11	2.21	0.40
1:B:209:TYR:O	1:B:213:VAL:HG22	2.21	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	357/359 (99%)	331 (93%)	21 (6%)	5 (1%)	14	24
1	B	358/359 (100%)	328 (92%)	26 (7%)	4 (1%)	17	31
All	All	715/718 (100%)	659 (92%)	47 (7%)	9 (1%)	15	26

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	362	VAL
1	A	155	ALA
1	A	359	GLU
1	B	155	ALA
1	A	213	VAL
1	A	372	ASN
1	B	227	TRP
1	B	364	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	362	VAL

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	324/324 (100%)	302 (93%)	22 (7%)	20	36
1	B	325/324 (100%)	301 (93%)	24 (7%)	17	31
All	All	649/648 (100%)	603 (93%)	46 (7%)	18	34

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	92	LEU
1	A	96	ARG
1	A	98	LEU
1	A	100	THR
1	A	101	LEU
1	A	124	ARG
1	A	132	ILE
1	A	153	ASN
1	A	160	TYR
1	A	183	GLU
1	A	213	VAL
1	A	220	LYS
1	A	298	ARG
1	A	341	HIS
1	A	342	ILE
1	A	358	LYS
1	A	374	ASN
1	A	378	SER
1	A	379	LYS
1	A	403	CYS
1	A	412	ARG
1	A	415	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	96	ARG
1	B	101	LEU
1	B	103	ILE
1	B	114	ARG
1	B	118	LYS
1	B	132	ILE
1	B	145	LYS
1	B	150	LEU
1	B	171	GLU
1	B	180	ASP
1	B	188	PHE
1	B	202	SER
1	B	220	LYS
1	B	240	HIS
1	B	246	ASP
1	B	254	SER
1	B	286	LYS
1	B	341	HIS
1	B	342	ILE
1	B	358	LYS
1	B	374	ASN
1	B	392	ASN
1	B	410	VAL
1	B	415	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	211	ASN
1	A	235	GLN
1	A	240	HIS
1	A	263	HIS
1	A	302	GLN
1	A	329	ASN
1	A	354	GLN
1	A	364	ASN
1	A	374	ASN
1	A	388	HIS
1	A	398	GLN
1	A	421	GLN
1	B	212	GLN
1	B	235	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	240	HIS
1	B	263	HIS
1	B	302	GLN
1	B	329	ASN
1	B	364	ASN
1	B	374	ASN
1	B	388	HIS
1	B	398	GLN
1	B	421	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](#)

8 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	FES	A	501	1	0,4,4	0.00	-	0,4,4	0.00	-
3	GOL	A	502	-	5,5,5	0.30	0	5,5,5	0.30	0
4	CHD	A	503	-	29,32,32	0.67	1 (3%)	48,51,51	1.26	4 (8%)
4	CHD	A	504	-	29,32,32	0.72	0	48,51,51	1.66	11 (22%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	FES	B	501	1	0,4,4	0.00	-	0,4,4	0.00	-
4	CHD	B	502	-	29,32,32	0.65	1 (3%)	48,51,51	2.18	16 (33%)
4	CHD	B	503	-	29,32,32	0.75	0	48,51,51	2.12	16 (33%)
4	CHD	B	504	-	29,32,32	0.53	0	48,51,51	1.57	8 (16%)

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	FES	A	501	1	-	0/0/4/4	0/1/1/1
3	GOL	A	502	-	-	0/4/4/4	0/0/0/0
4	CHD	A	503	-	-	0/7/74/74	0/4/4/4
4	CHD	A	504	-	-	0/7/74/74	0/4/4/4
2	FES	B	501	1	-	0/0/4/4	0/1/1/1
4	CHD	B	502	-	-	0/7/74/74	0/4/4/4
4	CHD	B	503	-	-	0/7/74/74	0/4/4/4
4	CHD	B	504	-	-	0/7/74/74	0/4/4/4

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	502	CHD	C13-C14	-2.13	1.51	1.55
4	A	503	CHD	C4-C3	2.10	1.55	1.51

All (55) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	503	CHD	C23-C22-C20	-5.07	108.78	114.75
4	B	503	CHD	C17-C13-C14	-4.70	95.30	100.05
4	B	502	CHD	C9-C8-C7	-4.59	106.50	111.92
4	B	503	CHD	C13-C17-C20	-4.53	113.98	119.50
4	B	502	CHD	C19-C10-C5	-4.15	102.92	110.25
4	B	504	CHD	C6-C5-C4	-4.09	106.48	111.05
4	A	504	CHD	C17-C13-C12	-3.85	114.27	117.68
4	B	502	CHD	C13-C17-C20	-3.66	115.04	119.50
4	B	502	CHD	C14-C13-C12	-3.45	104.30	107.39
4	A	504	CHD	C13-C17-C20	-3.38	115.39	119.50
4	A	504	CHD	C9-C8-C7	-3.32	108.00	111.92
4	B	504	CHD	C1-C10-C9	-3.20	106.29	111.45

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	504	CHD	C14-C8-C9	-2.55	106.11	109.62
4	B	503	CHD	C14-C8-C7	-2.45	108.35	111.74
4	B	503	CHD	C6-C5-C4	-2.45	108.31	111.05
4	B	503	CHD	C21-C20-C17	-2.34	109.06	112.96
4	B	502	CHD	C13-C14-C8	-2.19	111.93	114.75
4	B	503	CHD	C11-C12-C13	-2.13	109.03	111.20
4	A	504	CHD	C21-C20-C17	-2.12	109.42	112.96
4	B	503	CHD	C9-C11-C12	-2.03	111.80	114.36
4	A	503	CHD	C19-C10-C9	-2.01	108.17	111.18
4	A	503	CHD	C1-C10-C9	2.01	114.70	111.45
4	A	504	CHD	C4-C5-C10	2.03	114.89	112.66
4	B	503	CHD	C6-C5-C10	2.06	114.93	112.66
4	A	504	CHD	C5-C6-C7	2.27	116.96	114.44
4	B	502	CHD	C22-C20-C17	2.27	115.00	110.24
4	B	502	CHD	C11-C9-C8	2.28	113.97	110.73
4	B	502	CHD	C4-C5-C10	2.29	115.18	112.66
4	B	503	CHD	C11-C9-C8	2.35	114.07	110.73
4	B	504	CHD	C5-C4-C3	2.44	116.54	112.91
4	B	503	CHD	C15-C14-C13	2.46	106.05	103.60
4	B	502	CHD	C1-C10-C5	2.48	111.88	107.81
4	A	504	CHD	C18-C13-C12	2.49	111.52	109.09
4	A	504	CHD	C22-C20-C17	2.58	115.66	110.24
4	B	504	CHD	C17-C13-C14	2.58	102.66	100.05
4	A	503	CHD	C2-C1-C10	2.58	117.45	112.84
4	B	503	CHD	C5-C6-C7	2.61	117.35	114.44
4	B	503	CHD	O12-C12-C11	2.64	114.46	109.06
4	B	502	CHD	C18-C13-C12	2.64	111.66	109.09
4	B	504	CHD	C9-C10-C5	2.80	112.82	108.67
4	A	504	CHD	C6-C7-C8	2.83	114.47	111.47
4	B	504	CHD	C4-C5-C10	2.90	115.85	112.66
4	A	504	CHD	C15-C14-C8	2.93	122.57	118.32
4	B	502	CHD	C1-C2-C3	3.01	115.31	110.43
4	B	502	CHD	C9-C11-C12	3.01	118.16	114.36
4	A	503	CHD	C6-C7-C8	3.21	114.87	111.47
4	B	504	CHD	C10-C9-C8	3.23	115.43	111.88
4	B	502	CHD	C11-C12-C13	3.29	114.55	111.20
4	B	502	CHD	C9-C10-C5	3.48	113.83	108.67
4	B	503	CHD	C16-C17-C20	3.59	118.45	112.05
4	A	504	CHD	C16-C17-C20	3.78	118.80	112.05
4	B	503	CHD	C16-C17-C13	3.94	107.52	103.60
4	B	503	CHD	C14-C13-C12	4.49	111.41	107.39
4	B	502	CHD	C17-C13-C14	4.64	104.75	100.05

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	502	CHD	C6-C5-C10	5.18	118.37	112.66

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	503	CHD	3	0
4	A	504	CHD	3	0
2	B	501	FES	1	0
4	B	502	CHD	10	0
4	B	503	CHD	1	0
4	B	504	CHD	4	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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	359/359 (100%)	-0.14	3 (0%) 87 89	18, 34, 55, 66	0
1	B	359/359 (100%)	-0.12	5 (1%) 78 80	16, 33, 59, 69	0
All	All	718/718 (100%)	-0.13	8 (1%) 82 84	16, 33, 56, 69	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	362	VAL	4.8
1	B	98	LEU	2.9
1	B	100	THR	2.9
1	B	76	MET	2.3
1	A	361	GLY	2.1
1	B	358	LYS	2.1
1	A	100	THR	2.1
1	B	214	GLY	2.1

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains.

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	GOL	A	502	6/6	0.76	0.35	15.16	45,45,46,47	0
4	CHD	A	504	29/29	0.81	0.30	3.50	77,79,89,89	0
4	CHD	B	503	29/29	0.79	0.33	2.74	61,64,80,82	0
4	CHD	A	503	29/29	0.85	0.23	2.07	53,57,62,63	0
4	CHD	B	502	29/29	0.77	0.29	1.57	100,104,112,112	0
4	CHD	B	504	29/29	0.83	0.24	0.98	72,79,95,95	0
2	FES	A	501	4/4	0.99	0.06	-1.55	36,39,41,43	0
2	FES	B	501	4/4	0.99	0.06	-1.96	43,45,45,46	0

6.5 Other polymers [i](#)

There are no such residues in this entry.