Full wwPDB X-ray Structure Validation Report (i)

Sep 6, 2023 – 02:02 AM EDT

PDB ID	:	3VFL
Title	:	Structure, Function, Stability and Knockout Phenotype of Dihydrodipicolinate Synthase from Streptococcus pneumoniae
Authors	:	Gorman, M.A.
Deposited on	:	2012-01-10
Resolution	:	1.91 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:

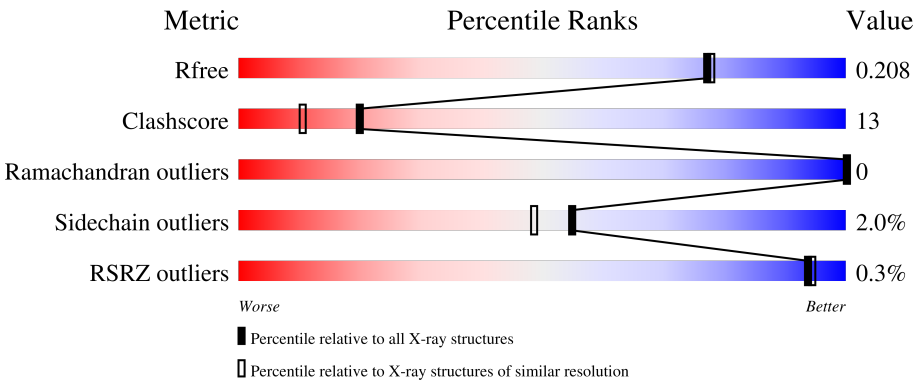
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, 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 1.91 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	7937 (1.94-1.90)
Clashscore	141614	8644 (1.94-1.90)
Ramachandran outliers	138981	8530 (1.94-1.90)
Sidechain outliers	138945	8530 (1.94-1.90)
RSRZ outliers	127900	7793 (1.94-1.90)

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	311	% 76% 22% ••
1	B	311	76% 22% ••

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	403	-	-	X	-
3	GOL	A	405	-	X	X	-
3	GOL	B	403	-	-	X	-

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 5014 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.

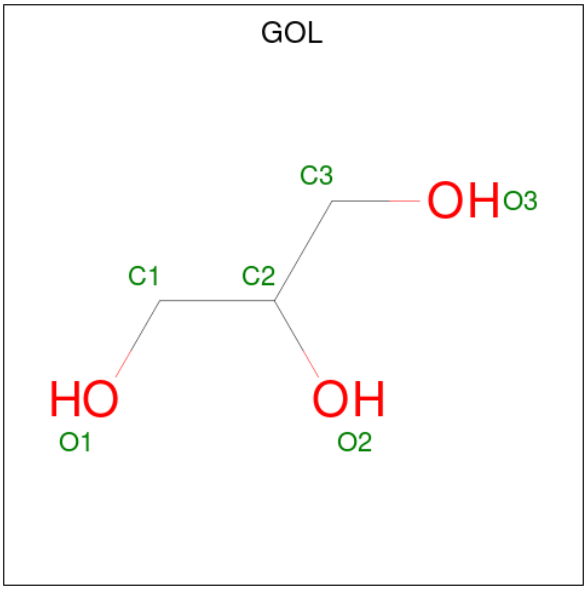
- Molecule 1 is a protein called Dihydrodipicolinate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	308	2366	1516	385	455	10	0	0	0
1	B	306	2359	1512	383	454	10	0	1	0

- Molecule 2 is POTASSIUM ION (three-letter code: K) (formula: K).

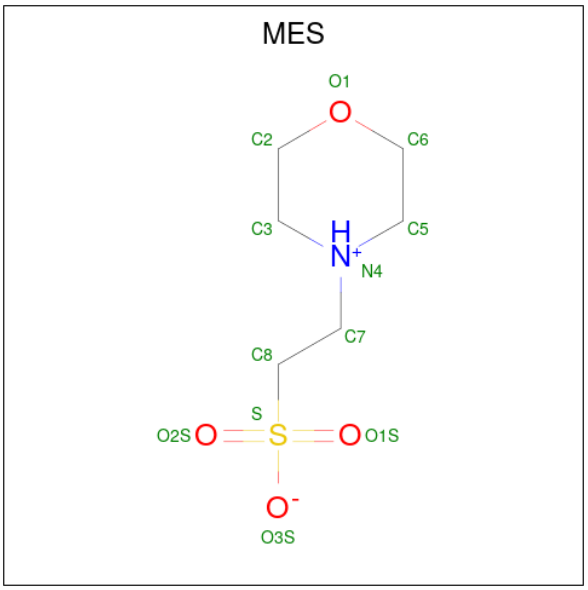
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	K		
2	A	2	2	2	0	0
2	B	2	2	2	0	0

- 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
			Total	C	O			
3	A	1	6	3	3		0	0
			Total	C	O			
3	A	1	6	3	3		0	0
			Total	C	O			
3	B	1	6	3	3		0	0

- Molecule 4 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: $C_6H_{13}NO_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
4	A	1	12	6	1	4	1	0	0
			Total	C	N	O	S		
4	B	1	12	6	1	4	1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	119	119	119	0	0
			Total	O		
5	B	118	118	118	0	0

3 Residue-property plots (i)

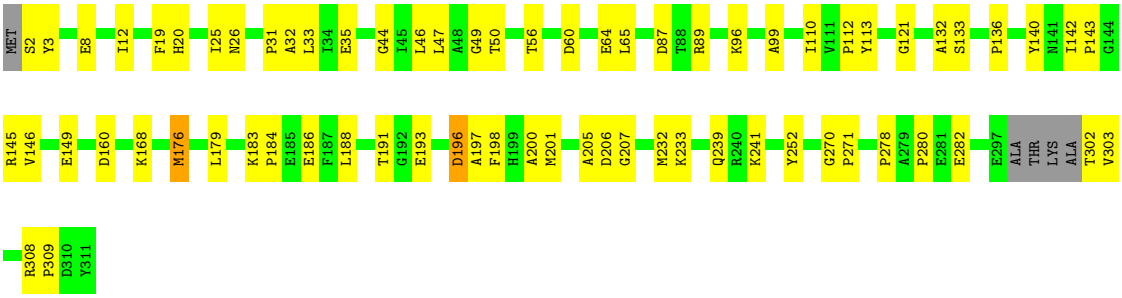
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Dihydrodipicolinate synthase

Chain A:

• Molecule 1: Dihydrodipicolinate synthase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, α, β, γ	62.36Å 105.33Å 105.51Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.83 – 1.91 47.83 – 1.91	Depositor EDS
% Data completeness (in resolution range)	99.1 (47.83-1.91) 99.3 (47.83-1.91)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.73 (at 1.91Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.165 , 0.212 0.166 , 0.208	Depositor DCC
R_{free} test set	2747 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å^2)	26.2	Xtriage
Anisotropy	0.658	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 16.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.40$, $< L^2 > = 0.22$	Xtriage
Estimated twinning fraction	0.459 for -h,l,k	Xtriage
Reported twinning fraction	0.541 for H, K, L 0.459 for -H, L, K	Depositor
Outliers	0 of 54037 reflections	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	5014	wwPDB-VP
Average B, all atoms (Å^2)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.83% 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: K, GOL, MES

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	0.99	0/2420	1.13	9/3293 (0.3%)
1	B	1.01	1/2416 (0.0%)	1.07	8/3287 (0.2%)
All	All	1.00	1/4836 (0.0%)	1.10	17/6580 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	193	GLU	CD-OE1	5.83	1.32	1.25

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	196[A]	ASP	CB-CG-OD2	7.86	125.37	118.30
1	B	196[B]	ASP	CB-CG-OD2	7.86	125.37	118.30
1	B	160	ASP	CB-CG-OD1	-7.83	111.25	118.30
1	B	33	LEU	CB-CG-CD1	-7.13	98.89	111.00
1	A	22	ASP	CB-CG-OD1	-7.09	111.92	118.30
1	A	196	ASP	CB-CA-C	-6.38	97.63	110.40
1	A	203	LEU	CB-CG-CD2	-6.22	100.43	111.00
1	A	60	ASP	CB-CG-OD2	6.13	123.82	118.30
1	B	188	LEU	CB-CG-CD2	-6.10	100.63	111.00
1	B	206	ASP	CB-CG-OD1	-6.04	112.87	118.30
1	A	131	ASP	CB-CG-OD1	6.02	123.72	118.30
1	A	274	LEU	CB-CG-CD1	-6.00	100.79	111.00
1	A	63	LEU	CB-CG-CD1	-5.85	101.06	111.00
1	A	22	ASP	N-CA-CB	-5.78	100.19	110.60
1	A	76	ARG	NE-CZ-NH1	-5.66	117.47	120.30
1	B	252	TYR	N-CA-C	-5.23	96.88	111.00
1	B	241	LYS	CD-CE-NZ	-5.12	99.93	111.70

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	2366	0	2336	58	0
1	B	2359	0	2328	66	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	18	0	24	17	0
3	B	6	0	8	16	0
4	A	12	0	13	1	0
4	B	12	0	13	1	0
5	A	119	0	0	4	0
5	B	118	0	0	6	0
All	All	5014	0	4722	120	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 13.

All (120) 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:49:GLY:CA	3:B:403:GOL:H31	1.83	1.07
1:B:50:THR:H	3:B:403:GOL:C3	1.67	1.07
1:B:50:THR:H	3:B:403:GOL:H32	1.19	1.02
1:B:50:THR:N	3:B:403:GOL:H32	1.78	0.97
1:B:49:GLY:HA2	3:B:403:GOL:H31	1.48	0.96
1:A:308:ARG:NH1	1:B:89:ARG:HG2	1.92	0.84
1:B:50:THR:N	3:B:403:GOL:C3	2.36	0.83
1:A:168:LYS:NZ	3:A:403:GOL:H2	1.99	0.77
1:A:168:LYS:HZ3	3:A:403:GOL:H2	1.49	0.77
1:B:176:MET:HE2	1:B:200:ALA:HB1	1.66	0.77
3:A:405:GOL:H2	1:B:89:ARG:HB3	1.68	0.76
1:A:56:THR:HB	3:A:405:GOL:H31	1.68	0.75
1:B:176:MET:HE3	1:B:205:ALA:HB2	1.70	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:176:MET:HE2	1:B:200:ALA:CB	2.18	0.73
1:B:176:MET:CE	1:B:200:ALA:HB1	2.20	0.72
1:B:233:LYS:HG3	5:B:530:HOH:O	1.88	0.72
1:A:168:LYS:NZ	3:A:403:GOL:C2	2.53	0.72
1:A:72:VAL:HG12	1:A:72:VAL:O	1.90	0.70
1:B:96:LYS:HE3	1:B:132:ALA:O	1.90	0.70
3:A:405:GOL:H11	1:B:87:ASP:OD1	1.92	0.70
1:A:182:HIS:HD2	5:A:580:HOH:O	1.75	0.68
1:B:302:THR:OG1	1:B:303:VAL:N	2.24	0.68
1:B:49:GLY:HA3	3:B:403:GOL:H12	1.76	0.67
1:A:182:HIS:HE1	5:A:525:HOH:O	1.77	0.66
3:A:405:GOL:H2	1:B:89:ARG:CB	2.25	0.66
1:B:50:THR:H	3:B:403:GOL:C2	2.08	0.65
1:A:20:HIS:C	1:A:22:ASP:H	2.01	0.64
1:B:278:PRO:O	1:B:303:VAL:HG11	1.98	0.64
1:B:282:GLU:H	1:B:282:GLU:CD	2.01	0.63
1:A:50:THR:N	3:A:403:GOL:H12	2.14	0.63
1:B:176:MET:CE	1:B:205:ALA:HB2	2.28	0.62
1:A:263:ASN:OD1	1:A:269:ALA:N	2.33	0.62
1:B:49:GLY:HA3	3:B:403:GOL:H31	1.75	0.62
1:A:72:VAL:O	1:A:72:VAL:CG1	2.47	0.62
1:A:193:GLU:OE1	4:A:406:MES:H21	2.00	0.62
1:A:58:THR:HG23	1:A:308:ARG:HE	1.65	0.61
1:A:140:TYR:CE1	1:A:168:LYS:HE2	2.37	0.59
1:A:46:LEU:HD22	3:A:403:GOL:O2	2.01	0.59
1:B:143:PRO:HB3	1:B:149:GLU:HB3	1.85	0.58
1:B:8:GLU:OE2	1:B:8:GLU:HA	2.04	0.57
1:A:223:MET:HB2	1:A:238:ILE:HG21	1.87	0.57
1:B:99:ALA:HB2	1:B:133:SER:OG	2.05	0.56
1:A:46:LEU:CD2	3:A:403:GOL:O2	2.53	0.56
3:A:405:GOL:C1	1:B:87:ASP:OD1	2.53	0.56
1:A:168:LYS:HZ2	3:A:403:GOL:C2	2.18	0.56
1:A:119:GLN:HE22	1:A:150:LEU:HA	1.71	0.55
1:A:56:THR:HB	3:A:405:GOL:H11	1.88	0.54
1:B:110:ILE:HA	1:B:140:TYR:HB3	1.90	0.54
1:A:58:THR:O	1:A:62:GLU:HG3	2.07	0.53
1:A:308:ARG:HG3	1:A:308:ARG:HH11	1.72	0.53
1:B:271:PRO:HA	5:B:517:HOH:O	2.09	0.53
1:B:49:GLY:CA	3:B:403:GOL:C3	2.74	0.52
1:B:184:PRO:HD3	5:B:562:HOH:O	2.12	0.50
1:B:280:PRO:HA	1:B:303:VAL:HG21	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:88:THR:O	1:A:92:ILE:HG12	2.12	0.50
1:B:113:TYR:HB3	1:B:146:VAL:HG12	1.94	0.50
1:A:191:THR:HG21	1:A:197:ALA:HA	1.93	0.49
1:A:92:ILE:HD11	1:A:128:ALA:HB1	1.93	0.49
1:B:50:THR:OG1	3:B:403:GOL:H32	2.13	0.49
1:A:41:HIS:HE1	5:A:568:HOH:O	1.94	0.49
1:A:289:LYS:O	1:A:293:ASP:HA	2.13	0.49
1:A:277:VAL:HG13	1:B:121:GLY:HA2	1.94	0.49
1:B:198:PHE:CG	1:B:239:GLN:HG2	2.47	0.49
1:A:80:ILE:HG12	1:A:106:ALA:HB3	1.94	0.49
1:B:198:PHE:CD2	1:B:239:GLN:HG2	2.48	0.48
1:A:20:HIS:C	1:A:22:ASP:N	2.63	0.48
1:A:74:ASN:CG	1:A:74:ASN:O	2.52	0.48
1:B:142:ILE:HG12	1:B:145:ARG:HB2	1.93	0.48
1:A:58:THR:CG2	1:A:308:ARG:HE	2.25	0.48
1:B:201:MET:SD	1:B:207:GLY:HA2	2.54	0.48
1:A:140:TYR:OH	1:A:142:ILE:HD13	2.14	0.48
1:B:308:ARG:NH1	5:B:609:HOH:O	2.42	0.47
1:B:133:SER:HB3	5:B:593:HOH:O	2.13	0.47
1:B:46:LEU:CD2	3:B:403:GOL:H11	2.45	0.47
1:B:60:ASP:O	1:B:64:GLU:HG3	2.15	0.47
1:A:26:ASN:OD1	1:A:28:ASP:HB2	2.15	0.46
1:B:140:TYR:OH	4:B:404:MES:O2S	2.34	0.46
1:A:50:THR:H	3:A:403:GOL:H12	1.80	0.46
1:A:116:LYS:HD3	1:B:145:ARG:HG2	1.97	0.46
1:B:25:ILE:HD13	1:B:65:LEU:HD11	1.98	0.46
1:A:142:ILE:HG12	1:A:145:ARG:HB2	1.99	0.45
1:A:110:ILE:HA	1:A:140:TYR:HB3	1.97	0.45
1:B:191:THR:HG21	1:B:197:ALA:HA	1.98	0.45
1:A:110:ILE:HB	3:A:404:GOL:H11	1.98	0.45
1:B:142:ILE:O	1:B:146:VAL:HG22	2.17	0.45
1:A:96:LYS:HA	1:A:133:SER:HB2	1.98	0.44
1:A:87:ASP:OD1	1:B:56:THR:HA	2.17	0.44
1:A:161:HIS:HB3	1:A:164:ILE:HG13	2.00	0.44
1:B:140:TYR:CD1	1:B:168:LYS:HD3	2.53	0.44
1:B:3:TYR:CE1	1:B:186:GLU:HA	2.53	0.44
1:B:46:LEU:HD23	3:B:403:GOL:H11	2.00	0.44
1:A:23:GLY:CA	1:A:306:VAL:CG1	2.96	0.44
1:B:12:ILE:HA	1:B:44:GLY:O	2.17	0.44
1:B:49:GLY:C	3:B:403:GOL:H31	2.35	0.44
1:B:3:TYR:CZ	1:B:186:GLU:HA	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:31:PRO:O	1:B:35:GLU:HG2	2.19	0.43
1:A:19:PHE:O	1:A:270:GLY:HA3	2.19	0.43
1:A:56:THR:O	3:A:405:GOL:C3	2.66	0.42
1:A:113:TYR:HB3	1:A:146:VAL:HG12	2.02	0.42
1:A:252:TYR:CD2	1:A:253:PRO:HD2	2.54	0.42
1:B:20:HIS:CE1	1:B:26:ASN:HA	2.54	0.42
1:A:140:TYR:CZ	1:A:168:LYS:HE2	2.55	0.41
1:B:19:PHE:O	1:B:270:GLY:HA3	2.20	0.41
1:B:50:THR:N	3:B:403:GOL:H12	2.34	0.41
1:A:43:ASP:O	1:A:78:PRO:HD2	2.19	0.41
1:A:259:LYS:HE3	5:A:511:HOH:O	2.18	0.41
1:B:32:ALA:N	5:B:532:HOH:O	2.41	0.41
1:A:108:LEU:HD21	1:A:140:TYR:HB2	2.02	0.41
1:A:119:GLN:NE2	1:A:151:THR:H	2.18	0.41
1:A:10:LYS:N	1:A:43:ASP:OD2	2.39	0.41
1:B:47:LEU:HA	1:B:47:LEU:HD23	1.85	0.41
1:B:49:GLY:CA	3:B:403:GOL:H12	2.49	0.41
1:A:6:LEU:HA	1:A:6:LEU:HD23	1.78	0.40
1:A:280:PRO:HA	1:A:303:VAL:HG21	2.03	0.40
1:B:179:LEU:O	1:B:183:LYS:HB3	2.22	0.40
1:A:273:ARG:CZ	1:A:308:ARG:HB3	2.51	0.40
3:A:405:GOL:H2	1:B:89:ARG:HB2	2.03	0.40
1:B:2:SER:HA	1:B:136:PRO:HG2	2.04	0.40
1:A:28:ASP:O	1:A:31:PRO:HD2	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	304/311 (98%)	290 (95%)	14 (5%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	303/311 (97%)	290 (96%)	13 (4%)	0	100	100
All	All	607/622 (98%)	580 (96%)	27 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	253/255 (99%)	247 (98%)	6 (2%)	49	41
1	B	253/255 (99%)	249 (98%)	4 (2%)	62	58
All	All	506/510 (99%)	496 (98%)	10 (2%)	55	49

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	ASP
1	A	108	LEU
1	A	133	SER
1	A	186	GLU
1	A	299	THR
1	A	309	PRO
1	B	112	PRO
1	B	176	MET
1	B	232	MET
1	B	309	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	41	HIS
1	A	74	ASN
1	A	119	GLN

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Mol	Chain	Res	Type
1	A	182	HIS
1	B	36	HIS
1	B	119	GLN
1	B	202	ASN
1	B	216	ASN

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 10 ligands modelled in this entry, 4 are monoatomic - leaving 6 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	$\#\|Z\| > 2$
4	MES	A	406	-	12,12,12	2.26	1 (8%)	14,16,16	6.31	10 (71%)
3	GOL	B	403	-	5,5,5	0.92	0	5,5,5	1.06	0
3	GOL	A	405	-	5,5,5	0.70	0	5,5,5	1.77	2 (40%)
4	MES	B	404	-	12,12,12	2.62	2 (16%)	14,16,16	6.82	8 (57%)
3	GOL	A	404	-	5,5,5	0.34	0	5,5,5	0.93	0
3	GOL	A	403	-	5,5,5	0.41	0	5,5,5	0.57	0

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
4	MES	A	406	-	-	5/6/14/14	0/1/1/1
3	GOL	B	403	-	-	4/4/4/4	-
3	GOL	A	405	-	-	4/4/4/4	-
4	MES	B	404	-	-	4/6/14/14	0/1/1/1
3	GOL	A	404	-	-	4/4/4/4	-
3	GOL	A	403	-	-	4/4/4/4	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	404	MES	C8-S	-8.53	1.65	1.77
4	A	406	MES	C8-S	-7.01	1.67	1.77
4	B	404	MES	O2S-S	2.00	1.50	1.45

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	B	404	MES	O1S-S-C8	15.43	125.50	106.92
4	B	404	MES	O3S-S-C8	-11.75	86.76	105.77
4	A	406	MES	O3S-S-O1S	-11.36	83.52	111.27
4	B	404	MES	O3S-S-O1S	-10.43	85.79	111.27
4	A	406	MES	O3S-S-O2S	-10.02	86.79	111.27
4	A	406	MES	O2S-S-C8	9.97	118.92	106.92
4	B	404	MES	O3S-S-O2S	-9.68	87.63	111.27
4	A	406	MES	O3S-S-C8	-9.35	90.64	105.77
4	A	406	MES	O1S-S-C8	9.34	118.17	106.92
4	B	404	MES	O2S-S-C8	5.44	113.46	106.92
4	A	406	MES	O1-C2-C3	4.25	121.16	111.80
4	B	404	MES	C6-C5-N4	-4.11	103.87	110.10
4	B	404	MES	C2-C3-N4	-3.64	104.58	110.10
4	A	406	MES	C2-C3-N4	3.39	115.25	110.10
3	A	405	GOL	O2-C2-C1	-2.94	96.20	109.12
4	A	406	MES	C6-C5-N4	-2.39	106.47	110.10
4	A	406	MES	C7-N4-C3	2.36	117.27	111.23
4	A	406	MES	O2S-S-O1S	2.35	122.06	113.95
4	B	404	MES	C6-O1-C2	2.15	117.08	109.89
3	A	405	GOL	C3-C2-C1	-2.13	103.40	111.70

There are no chirality outliers.

All (25) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	403	GOL	O1-C1-C2-C3
3	A	404	GOL	O1-C1-C2-C3
3	A	404	GOL	C1-C2-C3-O3
3	A	405	GOL	O1-C1-C2-O2
3	A	405	GOL	O1-C1-C2-C3
3	A	405	GOL	C1-C2-C3-O3
3	B	403	GOL	O1-C1-C2-O2
3	B	403	GOL	O1-C1-C2-C3
4	A	406	MES	C8-C7-N4-C3
4	A	406	MES	C8-C7-N4-C5
4	A	406	MES	N4-C7-C8-S
4	A	406	MES	C7-C8-S-O2S
4	B	404	MES	C8-C7-N4-C3
4	B	404	MES	C7-C8-S-O2S
3	A	405	GOL	O2-C2-C3-O3
4	B	404	MES	C7-C8-S-O3S
3	A	403	GOL	C1-C2-C3-O3
3	B	403	GOL	C1-C2-C3-O3
3	A	404	GOL	O1-C1-C2-O2
3	A	404	GOL	O2-C2-C3-O3
4	B	404	MES	C8-C7-N4-C5
3	B	403	GOL	O2-C2-C3-O3
3	A	403	GOL	O1-C1-C2-O2
4	A	406	MES	C7-C8-S-O1S
3	A	403	GOL	O2-C2-C3-O3

There are no ring outliers.

6 monomers are involved in 35 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	406	MES	1	0
3	B	403	GOL	16	0
3	A	405	GOL	8	0
4	B	404	MES	1	0
3	A	404	GOL	1	0
3	A	403	GOL	8	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	308/311 (99%)	-0.18	2 (0%) 89 90	16, 27, 44, 59	0
1	B	306/311 (98%)	-0.16	0 100 100	16, 27, 43, 65	0
All	All	614/622 (98%)	-0.17	2 (0%) 94 94	16, 27, 44, 65	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	299	THR	4.2
1	A	298	ALA	3.1

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
3	GOL	B	403	6/6	0.82	0.32	21,24,28,33	6
3	GOL	A	403	6/6	0.88	0.32	30,39,40,42	0
3	GOL	A	404	6/6	0.90	0.25	25,32,38,44	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	GOL	A	405	6/6	0.92	0.22	13,15,16,16	6
2	K	A	401	1/1	0.92	0.07	44,44,44,44	0
4	MES	B	404	12/12	0.92	0.12	33,43,49,52	0
2	K	B	402	1/1	0.93	0.15	48,48,48,48	0
4	MES	A	406	12/12	0.96	0.10	28,32,45,46	0
2	K	B	401	1/1	0.98	0.06	46,46,46,46	0
2	K	A	402	1/1	0.99	0.11	36,36,36,36	0

6.5 Other polymers (i)

There are no such residues in this entry.