



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 18, 2026 – 12:33 PM UTC

PDB ID : 2FSS / pdb_00002fss
Title : Candida boidinii formate dehydrogenase (FDH) K47E mutant
Authors : Schirwitz, K.; Schmidt, A.; Lamzin, V.S.
Deposited on : 2006-01-23
Resolution : 1.70 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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtrriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

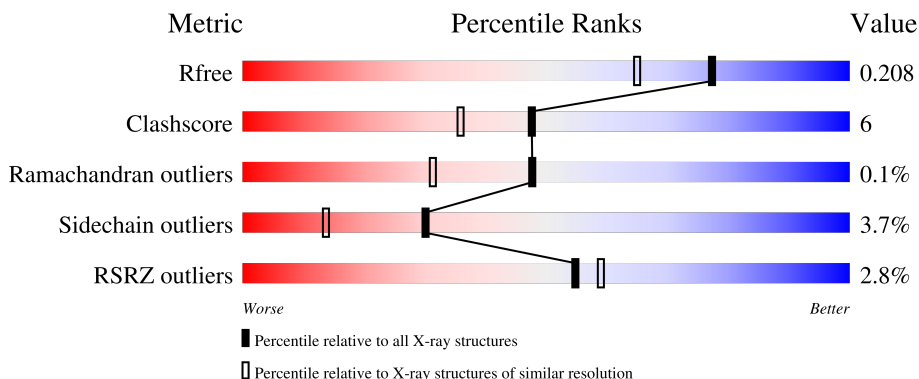
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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))
R_{free}	180053	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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 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 $\leq 5\%$. 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	365	 86% 9% ..
1	B	365	 3% 79% 15% . 5%
1	C	365	 2% 87% 9% ..
1	D	365	 4% 78% 13% . 8%

2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 11848 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 formate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	353	2761	1758	472	525	6	0	2	0
1	B	346	2707	1727	464	510	6	0	2	0
1	C	353	2755	1754	473	522	6	0	1	0
1	D	337	2653	1699	452	496	6	0	3	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	cloning artifact	GB 7657869
A	1	ALA	-	cloning artifact	GB 7657869
A	47	GLU	LYS	engineered mutation	GB 7657869
B	0	MET	-	cloning artifact	GB 7657869
B	1	ALA	-	cloning artifact	GB 7657869
B	47	GLU	LYS	engineered mutation	GB 7657869
C	0	MET	-	cloning artifact	GB 7657869
C	1	ALA	-	cloning artifact	GB 7657869
C	47	GLU	LYS	engineered mutation	GB 7657869
D	0	MET	-	cloning artifact	GB 7657869
D	1	ALA	-	cloning artifact	GB 7657869
D	47	GLU	LYS	engineered mutation	GB 7657869

- Molecule 2 is SULFATE ION (CCD ID: 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		
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		

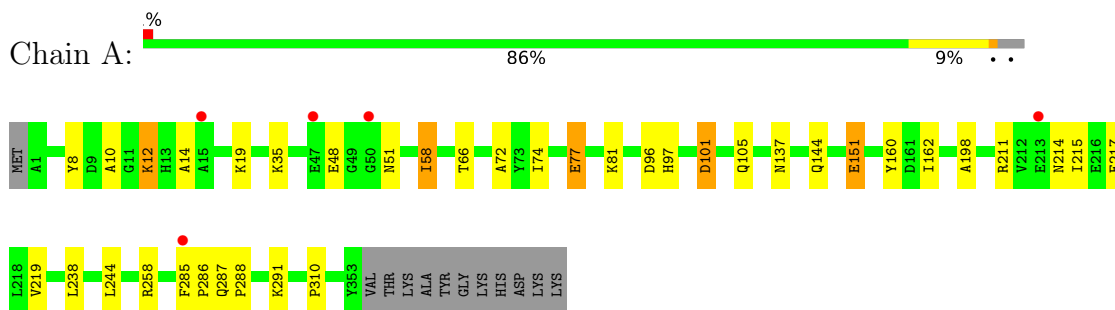
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	270	Total	O	0	0
			270	270		
3	B	224	Total	O	0	0
			224	224		
3	C	260	Total	O	0	0
			260	260		
3	D	193	Total	O	0	0
			193	193		

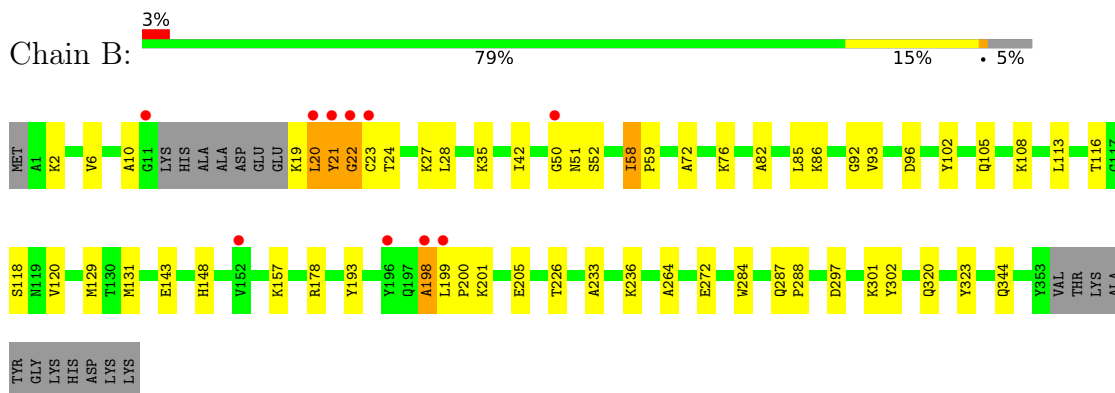
3 Residue-property plots

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.

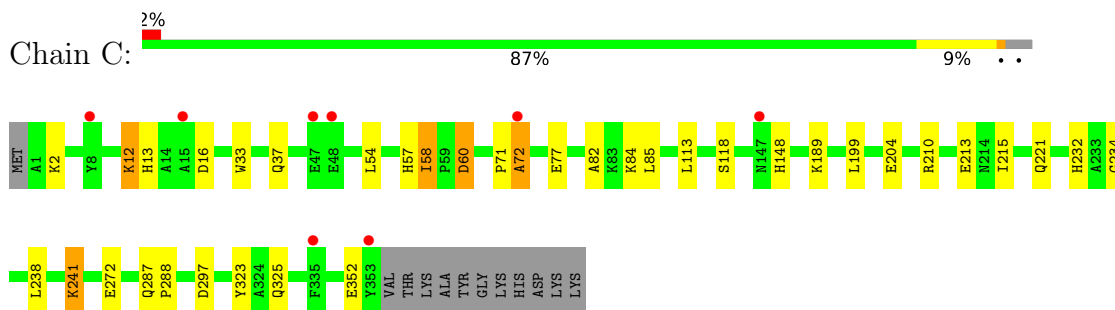
- Molecule 1: formate dehydrogenase



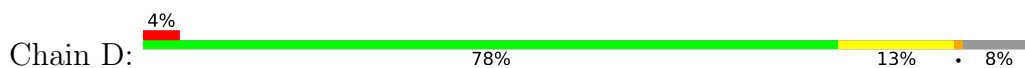
- Molecule 1: formate dehydrogenase

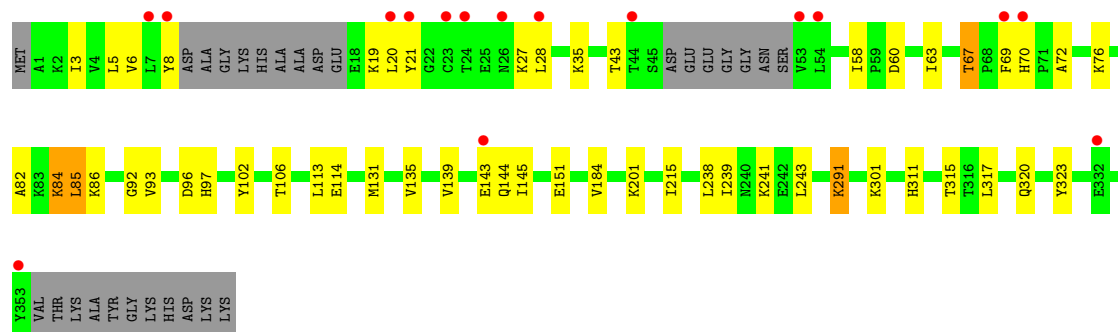


- Molecule 1: formate dehydrogenase



- Molecule 1: formate dehydrogenase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	53.43Å 68.39Å 109.37Å 77.93° 89.33° 81.34°	Depositor
Resolution (Å)	19.42 – 1.70 19.42 – 1.70	Depositor EDS
% Data completeness (in resolution range)	(Not available) (19.42-1.70) 96.5 (19.42-1.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.93 (at 1.70Å)	Xtrriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.204 , 0.249 0.205 , 0.208	Depositor DCC
R_{free} test set	7798 reflections (4.74%)	wwPDB-VP
Wilson B-factor (Å ²)	30.5	Xtrriage
Anisotropy	0.134	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 34.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11848	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.25% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ 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: 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	0.74	1/2824 (0.0%)	0.88	0/3834
1	B	0.72	0/2769	0.93	2/3759 (0.1%)
1	C	0.76	1/2815 (0.0%)	0.94	0/3822
1	D	0.70	1/2716 (0.0%)	0.88	5/3688 (0.1%)
All	All	0.73	3/11124 (0.0%)	0.91	7/15103 (0.0%)

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	3
1	D	0	1
All	All	0	4

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	184	VAL	CA-CB	9.17	1.58	1.54
1	A	58	ILE	CA-CB	7.48	1.57	1.54
1	C	58	ILE	CA-CB	5.16	1.56	1.54

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	10	ALA	N-CA-C	7.21	115.67	108.75
1	B	264	ALA	N-CA-C	5.28	117.45	111.11
1	D	184	VAL	N-CA-CB	5.14	113.96	110.52
1	D	67[A]	THR	CA-C-N	-5.09	114.80	120.45

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	67[A]	THR	C-N-CA	-5.09	114.80	120.45
1	D	67[B]	THR	CA-C-N	-5.09	114.80	120.45
1	D	67[B]	THR	C-N-CA	-5.09	114.80	120.45

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	198	ALA	Peptide
1	B	21	TYR	Peptide
1	B	22	GLY	Peptide
1	D	19	LYS	Peptide

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	2761	0	2757	28	0
1	B	2707	0	2711	39	0
1	C	2755	0	2753	31	0
1	D	2653	0	2675	30	0
2	A	10	0	0	0	0
2	C	10	0	0	0	0
2	D	5	0	0	0	0
3	A	270	0	0	5	0
3	B	224	0	0	3	0
3	C	260	0	0	3	0
3	D	193	0	0	4	0
All	All	11848	0	10896	122	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 6.

All (122) 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:19:LYS:HG2	3:A:2275:HOH:O	1.59	1.01
1:C:71:PRO:HA	1:C:72:ALA:HB3	1.40	1.00
1:C:71:PRO:HA	1:C:72:ALA:CB	1.92	0.99
1:A:12:LYS:H	1:A:12:LYS:HD2	1.25	0.97
1:D:291:LYS:HE3	1:D:291:LYS:H	1.35	0.89
1:A:66:THR:HG21	1:A:74:ILE:HD11	1.53	0.87
1:D:6:VAL:HG11	1:D:72:ALA:HB2	1.56	0.87
1:A:12:LYS:H	1:A:12:LYS:CD	1.88	0.84
1:B:120:VAL:HG12	1:B:178:ARG:HD2	1.63	0.80
1:D:60:ASP:HA	1:D:84:LYS:HG3	1.65	0.78
1:A:72:ALA:O	1:A:74:ILE:HD12	1.84	0.78
1:A:144:GLN:HE22	1:A:151:GLU:H	1.32	0.75
1:C:210:ARG:HH22	1:C:221:GLN:NE2	1.86	0.74
1:D:311:HIS:HD2	3:D:2072:HOH:O	1.73	0.71
1:D:93:VAL:HG13	1:D:114:GLU:OE1	1.90	0.71
1:C:232:HIS:HD2	1:C:234:GLY:H	1.37	0.70
1:D:67[B]:THR:HG22	1:D:69:PHE:H	1.57	0.69
1:C:241:LYS:HB2	1:C:241:LYS:NZ	2.07	0.68
1:A:8:TYR:HE2	3:A:2181:HOH:O	1.76	0.68
1:A:12:LYS:HD2	1:A:12:LYS:N	2.04	0.68
1:B:198:ALA:HB1	1:B:199:LEU:O	1.94	0.67
1:C:82:ALA:HB1	1:C:85:LEU:HB2	1.78	0.66
1:B:120:VAL:CG1	1:B:178:ARG:HD2	2.26	0.65
1:B:116:THR:HG21	1:B:344:GLN:O	1.97	0.64
1:C:60:ASP:HA	1:C:84:LYS:HD3	1.80	0.64
1:A:101:ASP:OD1	1:A:105:GLN:NE2	2.32	0.63
1:D:139:VAL:O	1:D:143:GLU:HG2	1.98	0.63
1:C:71:PRO:CA	1:C:72:ALA:HB3	2.21	0.63
1:A:144:GLN:NE2	1:A:151:GLU:H	1.97	0.63
1:C:210:ARG:HH22	1:C:221:GLN:HE22	1.47	0.62
1:A:287:GLN:HA	1:A:288:PRO:C	2.24	0.62
1:D:144:GLN:HE22	1:D:151:GLU:H	1.46	0.62
1:D:20:LEU:HD23	1:D:320:GLN:HE22	1.64	0.61
1:D:301:LYS:HD3	3:D:2031:HOH:O	2.00	0.60
1:A:51:ASN:HA	3:A:2156:HOH:O	2.01	0.60
1:C:325[B]:GLN:HG3	3:C:2075:HOH:O	2.03	0.59
1:B:120:VAL:HG12	1:B:178:ARG:CD	2.32	0.59
1:B:50:GLY:CA	1:B:51:ASN:HB3	2.34	0.58
1:B:50:GLY:HA3	1:B:52:SER:H	1.68	0.57
1:C:71:PRO:HA	1:C:72:ALA:HB2	1.85	0.57
1:D:102:TYR:O	1:D:106:THR:HG23	2.04	0.57
1:A:77:GLU:H	1:A:77:GLU:CD	2.11	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:131:MET:CE	1:B:226:THR:OG1	2.52	0.56
1:B:157:LYS:NZ	1:C:13:HIS:HD2	2.03	0.56
1:B:92:GLY:HA2	1:B:323:TYR:CE2	2.40	0.56
1:B:233:ALA:HA	1:B:236:LYS:HE3	1.88	0.56
1:D:8:TYR:HB2	1:D:70:HIS:CG	2.41	0.56
1:B:42:ILE:HD12	1:B:42:ILE:N	2.21	0.56
1:A:66:THR:HG21	1:A:74:ILE:CD1	2.33	0.55
1:B:118:SER:OG	1:B:323:TYR:HE2	1.90	0.55
1:B:20:LEU:HD23	1:B:320:GLN:OE1	2.08	0.54
1:B:50:GLY:HA2	1:B:51:ASN:HB3	1.90	0.54
1:C:241:LYS:HB2	1:C:241:LYS:HZ2	1.73	0.53
1:A:215:ILE:HD13	1:A:238:LEU:HD21	1.90	0.53
1:B:76:LYS:HG2	1:B:102:TYR:CG	2.43	0.53
1:B:287:GLN:HA	1:B:288:PRO:C	2.34	0.53
1:B:148:HIS:HE1	1:C:297:ASP:OD1	1.91	0.53
1:D:21:TYR:HB3	1:D:28:LEU:HA	1.92	0.52
1:B:118:SER:OG	1:B:323:TYR:CE2	2.62	0.52
1:C:241:LYS:NZ	1:C:241:LYS:CB	2.73	0.52
1:C:232:HIS:HE1	3:C:2043:HOH:O	1.92	0.52
1:A:310:PRO:HD3	1:D:145:ILE:HG12	1.92	0.51
1:A:10:ALA:HB3	1:A:14:ALA:HB2	1.93	0.51
1:D:92:GLY:HA2	1:D:323:TYR:CD1	2.46	0.51
1:B:82:ALA:HB1	1:B:85:LEU:HB2	1.94	0.50
1:D:58:ILE:HG23	1:D:82:ALA:HB2	1.94	0.50
1:D:3:ILE:HG12	1:D:63:ILE:HD12	1.94	0.49
1:A:19:LYS:HD3	3:D:2088:HOH:O	2.14	0.48
1:C:215:ILE:HD13	1:C:238:LEU:HD21	1.96	0.48
1:D:92:GLY:HA2	1:D:323:TYR:CE1	2.49	0.47
1:A:285:PHE:HA	1:A:286:PRO:C	2.38	0.47
1:C:232:HIS:HD2	1:C:234:GLY:N	2.10	0.47
1:C:241:LYS:HB2	1:C:241:LYS:HZ3	1.79	0.46
1:D:215:ILE:HD12	1:D:238:LEU:HD11	1.97	0.46
1:B:297:ASP:OD1	1:C:148:HIS:HE1	1.98	0.46
1:C:118:SER:OG	1:C:323:TYR:HE1	1.99	0.46
1:B:86:LYS:HE3	3:B:520:HOH:O	2.16	0.46
1:B:193:TYR:OH	1:B:199:LEU:HB2	2.15	0.46
1:C:287:GLN:HA	1:C:288:PRO:C	2.40	0.46
1:D:97:HIS:HB3	3:D:2079:HOH:O	2.15	0.46
1:D:82:ALA:HB1	1:D:85:LEU:HB2	1.99	0.45
1:B:284:TRP:H	1:B:287:GLN:NE2	2.14	0.45
1:B:108:LYS:NZ	3:B:531:HOH:O	2.48	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:199:LEU:HA	1:B:200:PRO:HD2	1.76	0.45
1:C:297:ASP:HB2	3:C:2050:HOH:O	2.15	0.45
1:A:215:ILE:O	1:A:219:VAL:HG22	2.17	0.45
1:C:2:LYS:NZ	1:C:57:HIS:HD2	2.14	0.45
1:B:302:TYR:OH	1:C:16:ASP:HB3	2.17	0.45
1:B:201:LYS:HD3	1:B:201:LYS:HA	1.73	0.44
1:B:2:LYS:HE2	1:B:42:ILE:CD1	2.48	0.44
1:B:6:VAL:HG11	1:B:72:ALA:HB2	1.99	0.44
1:B:24:THR:O	1:B:27:LYS:HG3	2.18	0.43
1:D:76:LYS:HG2	1:D:102:TYR:CE2	2.53	0.43
1:C:118:SER:OG	1:C:323:TYR:CE1	2.69	0.43
1:C:12:LYS:H	1:C:12:LYS:HZ3	1.66	0.43
1:C:33:TRP:CE2	1:C:37:GLN:OE1	2.71	0.43
1:C:58:ILE:HG23	1:C:82:ALA:HB2	2.01	0.43
1:D:27:LYS:NZ	1:D:43:THR:HG21	2.33	0.43
1:A:198:ALA:HB2	1:A:211:ARG:NH1	2.33	0.43
1:A:214:ASN:HD21	1:A:217:GLU:CD	2.26	0.43
1:A:77:GLU:O	1:A:81:LYS:HG2	2.18	0.43
1:D:131:MET:O	1:D:135:VAL:HG23	2.19	0.43
1:B:21:TYR:HB3	1:B:28:LEU:HA	2.02	0.42
1:D:144:GLN:NE2	1:D:151:GLU:H	2.14	0.42
1:D:239:ILE:HD13	1:D:243:LEU:HD23	2.01	0.42
1:B:58:ILE:N	1:B:59:PRO:CD	2.82	0.41
1:C:77:GLU:H	1:C:77:GLU:CD	2.28	0.41
1:D:8:TYR:HB2	1:D:70:HIS:CD2	2.55	0.41
1:D:20:LEU:CD2	1:D:320:GLN:HE22	2.30	0.41
1:B:301:LYS:HE3	1:B:302:TYR:CZ	2.56	0.41
1:A:160:TYR:CZ	1:D:317:LEU:HD12	2.55	0.41
1:D:20:LEU:HD23	1:D:20:LEU:HA	1.82	0.41
1:A:96:ASP:OD1	1:A:258:ARG:NH1	2.50	0.41
1:B:2:LYS:HE2	1:B:42:ILE:HD11	2.03	0.41
1:B:92:GLY:HA2	1:B:323:TYR:CD2	2.56	0.41
1:B:226:THR:HG22	3:B:380:HOH:O	2.20	0.41
1:C:199:LEU:HB2	1:C:204:GLU:HG3	2.02	0.41
1:A:8:TYR:CE2	3:A:2181:HOH:O	2.57	0.41
1:A:72:ALA:O	1:A:97:HIS:HD2	2.03	0.40
1:B:129:MET:HE2	1:B:129:MET:HB3	1.92	0.40
1:A:19:LYS:CG	3:A:2275:HOH:O	2.40	0.40
1:B:21:TYR:O	1:B:23:CYS:N	2.54	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	353/365 (97%)	339 (96%)	14 (4%)	0	100	100
1	B	344/365 (94%)	327 (95%)	16 (5%)	1 (0%)	36	22
1	C	352/365 (96%)	341 (97%)	10 (3%)	1 (0%)	36	22
1	D	334/365 (92%)	324 (97%)	10 (3%)	0	100	100
All	All	1383/1460 (95%)	1331 (96%)	50 (4%)	2 (0%)	48	31

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	22	GLY
1	C	72	ALA

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	290/298 (97%)	279 (96%)	11 (4%)	29	13
1	B	285/298 (96%)	273 (96%)	12 (4%)	26	11
1	C	289/298 (97%)	280 (97%)	9 (3%)	35	18
1	D	281/298 (94%)	270 (96%)	11 (4%)	28	12
All	All	1145/1192 (96%)	1102 (96%)	43 (4%)	30	13

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	LYS
1	A	35	LYS
1	A	48	GLU
1	A	58	ILE
1	A	77	GLU
1	A	101	ASP
1	A	137	ASN
1	A	151	GLU
1	A	162	ILE
1	A	244	LEU
1	A	291	LYS
1	B	19	LYS
1	B	20	LEU
1	B	35	LYS
1	B	58	ILE
1	B	93	VAL
1	B	96	ASP
1	B	105	GLN
1	B	113	LEU
1	B	143[A]	GLU
1	B	143[B]	GLU
1	B	205	GLU
1	B	272	GLU
1	C	12	LYS
1	C	54	LEU
1	C	60	ASP
1	C	113	LEU
1	C	189	LYS
1	C	213	GLU
1	C	241	LYS
1	C	272	GLU
1	C	352	GLU
1	D	5	LEU
1	D	35	LYS
1	D	84	LYS
1	D	85	LEU
1	D	86	LYS
1	D	96	ASP
1	D	113	LEU
1	D	201	LYS
1	D	241	LYS
1	D	291	LYS
1	D	315	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	137	ASN
1	A	144	GLN
1	B	37	GLN
1	B	105	GLN
1	B	126	HIS
1	B	148	HIS
1	B	275	GLN
1	B	287	GLN
1	B	311	HIS
1	C	13	HIS
1	C	37	GLN
1	C	57	HIS
1	C	105	GLN
1	C	126	HIS
1	C	147	ASN
1	C	148	HIS
1	C	221	GLN
1	C	232	HIS
1	C	275	GLN
1	C	287	GLN
1	C	311	HIS
1	D	56	GLN
1	D	126	HIS
1	D	144	GLN
1	D	311	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

5 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	SO4	C	2004	-	4,4,4	0.25	0	6,6,6	0.26	0
2	SO4	A	2003	-	4,4,4	0.32	0	6,6,6	0.38	0
2	SO4	C	2002	-	4,4,4	0.26	0	6,6,6	0.52	0
2	SO4	D	2001	-	4,4,4	0.23	0	6,6,6	0.17	0
2	SO4	A	2005	-	4,4,4	0.24	0	6,6,6	0.35	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.

No monomer is involved in short contacts.

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

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	353/365 (96%)	0.21	5 (1%) 73 77	22, 32, 45, 53	2 (0%)
1	B	346/365 (94%)	0.23	10 (2%) 53 58	17, 32, 49, 57	2 (0%)
1	C	353/365 (96%)	0.26	8 (2%) 61 65	20, 32, 50, 63	1 (0%)
1	D	337/365 (92%)	0.45	16 (4%) 36 40	22, 35, 58, 70	3 (0%)
All	All	1389/1460 (95%)	0.29	39 (2%) 55 59	17, 33, 51, 70	8 (0%)

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	20	LEU	4.6
1	D	8	TYR	4.3
1	D	53	VAL	4.1
1	C	47	GLU	3.5
1	D	24	THR	3.3
1	B	23	CYS	3.3
1	B	22	GLY	3.2
1	D	21	TYR	3.1
1	B	21	TYR	3.0
1	B	50	GLY	2.9
1	C	72	ALA	2.9
1	B	196	TYR	2.8
1	A	50	GLY	2.7
1	B	198	ALA	2.7
1	A	15	ALA	2.6
1	A	47	GLU	2.5
1	C	48	GLU	2.5
1	C	147	ASN	2.5
1	D	69	PHE	2.5
1	D	70	HIS	2.4
1	D	332	GLU	2.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	D	23	CYS	2.4
1	B	11	GLY	2.4
1	D	28	LEU	2.4
1	C	8	TYR	2.3
1	D	44	THR	2.3
1	A	285	PHE	2.3
1	B	199	LEU	2.3
1	D	7	LEU	2.3
1	C	335	PHE	2.2
1	C	353	TYR	2.2
1	D	353	TYR	2.2
1	A	213	GLU	2.2
1	D	26	ASN	2.2
1	B	20	LEU	2.1
1	D	143	GLU	2.1
1	D	54	LEU	2.1
1	C	15	ALA	2.0
1	B	152	VAL	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 oligosaccharides 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(\AA^2)	Q<0.9
2	SO4	A	2003	5/5	0.87	0.09	53,56,56,57	0
2	SO4	A	2005	5/5	0.94	0.07	49,51,52,52	0
2	SO4	C	2002	5/5	0.94	0.09	39,42,43,43	0
2	SO4	D	2001	5/5	0.94	0.06	56,56,57,58	0
2	SO4	C	2004	5/5	0.95	0.06	44,46,47,47	0

6.5 Other polymers [i](#)

There are no such residues in this entry.