Full wwPDB X-ray Structure Validation Report (i)

Aug 26, 2023 – 10:26 PM EDT

PDB ID	:	3GAK
Title	:	Structure of Giardia fructose-1,6-biphosphate aldolase
Authors	:	Galkin, A.; Herzberg, O.
Deposited on	:	2009-02-17
Resolution	:	2.90 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:

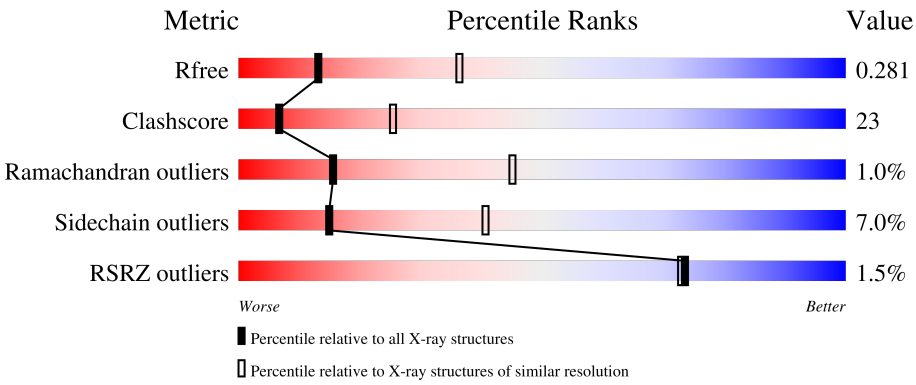
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

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 2.90 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	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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	323	.% 50% 35% • 10%
1	B	323	2% 52% 39% • 7%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 4536 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 Fructose-bisphosphate aldolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	290	2211	1403	379	412	17	0	0	0
1	B	302	2303	1461	395	430	17	0	0	0

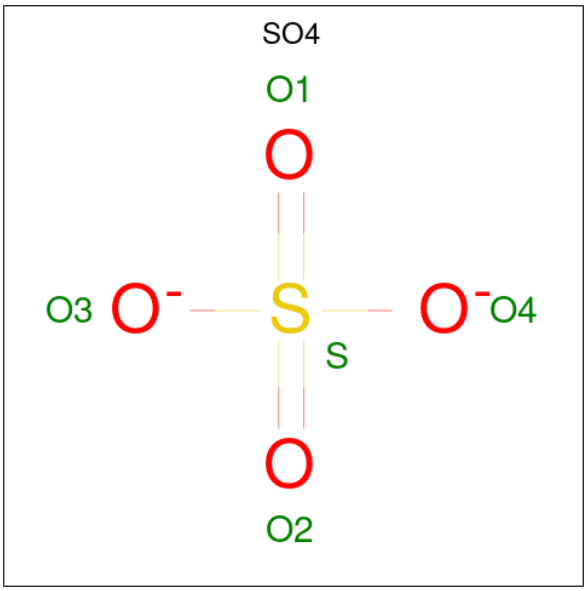
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	129	GLY	SER	SEE REMARK 999	UNP O97447
B	129	GLY	SER	SEE REMARK 999	UNP O97447

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Zn		
2	A	1	1	1	0	0
2	B	1	1	1	0	0

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



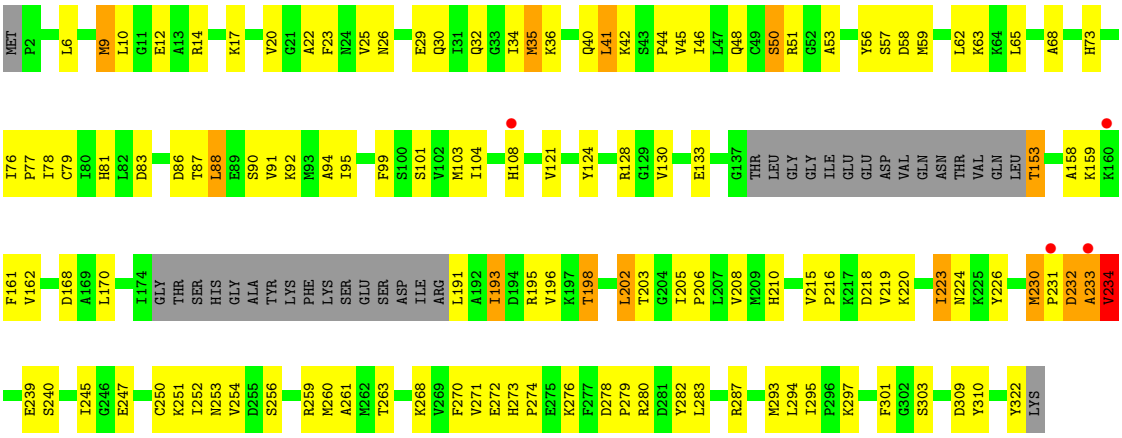
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0

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.

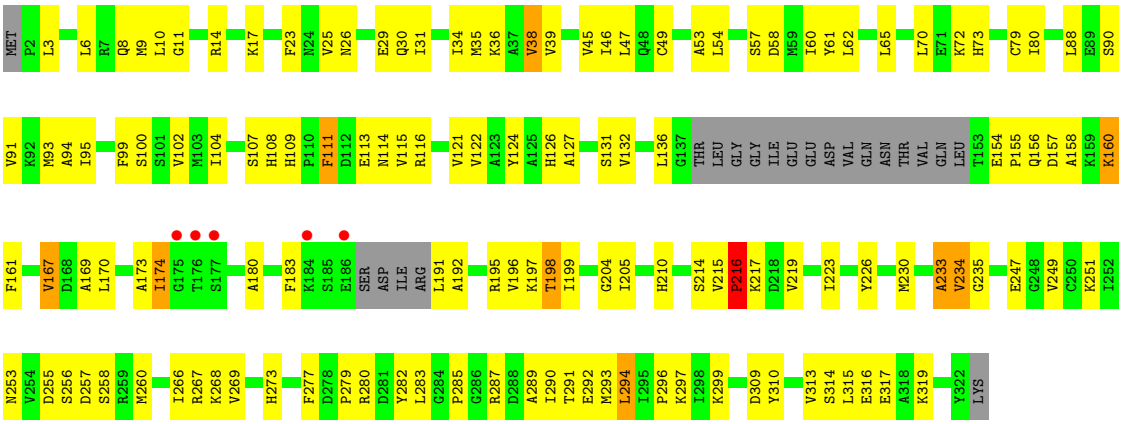
• Molecule 1: Fructose-bisphosphate aldolase

Chain A:



• Molecule 1: Fructose-bisphosphate aldolase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α, β, γ	62.92Å 62.92Å 318.59Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	10.00 – 2.90 9.97 – 2.90	Depositor EDS
% Data completeness (in resolution range)	86.5 (10.00-2.90) 86.5 (9.97-2.90)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.96 (at 2.89Å)	Xtriage
Refinement program	REFMAC 5.0	Depositor
R, R_{free}	0.220 , 0.286 0.216 , 0.281	Depositor DCC
R_{free} test set	1072 reflections (8.07%)	wwPDB-VP
Wilson B-factor (Å^2)	57.1	Xtriage
Anisotropy	0.271	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 53.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.53$, $< L^2 > = 0.37$	Xtriage
Estimated twinning fraction	0.069 for h,-h-k,-l	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	4536	wwPDB-VP
Average B, all atoms (Å^2)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.92% 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: ZN, 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	0.73	0/2250	0.82	2/3034 (0.1%)
1	B	0.73	0/2345	0.85	2/3160 (0.1%)
All	All	0.73	0/4595	0.84	4/6194 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
All	All	0	2

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	173	ALA	O-C-N	-5.70	113.58	122.70
1	A	309	ASP	CB-CG-OD1	5.60	123.34	118.30
1	B	233	ALA	N-CA-C	5.32	125.38	111.00
1	A	41	LEU	CB-CG-CD1	-5.18	102.19	111.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	233	ALA	Peptide
1	B	216	PRO	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	2211	0	2253	118	0
1	B	2303	0	2338	103	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	15	0	0	1	0
3	B	5	0	0	1	0
All	All	4536	0	4591	214	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 (214) 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:B:174:ILE:O	1:B:191:LEU:HD12	1.51	1.10
1:A:223:ILE:HG21	1:A:230:MET:HE3	1.36	1.07
1:B:111:PHE:O	1:B:115:VAL:HG23	1.61	1.00
1:B:183:PHE:HD2	1:B:234:VAL:HG11	1.28	0.96
1:A:57:SER:CB	1:A:62:LEU:HD13	1.97	0.94
1:B:38:VAL:HG11	1:B:45:VAL:HB	1.54	0.90
1:B:36:LYS:HD2	1:B:291:THR:HG21	1.54	0.87
1:B:94:ALA:O	1:B:99:PHE:HB2	1.78	0.83
1:B:183:PHE:HD2	1:B:234:VAL:CG1	1.92	0.82
1:B:174:ILE:O	1:B:191:LEU:CD1	2.29	0.79
1:A:198:THR:O	1:A:202:LEU:HD22	1.83	0.79
1:A:95:ILE:HG21	1:A:128:ARG:HD3	1.65	0.78
1:A:256:SER:HA	1:A:259:ARG:HD2	1.64	0.78
1:B:183:PHE:CD2	1:B:234:VAL:HG11	2.18	0.78
1:B:234:VAL:HG13	1:B:235:GLY:N	2.01	0.76
1:A:26:ASN:O	1:A:62:LEU:HD11	1.86	0.76
1:A:36:LYS:O	1:A:40:GLN:HB2	1.85	0.75
1:B:223:ILE:HG21	1:B:230:MET:HE2	1.69	0.73
1:B:279:PRO:HA	1:B:282:TYR:CZ	2.25	0.71
1:A:23:PHE:HB3	1:A:34:ILE:HD13	1.73	0.71
1:A:35:MET:HG3	1:A:73:HIS:CD2	2.25	0.70

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:215:VAL:HG13	1:B:260:MET:SD	2.32	0.69
1:B:49:CYS:SG	1:B:62:LEU:HD21	2.33	0.67
1:A:153:THR:O	1:A:195:ARG:NH1	2.26	0.67
1:A:219:VAL:CG1	1:A:260:MET:HB3	2.24	0.67
1:A:153:THR:O	1:A:195:ARG:HD3	1.93	0.67
1:A:88:LEU:O	1:A:91:VAL:HG22	1.94	0.66
1:A:191:LEU:HD23	1:A:193:ILE:CG1	2.24	0.66
1:A:282:TYR:CE1	1:A:283:LEU:HD12	2.31	0.66
1:A:191:LEU:HD23	1:A:193:ILE:HG12	1.78	0.66
1:B:282:TYR:C	1:B:285:PRO:HD2	2.15	0.65
1:B:234:VAL:CG1	1:B:235:GLY:N	2.60	0.65
1:A:50:SER:HA	1:A:83:ASP:HB3	1.80	0.64
1:A:263:THR:O	1:A:263:THR:HG22	1.97	0.63
1:A:57:SER:OG	1:A:62:LEU:HD13	1.98	0.63
1:A:68:ALA:HB1	1:B:60:ILE:HG13	1.80	0.63
1:B:169:ALA:HA	1:B:205:ILE:HD12	1.81	0.63
1:A:159:LYS:HD2	1:A:202:LEU:HB3	1.81	0.62
1:A:208:VAL:HG13	1:A:251:LYS:HG2	1.79	0.62
1:A:44:PRO:HD2	1:A:303:SER:HB2	1.82	0.61
1:A:233:ALA:C	1:A:234:VAL:CG2	2.69	0.61
1:B:282:TYR:O	1:B:285:PRO:HD2	2.01	0.61
1:A:226:TYR:CZ	1:A:268:LYS:HG3	2.35	0.60
1:A:208:VAL:HG13	1:A:251:LYS:HB3	1.82	0.60
1:B:195:ARG:O	1:B:199:ILE:HG13	2.00	0.60
1:A:12:GLU:OE2	1:A:310:TYR:CD2	2.53	0.60
1:A:245:ILE:HD11	1:A:252:ILE:HD11	1.83	0.60
1:A:23:PHE:CB	1:A:34:ILE:HD13	2.31	0.60
1:B:107:SER:HA	1:B:114:ASN:ND2	2.19	0.58
1:B:91:VAL:O	1:B:95:ILE:HG13	2.03	0.58
1:A:104:ILE:HB	1:A:121:VAL:HG21	1.85	0.58
1:A:158:ALA:O	1:A:162:VAL:HG23	2.04	0.57
1:A:81:HIS:HA	1:A:101:SER:O	2.03	0.57
1:A:259:ARG:NH2	3:A:328:SO4:O4	2.37	0.57
1:A:233:ALA:HA	1:A:234:VAL:HG22	1.87	0.57
1:A:245:ILE:HD11	1:A:252:ILE:CD1	2.34	0.57
1:B:154:GLU:O	1:B:157:ASP:HB2	2.05	0.57
1:A:297:LYS:HG2	1:A:301:PHE:CZ	2.40	0.56
1:B:14:ARG:NH2	1:B:204:GLY:HA2	2.19	0.56
1:A:92:LYS:HG3	1:A:124:TYR:CZ	2.40	0.56
1:B:122:VAL:HG13	1:B:126:HIS:CE1	2.40	0.56
1:A:57:SER:HB3	1:A:62:LEU:HD13	1.87	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:273:HIS:HB3	1:A:276:LYS:HG3	1.88	0.56
1:B:26:ASN:OD1	1:B:53:ALA:HB2	2.05	0.56
1:A:58:ASP:O	1:A:59:MET:HB2	2.04	0.56
1:A:278:ASP:OD1	1:A:280:ARG:N	2.38	0.56
1:B:196:VAL:HG11	1:B:249:VAL:HG23	1.86	0.56
1:A:57:SER:HB2	1:A:62:LEU:HD13	1.87	0.56
1:A:29:GLU:O	1:A:287:ARG:HG3	2.06	0.55
1:B:279:PRO:HA	1:B:282:TYR:CE2	2.41	0.55
1:A:12:GLU:OE2	1:A:310:TYR:HD2	1.90	0.55
1:B:23:PHE:CE1	1:B:294:LEU:HD11	2.41	0.55
1:A:233:ALA:HA	1:A:234:VAL:CG2	2.36	0.55
1:B:223:ILE:HG21	1:B:230:MET:CE	2.35	0.55
1:B:257:ASP:OD2	1:B:297:LYS:NZ	2.39	0.55
1:A:68:ALA:CB	1:B:60:ILE:HG13	2.37	0.55
1:B:183:PHE:CD2	1:B:234:VAL:CG1	2.83	0.55
1:A:282:TYR:CD1	1:A:283:LEU:HD12	2.43	0.54
1:B:10:LEU:N	1:B:10:LEU:HD23	2.22	0.54
1:A:91:VAL:HG23	1:A:92:LYS:N	2.22	0.54
1:B:313:VAL:HG12	1:B:314:SER:O	2.07	0.54
1:A:6:LEU:HD11	1:A:10:LEU:HD11	1.88	0.54
1:A:203:THR:O	1:A:205:ILE:HG23	2.08	0.54
1:B:25:VAL:HG11	1:B:31:ILE:HG12	1.90	0.54
1:A:208:VAL:HG13	1:A:251:LYS:CG	2.38	0.54
1:A:196:VAL:HG12	1:A:247:GLU:HB2	1.90	0.53
1:A:233:ALA:CA	1:A:234:VAL:HG22	2.38	0.53
1:B:289:ALA:O	1:B:290:ILE:C	2.45	0.53
1:B:113:GLU:OE1	1:B:116:ARG:NH2	2.30	0.53
1:A:282:TYR:CE1	1:A:283:LEU:CD1	2.92	0.53
1:A:198:THR:C	1:A:202:LEU:HD22	2.29	0.52
1:A:233:ALA:C	1:A:234:VAL:HG23	2.29	0.52
1:A:191:LEU:HD23	1:A:193:ILE:HG13	1.91	0.51
1:A:224:ASN:HD21	1:A:231:PRO:HA	1.75	0.51
1:B:6:LEU:HG	1:B:10:LEU:HG	1.92	0.51
1:A:51:ARG:NH2	1:A:86:ASP:OD2	2.43	0.51
1:A:159:LYS:CD	1:A:202:LEU:HB3	2.41	0.51
1:B:214:SER:HB2	1:B:216:PRO:HD3	1.93	0.51
1:A:245:ILE:CD1	1:A:252:ILE:HD11	2.41	0.51
1:B:102:VAL:HG23	1:B:132:VAL:HG13	1.92	0.50
1:A:30:GLN:O	1:A:34:ILE:HG13	2.11	0.50
1:B:124:TYR:O	1:B:127:ALA:HB3	2.11	0.50
1:A:86:ASP:OD1	1:A:87:THR:N	2.41	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:47:LEU:HD12	1:B:80:ILE:HD13	1.94	0.50
1:A:215:VAL:N	1:A:216:PRO:HD3	2.27	0.50
1:B:180:ALA:O	1:B:233:ALA:HA	2.11	0.49
1:A:65:LEU:HD21	1:B:65:LEU:HD21	1.93	0.49
1:A:196:VAL:CG1	1:A:247:GLU:HB2	2.43	0.49
1:B:226:TYR:CE2	1:B:268:LYS:HG3	2.47	0.49
1:B:26:ASN:OD1	1:B:53:ALA:CB	2.60	0.49
1:B:108:HIS:CD2	1:B:108:HIS:H	2.31	0.49
1:A:29:GLU:OE2	1:A:283:LEU:HD23	2.12	0.49
1:A:210:HIS:CD2	1:A:253:ASN:ND2	2.81	0.49
1:B:9:MET:HG2	1:B:10:LEU:HD23	1.94	0.48
1:A:56:TYR:CE1	1:B:280:ARG:HG2	2.48	0.48
1:B:35:MET:O	1:B:39:VAL:HG23	2.13	0.48
1:A:196:VAL:HG12	1:A:247:GLU:CB	2.43	0.48
1:A:208:VAL:HG13	1:A:251:LYS:CB	2.43	0.48
1:A:25:VAL:HG22	1:A:48:GLN:O	2.14	0.48
1:A:91:VAL:CG2	1:A:92:LYS:N	2.76	0.48
1:B:266:ILE:HG12	1:B:282:TYR:HD1	1.78	0.48
1:B:57:SER:OG	1:B:62:LEU:HB2	2.14	0.48
1:A:223:ILE:HG21	1:A:230:MET:CE	2.25	0.47
1:B:156:GLN:HB3	1:B:160:LYS:NZ	2.28	0.47
1:B:30:GLN:O	1:B:31:ILE:C	2.52	0.47
1:A:161:PHE:HE2	1:A:170:LEU:HD13	1.79	0.47
1:A:293:MET:HE3	1:A:294:LEU:HD12	1.96	0.47
1:B:309:ASP:HB2	1:B:310:TYR:CD2	2.49	0.47
1:A:35:MET:HG3	1:A:73:HIS:HD2	1.77	0.47
1:A:270:PHE:HE2	1:B:266:ILE:HG22	1.79	0.47
1:A:233:ALA:CA	1:A:234:VAL:CG2	2.93	0.47
1:B:31:ILE:O	1:B:35:MET:HG2	2.15	0.47
1:B:113:GLU:CD	1:B:116:ARG:HH21	2.17	0.46
1:B:170:LEU:HB2	1:B:205:ILE:HD11	1.96	0.46
1:A:94:ALA:O	1:A:99:PHE:HB2	2.15	0.46
1:A:104:ILE:HG23	1:A:104:ILE:O	2.15	0.46
1:B:316:GLU:O	1:B:319:LYS:HB2	2.16	0.46
1:A:32:GLN:HB2	1:A:287:ARG:HD2	1.97	0.46
1:B:219:VAL:CG1	1:B:260:MET:HB3	2.45	0.46
1:B:155:PRO:O	1:B:158:ALA:HB3	2.16	0.46
1:A:44:PRO:HB3	1:A:77:PRO:HG2	1.98	0.46
1:A:279:PRO:O	1:A:283:LEU:HB2	2.15	0.46
1:A:270:PHE:CE2	1:B:266:ILE:HG22	2.51	0.46
1:B:313:VAL:HG13	1:B:317:GLU:HB2	1.98	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:72:LYS:HG2	1:B:73:HIS:NE2	2.31	0.45
1:B:54:LEU:HD23	1:B:54:LEU:HA	1.66	0.45
1:A:168:ASP:OD2	1:A:322:TYR:OH	2.30	0.45
1:A:232:ASP:HB2	1:A:233:ALA:H	1.57	0.45
1:B:217:LYS:O	1:B:217:LYS:HG2	2.15	0.45
1:B:293:MET:SD	1:B:293:MET:C	2.95	0.45
1:B:46:ILE:CD1	1:B:251:LYS:HG3	2.47	0.45
1:A:62:LEU:O	1:A:63:LYS:C	2.54	0.45
1:A:205:ILE:HB	1:A:206:PRO:CD	2.47	0.45
1:B:109:HIS:HB3	1:B:113:GLU:HB3	1.99	0.45
1:A:23:PHE:HB3	1:A:34:ILE:CD1	2.45	0.45
1:A:95:ILE:HD13	1:A:124:TYR:CD2	2.52	0.45
1:A:104:ILE:CB	1:A:121:VAL:HG21	2.47	0.44
1:B:90:SER:O	1:B:93:MET:HB3	2.17	0.44
1:A:103:MET:HA	1:A:133:GLU:O	2.17	0.44
1:B:54:LEU:HD23	1:B:62:LEU:HD23	1.98	0.44
1:B:223:ILE:HD13	1:B:230:MET:HE1	1.98	0.44
1:B:226:TYR:HB3	1:B:267:ARG:HB3	2.00	0.44
1:B:136:LEU:HD13	1:B:161:PHE:CD1	2.53	0.44
1:A:9:MET:SD	1:A:20:VAL:HG11	2.58	0.44
1:A:32:GLN:CB	1:A:287:ARG:HD2	2.47	0.44
1:B:113:GLU:OE2	1:B:113:GLU:HA	2.16	0.44
1:A:208:VAL:CG1	1:A:251:LYS:HG2	2.48	0.43
1:A:293:MET:CE	1:A:294:LEU:HD12	2.48	0.43
1:A:218:ASP:OD2	1:A:219:VAL:N	2.51	0.43
1:B:29:GLU:O	1:B:287:ARG:HG3	2.18	0.43
1:A:263:THR:O	1:A:263:THR:CG2	2.66	0.43
1:B:14:ARG:HH22	1:B:204:GLY:HA2	1.83	0.43
1:B:223:ILE:CD1	1:B:230:MET:HE1	2.49	0.43
1:A:76:ILE:HG22	1:A:78:ILE:HG13	2.01	0.43
1:B:3:LEU:HB2	1:B:70:LEU:CD2	2.49	0.43
1:B:223:ILE:HD13	1:B:230:MET:CE	2.48	0.43
1:B:61:TYR:O	1:B:65:LEU:HG	2.19	0.43
1:A:87:THR:O	1:A:90:SER:OG	2.36	0.43
1:A:94:ALA:O	1:A:95:ILE:C	2.57	0.43
1:B:197:LYS:O	1:B:198:THR:C	2.55	0.42
1:B:156:GLN:O	1:B:157:ASP:C	2.55	0.42
1:A:95:ILE:HG12	1:A:130:VAL:HG21	2.01	0.42
1:B:34:ILE:O	1:B:38:VAL:HG22	2.19	0.42
1:A:103:MET:HE3	1:A:210:HIS:HE1	1.84	0.42
1:A:219:VAL:HG11	1:A:260:MET:HB3	1.99	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:230:MET:CE	1:B:277:PHE:HB3	2.50	0.42
1:A:278:ASP:OD1	1:A:278:ASP:C	2.58	0.42
1:A:50:SER:OG	1:A:53:ALA:N	2.46	0.42
1:A:271:VAL:O	1:A:274:PRO:HD3	2.20	0.42
1:B:210:HIS:HA	1:B:253:ASN:HB2	2.02	0.42
1:A:124:TYR:CZ	1:A:128:ARG:CZ	3.03	0.42
1:B:10:LEU:HD11	1:B:251:LYS:HB2	2.02	0.41
1:B:266:ILE:HG12	1:B:282:TYR:CD1	2.54	0.41
1:B:292:GLU:O	1:B:296:PRO:HD2	2.20	0.41
1:A:9:MET:SD	1:A:44:PRO:HB2	2.60	0.41
1:A:254:VAL:HG21	1:A:297:LYS:HD3	2.02	0.41
1:B:136:LEU:HD21	1:B:157:ASP:HB3	2.03	0.41
1:B:255:ASP:HB3	3:B:329:SO4:O1	2.20	0.41
1:A:14:ARG:HB2	1:A:250:CYS:SG	2.61	0.41
1:A:22:ALA:HA	1:A:46:ILE:HB	2.03	0.41
1:B:196:VAL:HG12	1:B:247:GLU:O	2.21	0.41
1:B:256:SER:O	1:B:257:ASP:C	2.59	0.41
1:A:40:GLN:OE1	1:A:295:ILE:HD11	2.21	0.41
1:B:174:ILE:HG21	1:B:192:ALA:O	2.20	0.41
1:B:100:SER:HB2	1:B:315:LEU:HD21	2.02	0.40
1:B:104:ILE:HD13	1:B:121:VAL:HG21	2.02	0.40
1:A:41:LEU:O	1:A:42:LYS:C	2.59	0.40
1:A:45:VAL:CG1	1:A:78:ILE:HG12	2.51	0.40
1:B:122:VAL:HG13	1:B:126:HIS:ND1	2.37	0.40
1:B:223:ILE:HD11	1:B:260:MET:HA	2.02	0.40
1:B:234:VAL:HG13	1:B:235:GLY:H	1.82	0.40
1:A:205:ILE:HB	1:A:206:PRO:HD2	2.03	0.40
1:B:8:GLN:O	1:B:11:GLY:N	2.54	0.40
1:B:269:VAL:HG11	1:B:282:TYR:HB3	2.03	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	284/323 (88%)	243 (86%)	38 (13%)	3 (1%)	14	42
1	B	296/323 (92%)	260 (88%)	33 (11%)	3 (1%)	15	45
All	All	580/646 (90%)	503 (87%)	71 (12%)	6 (1%)	15	45

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	261	ALA
1	B	216	PRO
1	B	111	PHE
1	B	167	VAL
1	A	234	VAL
1	A	223	ILE

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	240/268 (90%)	222 (92%)	18 (8%)	13	37
1	B	249/268 (93%)	233 (94%)	16 (6%)	17	45
All	All	489/536 (91%)	455 (93%)	34 (7%)	15	41

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	MET
1	A	17	LYS
1	A	35	MET
1	A	50	SER
1	A	79	CYS
1	A	88	LEU
1	A	108	HIS
1	A	153	THR
1	A	193	ILE
1	A	198	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	202	LEU
1	A	220	LYS
1	A	230	MET
1	A	232	ASP
1	A	234	VAL
1	A	239	GLU
1	A	240	SER
1	A	272	GLU
1	B	17	LYS
1	B	38	VAL
1	B	58	ASP
1	B	79	CYS
1	B	88	LEU
1	B	131	SER
1	B	160	LYS
1	B	167	VAL
1	B	174	ILE
1	B	198	THR
1	B	234	VAL
1	B	258	SER
1	B	273	HIS
1	B	283	LEU
1	B	294	LEU
1	B	299	LYS

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

Mol	Chain	Res	Type
1	A	73	HIS
1	A	109	HIS
1	A	156	GLN
1	A	224	ASN
1	B	108	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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	$\#\|Z\| > 2$
3	SO4	A	327	-	4,4,4	0.57	0	6,6,6	0.52	0
3	SO4	A	328	-	4,4,4	0.55	0	6,6,6	0.39	0
3	SO4	B	329	-	4,4,4	0.36	0	6,6,6	0.30	0
3	SO4	A	330	-	4,4,4	0.50	0	6,6,6	0.84	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.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	328	SO4	1	0
3	B	329	SO4	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	290/323 (89%)	-0.44	4 (1%)	75	75	25, 46, 73, 95	0
1	B	302/323 (93%)	-0.47	5 (1%)	70	69	19, 44, 69, 95	0
All	All	592/646 (91%)	-0.45	9 (1%)	73	73	19, 45, 73, 95	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	108	HIS	3.8
1	B	184	LYS	3.2
1	B	176	THR	2.8
1	A	160	LYS	2.6
1	B	177	SER	2.5
1	A	231	PRO	2.5
1	B	175	GLY	2.3
1	B	186	GLU	2.2
1	A	233	ALA	2.2

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
3	SO4	A	330	5/5	0.96	0.23	44,44,44,44	0
3	SO4	A	327	5/5	0.97	0.16	44,44,44,44	0
2	ZN	B	332	1/1	0.97	0.17	44,44,44,44	0
3	SO4	A	328	5/5	0.98	0.23	44,44,44,44	0
2	ZN	A	331	1/1	0.98	0.27	44,44,44,44	0
3	SO4	B	329	5/5	0.98	0.13	44,44,44,44	0

6.5 Other polymers (i)

There are no such residues in this entry.