Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 07:52 AM GMT

PDB ID	:	3CHI
Title	:	Crystal Structure of Di-iron AurF (Monoclinic form)
Authors	:	Zhang, H.; Brunzelle, J.S.; Nair, S.K.
Deposited on	:	2008-03-09
Resolution	:	2.10 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 (i) symbol.

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

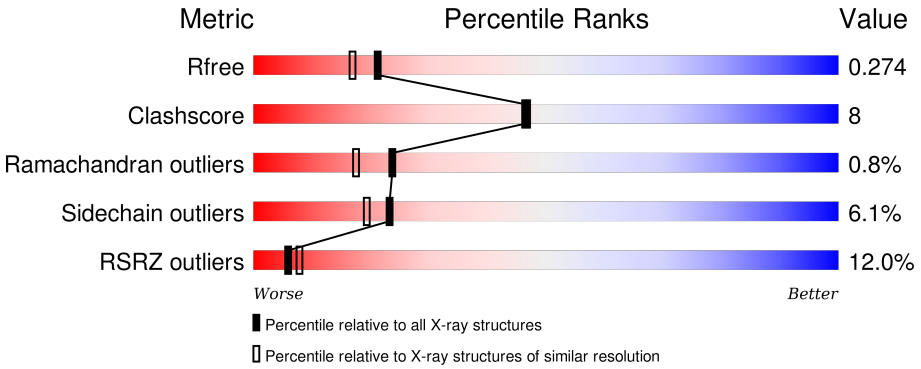
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 2.10 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	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (2.10-2.10)

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	336	8% 73% 14% • 10%
1	B	336	13% 71% 17% • 10%

2 Entry composition

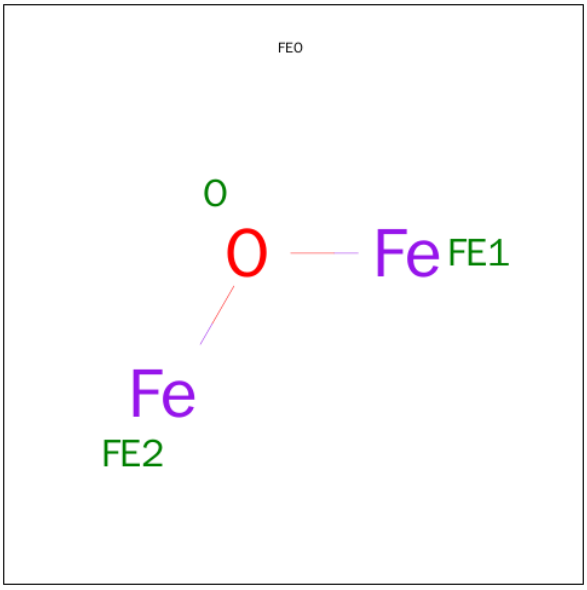
There are 3 unique types of molecules in this entry. The entry contains 5019 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 p-Aminobenzoate N-Oxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	301	Total	C	N	O	S	0	0	0
			2404	1525	431	438	10			
1	B	301	Total	C	N	O	S	0	0	0
			2404	1525	431	438	10			

- Molecule 2 is MU-OXO-DIIRON (three-letter code: FEO) (formula: Fe_2O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	Fe	O		0	0
			3	2	1			
2	B	1	Total	Fe	O		0	0
			3	2	1			

- Molecule 3 is water.

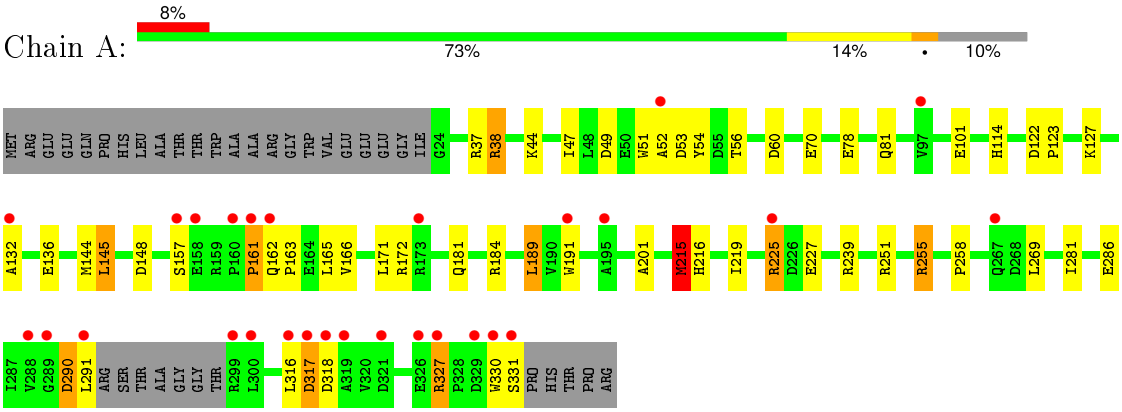
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	125	Total O 125 125	0	0
3	B	80	Total O 80 80	0	0

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.

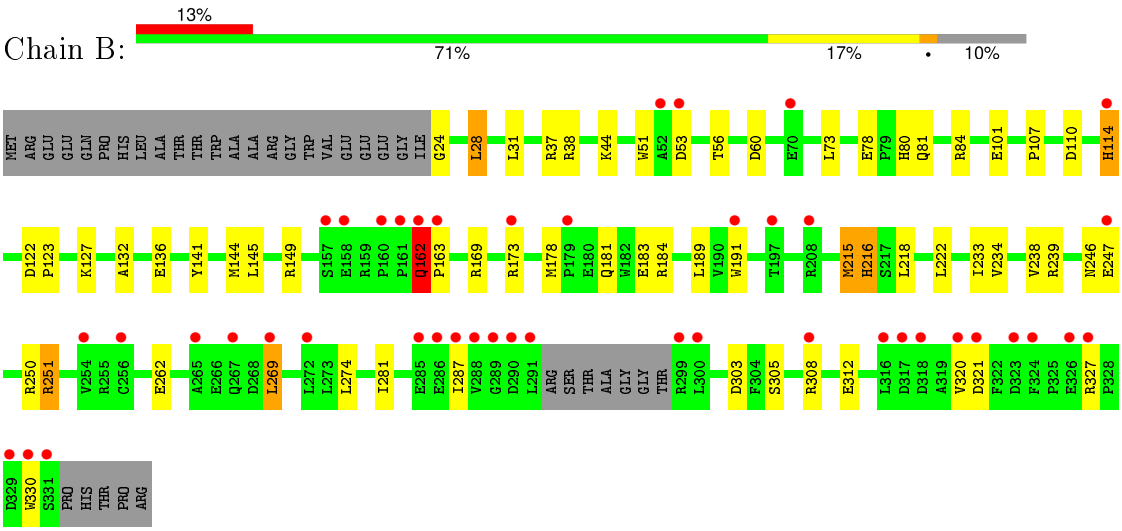
- Molecule 1: p-Aminobenzoate N-Oxygenase

Chain A:



- Molecule 1: p-Aminobenzoate N-Oxygenase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	58.42Å 73.32Å 74.69Å 90.00° 100.92° 90.00°	Depositor
Resolution (Å)	25.00 – 2.10 32.80 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.6 (25.00-2.10) 99.4 (32.80-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$< I/\sigma(I) >$ [1]	3.66 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.220 , 0.277 0.219 , 0.274	Depositor DCC
R_{free} test set	1800 reflections (5.24%)	DCC
Wilson B-factor (Å^2)	29.1	Xtriage
Anisotropy	0.235	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 52.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Outliers	2 of 36189 reflections (0.006%)	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	5019	wwPDB-VP
Average B, all atoms (Å^2)	36.0	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 26.88 % 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 2.4318e-03.The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[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: FEO

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	0.65	0/2454	0.75	6/3338 (0.2%)
1	B	0.64	3/2454 (0.1%)	0.74	5/3338 (0.1%)
All	All	0.65	3/4908 (0.1%)	0.74	11/6676 (0.2%)

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	B	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	251	ARG	CZ-NH2	6.81	1.42	1.33
1	B	251	ARG	NE-CZ	6.07	1.41	1.33
1	B	321	ASP	CG-OD2	5.02	1.36	1.25

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	38	ARG	NE-CZ-NH2	-9.51	115.55	120.30
1	B	251	ARG	NE-CZ-NH1	-7.96	116.32	120.30
1	B	251	ARG	NE-CZ-NH2	7.58	124.09	120.30
1	B	38	ARG	NE-CZ-NH2	-6.25	117.18	120.30
1	A	239	ARG	NE-CZ-NH1	-6.05	117.28	120.30
1	B	321	ASP	CB-CG-OD2	-5.95	112.95	118.30
1	A	189	LEU	CA-CB-CG	5.63	128.25	115.30
1	B	239	ARG	NE-CZ-NH1	-5.40	117.60	120.30

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	145	LEU	CA-CB-CG	5.39	127.71	115.30
1	A	165	LEU	CA-CB-CG	5.23	127.32	115.30
1	A	215	MET	CG-SD-CE	-5.05	92.12	100.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	162	GLN	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	2404	0	2403	33	1
1	B	2404	0	2403	42	0
2	A	3	0	0	0	0
2	B	3	0	0	1	0
3	A	125	0	0	8	0
3	B	80	0	0	7	0
All	All	5019	0	4806	73	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 8.

All (73) 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:225:ARG:HD3	3:A:719:HOH:O	1.70	0.89
1:B:78:GLU:H	1:B:81:GLN:HE21	1.20	0.85
1:A:148:ASP:OD2	3:A:721:HOH:O	1.95	0.83
1:A:38:ARG:O	1:A:225:ARG:NH2	2.12	0.83
1:B:181:GLN:NE2	1:B:184:ARG:HH11	1.76	0.82
1:B:169:ARG:O	1:B:173:ARG:HG3	1.82	0.80
1:A:49:ASP:HB3	3:A:708:HOH:O	1.80	0.80

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:101:GLU:OE2	1:A:136:GLU:OE1	2.04	0.74
1:A:78:GLU:H	1:A:81:GLN:HE21	1.32	0.74
1:B:78:GLU:H	1:B:81:GLN:NE2	1.85	0.73
1:B:101:GLU:OE2	2:B:501:FEO:FE2	1.41	0.73
1:A:144:MET:SD	3:B:672:HOH:O	2.48	0.72
1:A:37:ARG:HG2	3:A:716:HOH:O	1.92	0.69
1:B:262:GLU:HB3	3:B:656:HOH:O	1.93	0.68
1:B:162:GLN:HB3	1:B:163:PRO:HD3	1.74	0.67
1:A:225:ARG:CD	3:A:719:HOH:O	2.32	0.64
1:A:37:ARG:NE	3:A:716:HOH:O	2.31	0.64
1:B:162:GLN:HB3	1:B:163:PRO:CD	2.27	0.62
1:A:145:LEU:HD22	1:B:127:LYS:HE3	1.80	0.62
1:B:73:LEU:HB3	3:B:614:HOH:O	2.01	0.61
1:B:81:GLN:HG2	1:B:84:ARG:HH11	1.66	0.61
1:B:31:LEU:HD13	1:B:233:ILE:HG13	1.83	0.61
1:A:78:GLU:H	1:A:81:GLN:NE2	2.00	0.59
1:A:181:GLN:NE2	1:A:184:ARG:HH11	2.01	0.58
1:B:132:ALA:O	1:B:136:GLU:HG2	2.03	0.57
1:B:101:GLU:OE2	1:B:136:GLU:OE1	2.22	0.56
1:B:107:PRO:O	3:B:680:HOH:O	2.18	0.56
1:B:24:GLY:N	3:B:671:HOH:O	2.39	0.55
1:B:216:HIS:HE1	3:B:609:HOH:O	1.89	0.54
1:A:127:LYS:HE3	1:B:145:LEU:HD22	1.91	0.53
1:B:181:GLN:HE21	1:B:184:ARG:HH11	1.53	0.53
1:B:162:GLN:CB	1:B:163:PRO:HD3	2.38	0.53
1:B:269:LEU:H	1:B:269:LEU:CD2	2.22	0.53
1:A:51:TRP:NE1	1:A:215:MET:HE3	2.23	0.52
1:A:38:ARG:HA	1:A:225:ARG:HH22	1.75	0.52
1:A:225:ARG:CZ	1:A:225:ARG:HB3	2.43	0.48
1:B:178:MET:SD	1:B:183:GLU:HG2	2.53	0.48
1:A:286:GLU:O	1:A:290:ASP:HB3	2.13	0.48
1:B:308:ARG:O	1:B:312:GLU:HG3	2.13	0.47
1:A:227:GLU:OE2	1:A:227:GLU:HA	2.14	0.47
3:A:622:HOH:O	1:B:144:MET:HG3	2.14	0.47
1:A:54:TYR:HD2	1:A:215:MET:HE1	1.79	0.46
1:B:80:HIS:O	1:B:84:ARG:HG3	2.16	0.46
1:A:290:ASP:OD2	1:A:290:ASP:C	2.53	0.46
1:B:51:TRP:CD1	1:B:215:MET:HE3	2.50	0.46
1:B:246:ASN:O	1:B:250:ARG:HG2	2.16	0.46
1:B:234:VAL:O	1:B:238:VAL:HG23	2.14	0.46
1:B:269:LEU:H	1:B:269:LEU:HD23	1.81	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:44:LYS:HE3	1:B:218:LEU:HD21	1.98	0.45
1:B:303:ASP:HB2	1:B:330:TRP:CH2	2.51	0.45
1:B:51:TRP:CE2	1:B:215:MET:HG3	2.51	0.45
1:B:305:SER:OG	1:B:327:ARG:NE	2.49	0.45
1:A:316:LEU:C	1:A:318:ASP:H	2.20	0.44
1:A:78:GLU:N	1:A:81:GLN:HE21	2.09	0.44
1:B:149:ARG:HD2	1:B:149:ARG:HA	1.83	0.44
1:B:141:TYR:CZ	1:B:145:LEU:HD11	2.52	0.43
1:A:54:TYR:HE1	1:A:56:THR:HG22	1.83	0.43
1:A:255:ARG:O	1:A:258:PRO:HD2	2.18	0.43
1:B:122:ASP:HA	1:B:123:PRO:HD3	1.87	0.43
1:B:287:ILE:HD11	3:B:614:HOH:O	2.16	0.43
1:A:201:ALA:HA	1:A:330:TRP:CZ2	2.54	0.43
1:B:44:LYS:HG2	1:B:222:LEU:HD11	2.00	0.43
1:B:110:ASP:O	1:B:114:HIS:ND1	2.51	0.42
1:A:161:PRO:O	1:A:163:PRO:CD	2.67	0.42
1:B:247:GLU:HB3	1:B:251:ARG:NH2	2.34	0.42
1:A:60:ASP:O	1:A:216:HIS:HD2	2.03	0.42
1:A:132:ALA:O	1:A:136:GLU:HG2	2.20	0.41
1:B:60:ASP:O	1:B:216:HIS:HD2	2.03	0.41
1:A:166:VAL:HG12	3:A:678:HOH:O	2.20	0.41
1:B:28:LEU:HD12	1:B:28:LEU:HA	1.97	0.41
1:A:47:ILE:HD11	1:A:219:ILE:HG13	2.04	0.40
1:A:327:ARG:HD3	1:A:331:SER:HB3	2.03	0.40
1:A:122:ASP:HA	1:A:123:PRO:HD3	1.94	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)
1:A:225:ARG:NH1	1:A:251:ARG:NH2[2_646]	1.90	0.30

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	297/336 (88%)	282 (95%)	11 (4%)	4 (1%)	15 9
1	B	297/336 (88%)	290 (98%)	6 (2%)	1 (0%)	46 45
All	All	594/672 (88%)	572 (96%)	17 (3%)	5 (1%)	24 17

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	162	GLN
1	B	162	GLN
1	A	52	ALA
1	A	317	ASP
1	A	161	PRO

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	253/281 (90%)	235 (93%)	18 (7%)	18 14
1	B	253/281 (90%)	240 (95%)	13 (5%)	29 26
All	All	506/562 (90%)	475 (94%)	31 (6%)	23 19

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	44	LYS
1	A	53	ASP
1	A	70	GLU
1	A	114	HIS
1	A	157	SER
1	A	171	LEU
1	A	172	ARG
1	A	189	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	191	TRP
1	A	215	MET
1	A	225	ARG
1	A	255	ARG
1	A	269	LEU
1	A	281	ILE
1	A	290	ASP
1	A	291	LEU
1	A	317	ASP
1	A	327	ARG
1	B	28	LEU
1	B	37	ARG
1	B	53	ASP
1	B	56	THR
1	B	114	HIS
1	B	189	LEU
1	B	191	TRP
1	B	215	MET
1	B	216	HIS
1	B	269	LEU
1	B	274	LEU
1	B	281	ILE
1	B	320	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	81	GLN
1	A	181	GLN
1	A	213	GLN
1	A	216	HIS
1	A	277	ASN
1	B	81	GLN
1	B	181	GLN
1	B	213	GLN
1	B	216	HIS
1	B	248	GLN
1	B	277	ASN

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

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	FEO	A	501	1,3	0,2,2	0.00	-	0,1,1	0.00	-
2	FEO	B	501	1	0,2,2	0.00	-	0,1,1	0.00	-

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	FEO	A	501	1,3	-	0/0/0/0	0/0/0/0
2	FEO	B	501	1	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	501	FEO	1	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	301/336 (89%)	0.77	28 (9%) 11 15	19, 34, 55, 62	0
1	B	301/336 (89%)	0.90	44 (14%) 3 5	17, 34, 52, 59	0
All	All	602/672 (89%)	0.84	72 (11%) 6 8	17, 34, 53, 62	0

All (72) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	161	PRO	8.2
1	B	291	LEU	7.5
1	A	161	PRO	6.9
1	A	299	ARG	6.4
1	A	319	ALA	6.1
1	B	162	GLN	5.6
1	B	317	ASP	5.5
1	A	162	GLN	5.3
1	A	317	ASP	5.2
1	A	330	TRP	5.0
1	A	160	PRO	4.9
1	B	157	SER	4.9
1	B	320	VAL	4.9
1	B	160	PRO	4.5
1	A	321	ASP	4.4
1	A	329	ASP	4.3
1	B	321	ASP	4.3
1	B	329	ASP	4.2
1	A	327	ARG	4.1
1	B	267	GLN	4.0
1	A	291	LEU	4.0
1	B	179	PRO	3.9
1	B	286	GLU	3.8
1	B	208	ARG	3.8

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	157	SER	3.7
1	B	287	ILE	3.6
1	B	70	GLU	3.6
1	B	52	ALA	3.6
1	B	288	VAL	3.5
1	A	158	GLU	3.5
1	B	158	GLU	3.4
1	B	327	ARG	3.4
1	B	272	LEU	3.4
1	A	331	SER	3.3
1	A	52	ALA	3.3
1	B	318	ASP	3.2
1	B	299	ARG	3.2
1	B	330	TRP	3.2
1	A	300	LEU	3.2
1	B	326	GLU	3.1
1	B	269	LEU	3.0
1	B	290	ASP	3.0
1	B	53	ASP	2.9
1	B	191	TRP	2.9
1	A	191	TRP	2.8
1	A	318	ASP	2.8
1	B	331	SER	2.7
1	B	254	VAL	2.7
1	B	324	PHE	2.7
1	B	300	LEU	2.7
1	A	289	GLY	2.7
1	A	316	LEU	2.6
1	B	289	GLY	2.6
1	B	197	THR	2.5
1	A	267	GLN	2.5
1	A	326	GLU	2.4
1	B	163	PRO	2.4
1	B	308	ARG	2.4
1	B	323	ASP	2.4
1	B	285	GLU	2.3
1	B	114	HIS	2.3
1	A	173	ARG	2.3
1	A	225	ARG	2.3
1	A	97	VAL	2.2
1	B	173	ARG	2.2
1	A	288	VAL	2.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	247	GLU	2.2
1	A	195	ALA	2.1
1	B	316	LEU	2.1
1	B	256	CYS	2.1
1	A	132	ALA	2.0
1	B	265	ALA	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^{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
2	FEO	B	501	3/3	0.99	0.03	-5.65	26,26,29,30	0
2	FEO	A	501	3/3	1.00	0.03	-7.83	19,19,20,22	0

6.5 Other polymers (i)

There are no such residues in this entry.