Full wwPDB X-ray Structure Validation Report (i)

Mar 5, 2024 – 12:07 PM EST

PDB ID	:	2ZYL
Title	:	Crystal structure of 3-ketosteroid-9-alpha-hydroxylase (KshA) from M. tuberculosis
Authors	:	D'Angelo, I.; Capyk, J.; Strynadka, N.; Eltis, L.
Deposited on	:	2009-01-27
Resolution	:	2.30 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
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

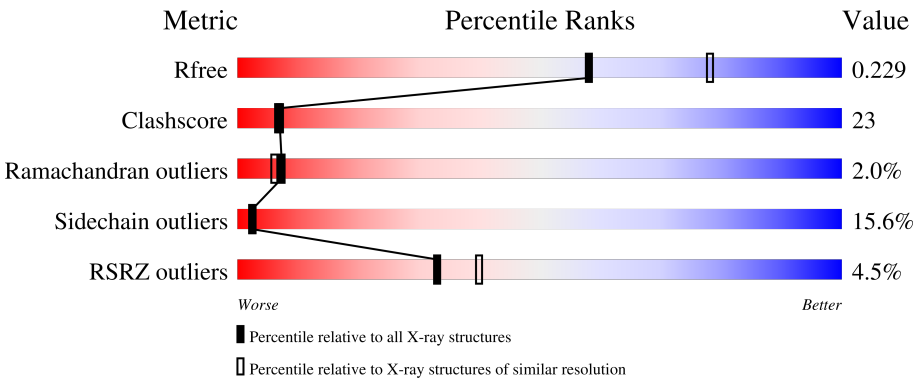
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

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

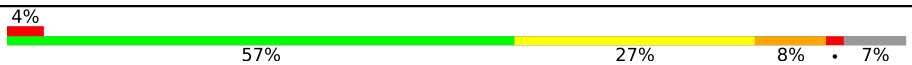
The reported resolution of this entry is 2.30 Å.

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}	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. 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	386	

2 Entry composition (i)

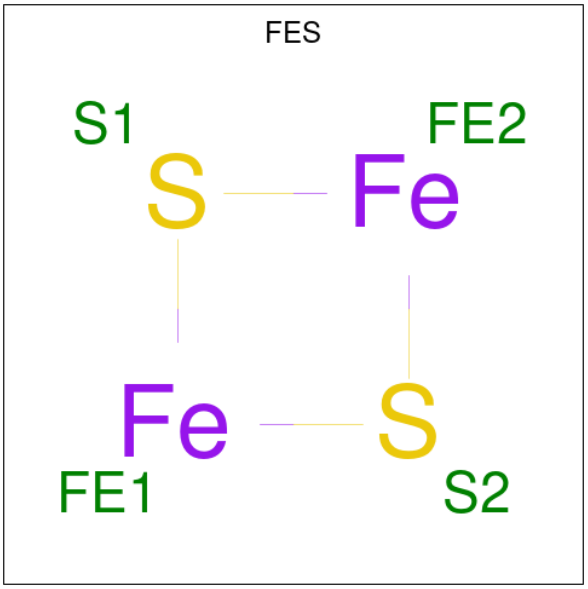
There are 5 unique types of molecules in this entry. The entry contains 3094 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 POSSIBLE OXIDOREDUCTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	359	2916	1857	509	537	13	0	0	0

- Molecule 2 is FE2/S2 (INORGANIC) CLUSTER (three-letter code: FES) (formula: Fe_2S_2).



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

- Molecule 3 is FE (II) ION (three-letter code: FE2) (formula: Fe).

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

- Molecule 4 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Na		
4	A	1	1	1	0	0

- Molecule 5 is water.

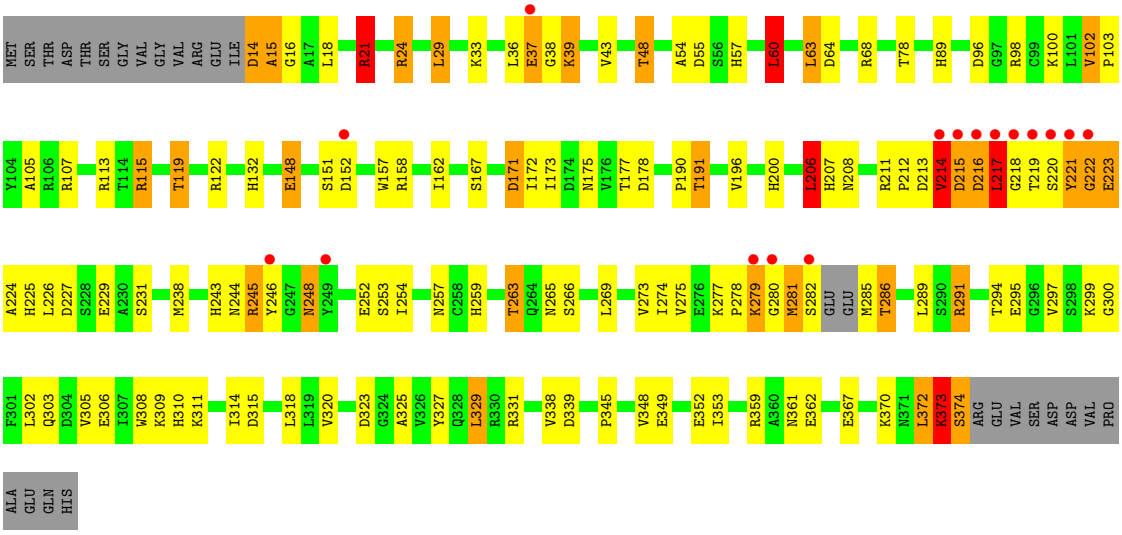
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	172	172	172	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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: POSSIBLE OXIDOREDUCTASE

Chain A: 4% 57% 27% 8% • 7%



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 3 2 1	Depositor
Cell constants a, b, c, α, β, γ	116.20Å 116.20Å 81.00Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.85 – 2.30 19.85 – 2.30	Depositor EDS
% Data completeness (in resolution range)	100.0 (19.85-2.30) 99.9 (19.85-2.30)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$< I/\sigma(I) >$ [1]	4.12 (at 2.30Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.197 , 0.235 0.195 , 0.229	Depositor DCC
R_{free} test set	1417 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	40.5	Xtriage
Anisotropy	0.286	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 42.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.026 for -h,-k,l	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	3094	wwPDB-VP
Average B, all atoms (Å^2)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.87% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 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: FE2, NA, 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	1.24	7/3004 (0.2%)	1.15	17/4079 (0.4%)

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	3

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	171	ASP	CB-CG	6.18	1.64	1.51
1	A	21	ARG	CB-CG	-5.71	1.37	1.52
1	A	367	GLU	CG-CD	5.69	1.60	1.51
1	A	327	TYR	CD1-CE1	5.59	1.47	1.39
1	A	320	VAL	CB-CG2	5.49	1.64	1.52
1	A	295	GLU	CG-CD	5.45	1.60	1.51
1	A	107	ARG	CG-CD	-5.36	1.38	1.51

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	24	ARG	NE-CZ-NH2	-11.12	114.74	120.30
1	A	178	ASP	CB-CG-OD1	-10.68	108.69	118.30
1	A	24	ARG	NE-CZ-NH1	6.94	123.77	120.30
1	A	171	ASP	CB-CG-OD1	6.79	124.41	118.30
1	A	21	ARG	CB-CA-C	-6.61	97.18	110.40
1	A	63	LEU	CB-CG-CD1	6.49	122.03	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	29	LEU	CB-CG-CD1	6.43	121.94	111.00
1	A	15	ALA	N-CA-C	6.37	128.20	111.00
1	A	113	ARG	NE-CZ-NH2	-6.06	117.27	120.30
1	A	60	LEU	CA-CB-CG	5.89	128.85	115.30
1	A	36	LEU	CA-CB-CG	5.86	128.78	115.30
1	A	102	VAL	CG1-CB-CG2	5.69	120.01	110.90
1	A	323	ASP	CB-CG-OD2	5.63	123.37	118.30
1	A	206	LEU	CA-CB-CG	5.60	128.18	115.30
1	A	68	ARG	NE-CZ-NH2	-5.55	117.52	120.30
1	A	222	GLY	N-CA-C	-5.46	99.46	113.10
1	A	96	ASP	CB-CG-OD1	5.25	123.02	118.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	14	ASP	Peptide
1	A	214	VAL	Peptide
1	A	37	GLU	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	2916	0	2752	130	0
2	A	4	0	0	1	0
3	A	1	0	0	0	0
4	A	1	0	0	0	0
5	A	172	0	0	35	0
All	All	3094	0	2752	130	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 23.

All (130) 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:280:GLY:HA2	1:A:281:MET:CB	1.47	1.40
1:A:280:GLY:CA	1:A:281:MET:HB2	1.61	1.30
1:A:286:THR:HA	5:A:462:HOH:O	1.33	1.28
1:A:282:SER:HB2	5:A:494:HOH:O	1.33	1.23
1:A:217:LEU:HD22	1:A:218:GLY:N	1.58	1.18
1:A:222:GLY:O	1:A:223:GLU:HB3	1.49	1.06
1:A:64:ASP:OD1	1:A:115:ARG:NH1	1.91	1.03
1:A:217:LEU:HD22	1:A:218:GLY:H	1.20	1.00
1:A:15:ALA:HB1	1:A:315:ASP:H	1.23	1.00
1:A:100:LYS:HE2	5:A:438:HOH:O	1.61	1.00
1:A:280:GLY:HA2	1:A:281:MET:HB3	1.44	0.99
1:A:280:GLY:HA2	1:A:281:MET:HB2	0.99	0.99
1:A:374:SER:HB2	5:A:558:HOH:O	1.68	0.92
1:A:263:THR:HG22	1:A:265:ASN:H	1.32	0.92
1:A:217:LEU:CD2	1:A:218:GLY:N	2.33	0.91
1:A:217:LEU:CD2	1:A:218:GLY:H	1.84	0.90
1:A:171:ASP:HB3	5:A:481:HOH:O	1.71	0.90
1:A:38:GLY:HA2	5:A:473:HOH:O	1.73	0.87
1:A:280:GLY:CA	1:A:281:MET:CB	2.29	0.84
1:A:238:MET:HE3	5:A:512:HOH:O	1.76	0.84
1:A:285:MET:HA	5:A:494:HOH:O	1.76	0.83
1:A:39:LYS:HE2	5:A:461:HOH:O	1.79	0.82
1:A:158:ARG:HH12	1:A:291:ARG:HG2	1.44	0.82
1:A:280:GLY:N	1:A:281:MET:HE3	1.95	0.82
1:A:60:LEU:O	1:A:119:THR:HG21	1.79	0.82
1:A:285:MET:HG2	1:A:286:THR:H	1.46	0.81
1:A:158:ARG:HD2	1:A:294:THR:HG21	1.63	0.81
1:A:243:HIS:HD2	1:A:252:GLU:OE1	1.62	0.81
1:A:285:MET:HG3	5:A:526:HOH:O	1.80	0.81
1:A:248:ASN:HB3	5:A:429:HOH:O	1.80	0.81
1:A:280:GLY:N	1:A:281:MET:CE	2.43	0.80
1:A:216:ASP:O	1:A:222:GLY:CA	2.30	0.79
1:A:223:GLU:CG	1:A:223:GLU:O	2.30	0.78
1:A:265:ASN:HB2	5:A:470:HOH:O	1.83	0.77
1:A:359:ARG:HB3	5:A:532:HOH:O	1.84	0.76
1:A:285:MET:CG	1:A:286:THR:H	2.00	0.75
1:A:280:GLY:H	1:A:281:MET:HE2	1.52	0.74
1:A:374:SER:HB2	5:A:440:HOH:O	1.87	0.73
1:A:280:GLY:N	1:A:281:MET:HB2	2.05	0.70
1:A:54:ALA:HB2	1:A:60:LEU:HD13	1.73	0.69
1:A:15:ALA:CB	1:A:315:ASP:H	2.02	0.68
1:A:151:SER:O	5:A:560:HOH:O	2.11	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:229:GLU:OE1	1:A:243:HIS:HE1	1.76	0.68
1:A:15:ALA:HB1	1:A:315:ASP:N	2.05	0.67
1:A:211:ARG:HB2	1:A:214:VAL:CG2	2.24	0.67
1:A:200:HIS:HE1	5:A:431:HOH:O	1.79	0.66
1:A:223:GLU:O	1:A:223:GLU:HG2	1.95	0.65
1:A:224:ALA:HA	1:A:245:ARG:O	1.97	0.65
1:A:217:LEU:CG	1:A:218:GLY:H	2.05	0.64
1:A:216:ASP:N	1:A:216:ASP:OD1	2.30	0.63
1:A:38:GLY:CA	5:A:473:HOH:O	2.37	0.63
1:A:102:VAL:HG22	1:A:105:ALA:CB	2.29	0.63
1:A:191:THR:CG2	5:A:547:HOH:O	2.46	0.62
1:A:345:PRO:O	1:A:349:GLU:HG2	1.99	0.61
1:A:216:ASP:O	1:A:222:GLY:HA3	2.00	0.61
1:A:222:GLY:O	1:A:223:GLU:CB	2.32	0.61
1:A:215:ASP:HB2	5:A:525:HOH:O	2.00	0.60
1:A:175:ASN:ND2	1:A:308:TRP:HE1	1.99	0.60
1:A:207:HIS:HD2	1:A:227:ASP:OD1	1.83	0.60
1:A:14:ASP:HB3	1:A:16:GLY:N	2.16	0.59
1:A:238:MET:CE	5:A:512:HOH:O	2.42	0.59
1:A:218:GLY:HA3	1:A:219:THR:C	2.20	0.59
1:A:211:ARG:HD2	1:A:213:ASP:OD2	2.03	0.59
1:A:102:VAL:HG22	1:A:105:ALA:HB3	1.85	0.57
1:A:216:ASP:O	1:A:222:GLY:HA2	2.02	0.57
1:A:98:ARG:HB2	1:A:100:LYS:HE3	1.84	0.57
1:A:24:ARG:NH2	1:A:338:VAL:O	2.37	0.57
1:A:217:LEU:HD22	1:A:218:GLY:CA	2.31	0.56
1:A:359:ARG:HD2	5:A:411:HOH:O	2.06	0.56
1:A:225:HIS:O	1:A:244:ASN:HB2	2.07	0.55
1:A:217:LEU:HD22	1:A:219:THR:HA	1.88	0.54
1:A:263:THR:HB	1:A:266:SER:OG	2.08	0.53
1:A:191:THR:HG22	5:A:547:HOH:O	2.08	0.53
1:A:238:MET:HE2	1:A:257:ASN:HB3	1.91	0.52
1:A:212:PRO:O	1:A:217:LEU:HB2	2.10	0.52
1:A:285:MET:CG	1:A:286:THR:N	2.71	0.52
1:A:325:ALA:HB2	1:A:353:ILE:HG21	1.92	0.52
1:A:374:SER:CB	5:A:558:HOH:O	2.40	0.52
1:A:259:HIS:HE1	5:A:400:HOH:O	1.91	0.52
1:A:285:MET:HG2	1:A:286:THR:N	2.22	0.51
1:A:216:ASP:O	1:A:217:LEU:O	2.30	0.50
1:A:280:GLY:H	1:A:281:MET:CE	2.12	0.49
1:A:64:ASP:CG	1:A:115:ARG:NH1	2.64	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:329:LEU:HD12	5:A:432:HOH:O	2.11	0.49
1:A:15:ALA:CB	1:A:314:ILE:HA	2.43	0.48
1:A:119:THR:HG23	5:A:553:HOH:O	2.13	0.48
1:A:214:VAL:O	1:A:216:ASP:N	2.37	0.48
1:A:14:ASP:HB3	1:A:15:ALA:C	2.34	0.48
1:A:373:LYS:C	5:A:440:HOH:O	2.51	0.48
1:A:102:VAL:HG22	1:A:105:ALA:HB2	1.96	0.47
1:A:281:MET:HB3	1:A:286:THR:HG22	1.96	0.47
1:A:359:ARG:CD	5:A:411:HOH:O	2.62	0.47
1:A:300:GLY:O	1:A:303:GLN:HG2	2.14	0.47
1:A:21:ARG:NH2	1:A:339:ASP:OD2	2.34	0.47
1:A:217:LEU:O	1:A:222:GLY:N	2.48	0.47
1:A:259:HIS:HD2	5:A:406:HOH:O	1.96	0.47
1:A:223:GLU:O	1:A:223:GLU:HG3	2.13	0.46
1:A:211:ARG:CB	1:A:214:VAL:HG22	2.46	0.46
1:A:372:LEU:C	1:A:374:SER:H	2.19	0.46
1:A:14:ASP:HB3	1:A:16:GLY:H	1.80	0.46
1:A:57:HIS:HE1	5:A:439:HOH:O	1.98	0.46
1:A:279:LYS:C	1:A:281:MET:HE3	2.36	0.45
1:A:277:LYS:HA	1:A:278:PRO:HD3	1.70	0.45
1:A:102:VAL:HA	1:A:103:PRO:HD3	1.82	0.45
1:A:217:LEU:CD2	1:A:219:THR:HA	2.46	0.45
1:A:21:ARG:HG2	1:A:132:HIS:CG	2.52	0.45
1:A:200:HIS:CD2	1:A:200:HIS:H	2.34	0.44
1:A:281:MET:HB3	1:A:286:THR:CG2	2.46	0.44
1:A:305:VAL:HG12	1:A:309:LYS:HD2	1.99	0.44
1:A:374:SER:C	5:A:471:HOH:O	2.55	0.44
1:A:98:ARG:HD2	5:A:438:HOH:O	2.17	0.44
1:A:211:ARG:HB2	1:A:214:VAL:HG22	2.00	0.43
1:A:196:VAL:HG22	1:A:352:GLU:HG2	1.99	0.43
1:A:190:PRO:O	1:A:361:ASN:ND2	2.46	0.43
1:A:157:TRP:CE3	1:A:274:ILE:HG23	2.54	0.43
1:A:98:ARG:NH2	5:A:559:HOH:O	2.45	0.43
1:A:211:ARG:CB	1:A:214:VAL:CG2	2.96	0.43
1:A:215:ASP:CB	5:A:525:HOH:O	2.64	0.43
1:A:89:HIS:HB2	2:A:451:FES:S2	2.59	0.42
1:A:18:LEU:HD23	1:A:18:LEU:HA	1.89	0.41
1:A:206:LEU:HD13	1:A:208:ASN:HB2	2.02	0.41
1:A:359:ARG:O	1:A:362:GLU:HG2	2.20	0.41
1:A:253:SER:HB3	1:A:275:VAL:HG12	2.02	0.41
1:A:148:GLU:HG2	1:A:254:ILE:HD12	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:162:ILE:HG21	1:A:162:ILE:HD13	1.87	0.41
1:A:231:SER:O	1:A:238:MET:HA	2.21	0.41
1:A:172:ILE:HD13	1:A:172:ILE:HG21	1.85	0.41
1:A:48:THR:HG21	5:A:556:HOH:O	2.21	0.41
1:A:243:HIS:CD2	1:A:252:GLU:OE1	2.55	0.41
1:A:306:GLU:O	1:A:310:HIS:HD2	2.04	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	355/386 (92%)	337 (95%)	11 (3%)	7 (2%)	7 6

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	55	ASP
1	A	215	ASP
1	A	217	LEU
1	A	223	GLU
1	A	281	MET
1	A	221	TYR
1	A	373	LYS

5.3.2 Protein sidechains (i)

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

The Analysed column shows the number of residues for which the sidechain conformation was

analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	308/333 (92%)	260 (84%)	48 (16%)	2 2

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	ARG
1	A	29	LEU
1	A	33	LYS
1	A	37	GLU
1	A	39	LYS
1	A	43	VAL
1	A	48	THR
1	A	60	LEU
1	A	63	LEU
1	A	78	THR
1	A	115	ARG
1	A	119	THR
1	A	122	ARG
1	A	148	GLU
1	A	152	ASP
1	A	167	SER
1	A	173	ILE
1	A	177	THR
1	A	191	THR
1	A	206	LEU
1	A	214	VAL
1	A	216	ASP
1	A	217	LEU
1	A	220	SER
1	A	221	TYR
1	A	226	LEU
1	A	245	ARG
1	A	246	TYR
1	A	248	ASN
1	A	263	THR
1	A	269	LEU
1	A	273	VAL
1	A	279	LYS
1	A	286	THR
1	A	289	LEU
1	A	291	ARG

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Mol	Chain	Res	Type
1	A	297	VAL
1	A	299	LYS
1	A	302	LEU
1	A	311	LYS
1	A	318	LEU
1	A	329	LEU
1	A	331	ARG
1	A	348	VAL
1	A	370	LYS
1	A	372	LEU
1	A	373	LYS
1	A	374	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	175	ASN
1	A	200	HIS
1	A	207	HIS
1	A	225	HIS
1	A	243	HIS
1	A	257	ASN
1	A	259	HIS
1	A	310	HIS
1	A	335	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	FES	A	451	1	0,4,4	-	-	-		

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	451	1	-	-	0/1/1/1

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	A	451	FES	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 (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	359/386 (93%)	-0.12	16 (4%) 33 40	19, 36, 73, 110	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	218	GLY	13.9
1	A	220	SER	7.6
1	A	217	LEU	7.0
1	A	219	THR	6.7
1	A	222	GLY	6.6
1	A	216	ASP	5.3
1	A	221	TYR	5.1
1	A	215	ASP	5.1
1	A	246	TYR	4.6
1	A	280	GLY	3.1
1	A	249	TYR	3.1
1	A	37	GLU	2.7
1	A	282	SER	2.5
1	A	214	VAL	2.4
1	A	279	LYS	2.3
1	A	152	ASP	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 monosaccharides 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. 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	B-factors(Å^2)	Q<0.9
4	NA	A	387	1/1	0.96	0.06	41,41,41,41	0
2	FES	A	451	4/4	0.99	0.11	26,29,29,30	0
3	FE2	A	452	1/1	1.00	0.09	34,34,34,34	0

6.5 Other polymers (i)

There are no such residues in this entry.