Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 03:17 pm BST

PDB ID	:	3DBN
Title	:	Crystal structure of the Streptoccocus suis serotype2 D-mannonate dehydratase in complex with its substrate
Authors	:	Peng, H.; Zhang, Q.; Gao, F.; Gao, G.F.
Deposited on	:	2008-06-02
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 ⓘ symbol.

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

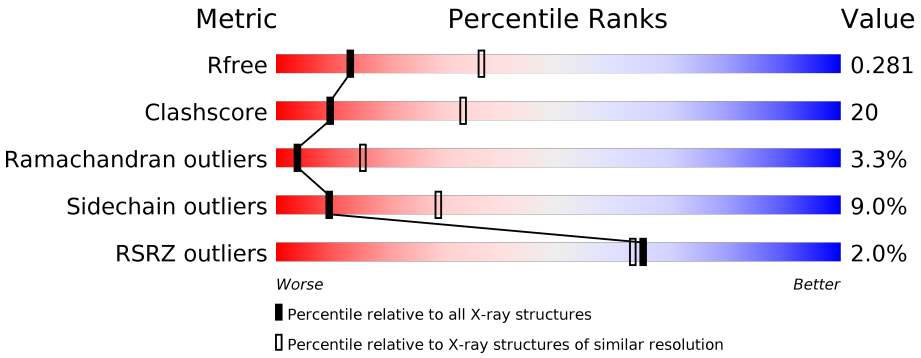
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

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 on 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% 54% 30% 6% 10%
1	B	386	2% 53% 33% 5% 10%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 5484 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 Mannonate dehydratase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	349	Total	C	N	O	S	0	0	0
			2737	1745	465	513	14			
1	B	349	Total	C	N	O	S	0	0	0
			2732	1742	465	512	13			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	EXPRESSION TAG	UNP A4VVI4
A	2	GLY	-	EXPRESSION TAG	UNP A4VVI4
A	3	SER	-	EXPRESSION TAG	UNP A4VVI4
A	4	SER	-	EXPRESSION TAG	UNP A4VVI4
A	5	HIS	-	EXPRESSION TAG	UNP A4VVI4
A	6	HIS	-	EXPRESSION TAG	UNP A4VVI4
A	7	HIS	-	EXPRESSION TAG	UNP A4VVI4
A	8	HIS	-	EXPRESSION TAG	UNP A4VVI4
A	9	HIS	-	EXPRESSION TAG	UNP A4VVI4
A	10	HIS	-	EXPRESSION TAG	UNP A4VVI4
A	11	SER	-	EXPRESSION TAG	UNP A4VVI4
A	12	SER	-	EXPRESSION TAG	UNP A4VVI4
A	13	GLY	-	EXPRESSION TAG	UNP A4VVI4
A	14	LEU	-	EXPRESSION TAG	UNP A4VVI4
A	15	VAL	-	EXPRESSION TAG	UNP A4VVI4
A	16	PRO	-	EXPRESSION TAG	UNP A4VVI4
A	17	ARG	-	EXPRESSION TAG	UNP A4VVI4
A	18	GLY	-	EXPRESSION TAG	UNP A4VVI4
A	19	SER	-	EXPRESSION TAG	UNP A4VVI4
A	20	HIS	-	EXPRESSION TAG	UNP A4VVI4
B	1	MET	-	EXPRESSION TAG	UNP A4VVI4
B	2	GLY	-	EXPRESSION TAG	UNP A4VVI4
B	3	SER	-	EXPRESSION TAG	UNP A4VVI4
B	4	SER	-	EXPRESSION TAG	UNP A4VVI4
B	5	HIS	-	EXPRESSION TAG	UNP A4VVI4

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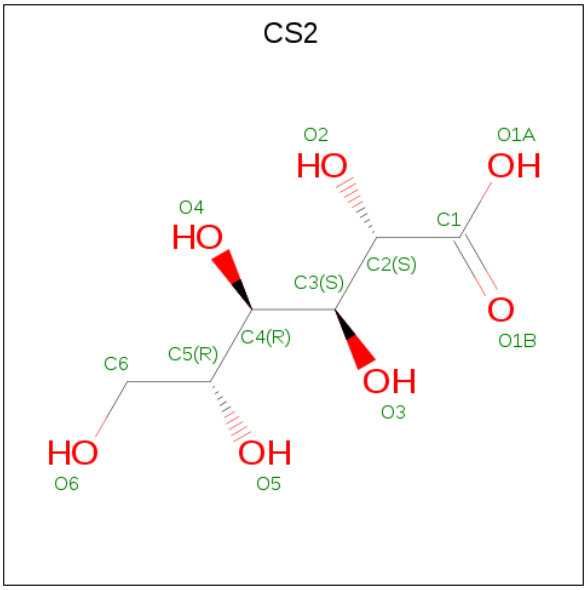
Chain	Residue	Modelled	Actual	Comment	Reference
B	6	HIS	-	EXPRESSION TAG	UNP A4VVI4
B	7	HIS	-	EXPRESSION TAG	UNP A4VVI4
B	8	HIS	-	EXPRESSION TAG	UNP A4VVI4
B	9	HIS	-	EXPRESSION TAG	UNP A4VVI4
B	10	HIS	-	EXPRESSION TAG	UNP A4VVI4
B	11	SER	-	EXPRESSION TAG	UNP A4VVI4
B	12	SER	-	EXPRESSION TAG	UNP A4VVI4
B	13	GLY	-	EXPRESSION TAG	UNP A4VVI4
B	14	LEU	-	EXPRESSION TAG	UNP A4VVI4
B	15	VAL	-	EXPRESSION TAG	UNP A4VVI4
B	16	PRO	-	EXPRESSION TAG	UNP A4VVI4
B	17	ARG	-	EXPRESSION TAG	UNP A4VVI4
B	18	GLY	-	EXPRESSION TAG	UNP A4VVI4
B	19	SER	-	EXPRESSION TAG	UNP A4VVI4
B	20	HIS	-	EXPRESSION TAG	UNP A4VVI4

- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

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

- Molecule 3 is D-MANNONIC ACID (three-letter code: CS2) (formula: $C_6H_{12}O_7$).



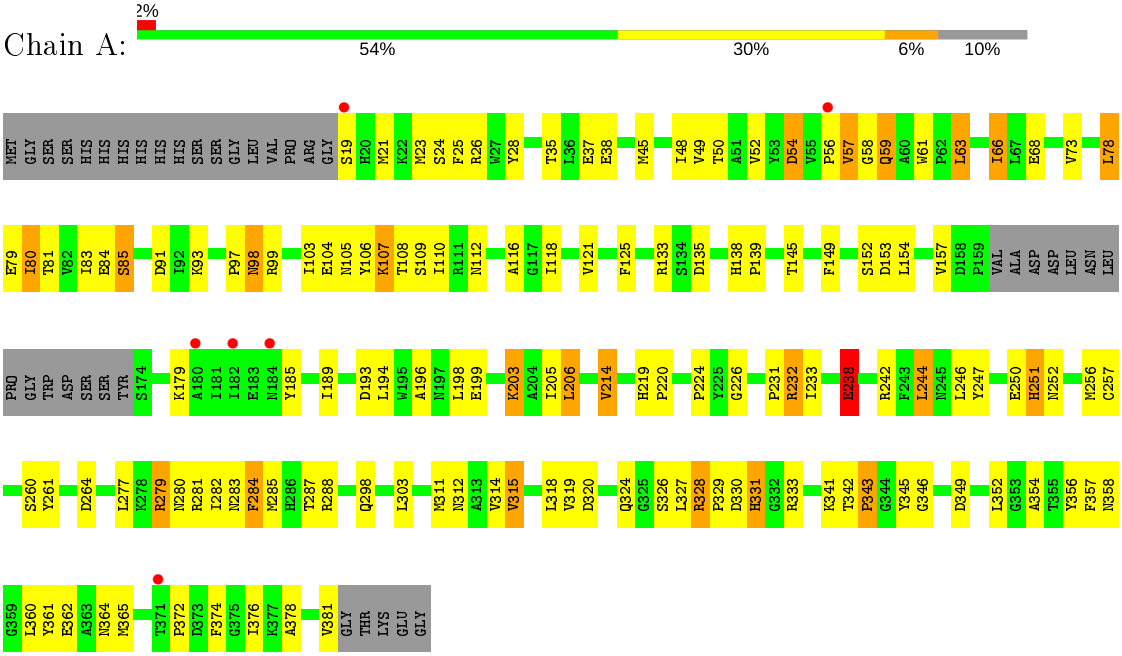
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			13	6	7		

3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

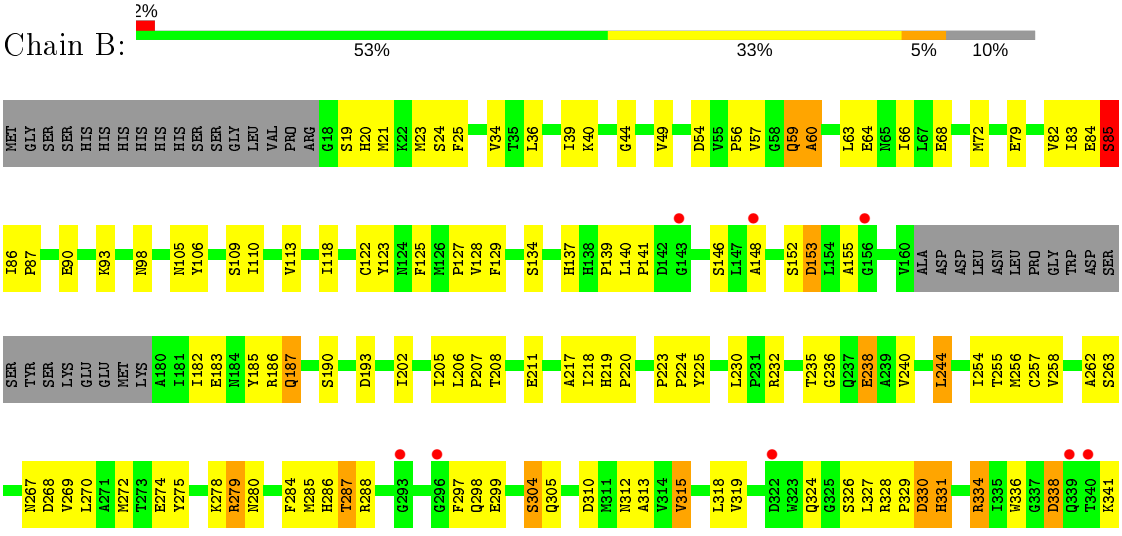
- Molecule 1: Mannonate dehydratase

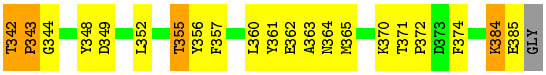
Chain A:



- Molecule 1: Mannonate dehydratase

Chain B:




T342
P343
G344
Y348
D349
L352
T355
Y356
F357
L360
Y361
E362
A363
N364
M365
K370
T371
P372
D373
F374
K384
E385
GLY

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α, β, γ	105.52Å 105.52Å 160.08Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.90 25.87 – 2.90	Depositor EDS
% Data completeness (in resolution range)	95.8 (50.00-2.90) 96.0 (25.87-2.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	8.03 (at 2.89Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.231 , 0.280 0.230 , 0.281	Depositor DCC
R_{free} test set	981 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å^2)	33.4	Xtriage
Anisotropy	0.019	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 39.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.31$, $< L^2 > = 0.14$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.82	EDS
Total number of atoms	5484	wwPDB-VP
Average B, all atoms (Å^2)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.74% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MN, CS2

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#$\|Z\|>5$	RMSZ	#$\|Z\|>5$
1	A	0.65	1/2804 (0.0%)	0.76	0/3807
1	B	0.64	0/2799	0.76	0/3801
All	All	0.65	1/5603 (0.0%)	0.76	0/7608

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	238	GLU	CB-CG	5.90	1.63	1.52

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	85	SER	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	2737	0	2676	115	0
1	B	2732	0	2678	106	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	13	0	10	2	0
All	All	5484	0	5364	212	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 20.

All (212) 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:182:ILE:HG22	1:B:186:ARG:HD2	1.31	1.06
1:B:287:THR:HG23	1:B:287:THR:O	1.57	1.01
1:A:226:GLY:HA2	1:A:232:ARG:HD3	1.48	0.94
1:A:250:GLU:HG2	1:A:280:ASN:HD22	1.31	0.92
1:A:57:VAL:HG13	1:A:58:GLY:H	1.30	0.92
1:A:194:LEU:HD23	1:A:233:ILE:HG12	1.51	0.91
1:B:56:PRO:O	1:B:59:GLN:HB2	1.71	0.90
1:B:287:THR:O	1:B:287:THR:CG2	2.25	0.84
1:B:57:VAL:HG13	1:B:87:PRO:HB2	1.61	0.81
1:B:36:LEU:HA	1:B:39:ILE:HD12	1.64	0.78
1:B:236:GLY:HA3	1:B:267:ASN:HD21	1.52	0.75
1:A:330:ASP:O	1:A:331:HIS:O	2.06	0.73
1:A:57:VAL:HG13	1:A:58:GLY:N	2.04	0.73
1:A:21:MET:HE1	1:A:23:MET:HE1	1.70	0.73
1:A:19:SER:N	1:A:324:GLN:HE21	1.87	0.72
1:A:104:GLU:HA	1:A:107:LYS:HG3	1.73	0.71
1:A:250:GLU:HG3	1:A:281:ARG:NH1	2.06	0.70
1:B:182:ILE:HG22	1:B:186:ARG:CD	2.18	0.70
1:A:49:VAL:HG13	1:A:84:GLU:HB2	1.75	0.69
1:B:140:LEU:HD11	1:B:146:SER:HB3	1.73	0.69
1:A:121:VAL:HG23	1:A:214:VAL:CG1	2.23	0.68
1:A:121:VAL:HG23	1:A:214:VAL:HG13	1.76	0.68
1:B:236:GLY:O	1:B:240:VAL:HG23	1.95	0.67
1:B:287:THR:HG21	1:B:327:LEU:HD21	1.77	0.67
1:A:312:ASN:HD21	1:A:364:ASN:HD21	1.42	0.66
1:A:220:PRO:HB3	1:A:261:TYR:CZ	2.30	0.66
1:B:49:VAL:HG22	1:B:82:VAL:HB	1.77	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:283:ASN:O	1:A:284:PHE:HB2	1.96	0.66
1:B:236:GLY:HA3	1:B:267:ASN:ND2	2.10	0.65
1:A:48:ILE:HG21	1:A:73:VAL:HG11	1.79	0.65
1:A:57:VAL:CG1	1:A:58:GLY:H	2.07	0.65
1:B:287:THR:HG23	1:B:357:PHE:CZ	2.33	0.64
1:B:286:HIS:CE1	1:B:328:ARG:HH11	2.16	0.64
1:A:312:ASN:ND2	1:A:364:ASN:HD21	1.95	0.64
1:A:26:ARG:HA	1:A:49:VAL:HB	1.79	0.64
1:A:250:GLU:HG3	1:A:281:ARG:HH12	1.63	0.63
1:A:352:LEU:HD22	1:B:355:THR:CG2	2.29	0.63
1:A:91:ASP:HB3	1:A:98:ASN:HD21	1.62	0.63
1:A:199:GLU:O	1:A:203:LYS:HG2	1.99	0.63
1:A:288:ARG:HD3	1:A:330:ASP:HB3	1.80	0.63
1:B:105:ASN:O	1:B:109:SER:OG	2.16	0.62
1:A:319:VAL:HG13	1:A:365:MET:HE3	1.82	0.62
1:A:66:ILE:HD11	1:A:116:ALA:CB	2.29	0.61
1:A:73:VAL:HG13	1:A:78:LEU:HB2	1.81	0.61
1:A:84:GLU:O	1:A:85:SER:HB2	2.01	0.61
1:B:205:ILE:O	1:B:208:THR:HB	1.99	0.61
1:B:182:ILE:CG2	1:B:186:ARG:HD2	2.19	0.60
1:B:90:GLU:OE2	1:B:128:VAL:N	2.29	0.60
1:A:206:LEU:HB3	1:A:247:TYR:CZ	2.36	0.60
1:B:330:ASP:O	1:B:331:HIS:HB2	2.01	0.60
1:B:183:GLU:HA	1:B:186:ARG:HB2	1.85	0.59
1:A:193:ASP:O	1:A:196:ALA:HB3	2.01	0.59
1:B:106:TYR:O	1:B:110:ILE:HG13	2.02	0.59
1:B:235:THR:O	1:B:267:ASN:ND2	2.36	0.59
1:A:21:MET:HE1	1:A:23:MET:CE	2.34	0.58
1:A:125:PHE:CG	1:A:198:LEU:HD13	2.38	0.58
1:A:311:MET:O	1:A:315:VAL:HG12	2.04	0.58
1:A:358:ASN:HD22	1:A:374:PHE:HE1	1.53	0.57
1:A:25:PHE:HB2	1:A:45:MET:SD	2.45	0.57
1:A:314:VAL:O	1:A:318:LEU:HG	2.04	0.57
1:B:331:HIS:HA	1:B:344:GLY:O	2.05	0.57
1:B:40:LYS:O	1:B:40:LYS:HD3	2.04	0.56
1:A:49:VAL:CG1	1:A:84:GLU:HB2	2.35	0.56
1:A:352:LEU:HD22	1:B:355:THR:HG21	1.85	0.56
1:B:85:SER:HA	1:B:122:CYS:O	2.06	0.56
1:A:135:ASP:OD2	1:A:138:HIS:ND1	2.33	0.56
1:B:123:TYR:CE1	1:B:218:ILE:HD12	2.40	0.56
1:A:21:MET:CE	1:A:23:MET:CE	2.84	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:83:ILE:HG12	1:B:118:ILE:HG21	1.88	0.56
1:B:312:ASN:ND2	1:B:364:ASN:HD21	2.04	0.55
1:A:35:THR:HB	1:A:38:GLU:HG3	1.88	0.55
1:B:310:ASP:OD1	1:B:313:ALA:HB2	2.05	0.55
1:A:372:PRO:HG2	1:B:305:GLN:HE22	1.72	0.55
1:B:284:PHE:CE1	1:B:286:HIS:HE1	2.25	0.55
1:A:54:ASP:N	1:A:54:ASP:OD1	2.39	0.55
1:A:80:ILE:HG21	1:A:118:ILE:HD13	1.88	0.55
1:A:356:TYR:OH	1:B:362:GLU:OE1	2.24	0.55
1:A:21:MET:HE3	1:A:23:MET:HE2	1.89	0.54
1:A:312:ASN:HD22	1:A:360:LEU:HD22	1.72	0.54
1:B:356:TYR:CZ	1:B:360:LEU:HD11	2.41	0.54
1:A:206:LEU:HD11	1:A:252:ASN:HB2	1.89	0.54
1:B:20:HIS:H	1:B:324:GLN:HG2	1.73	0.53
1:A:99:ARG:O	1:A:103:ILE:HG13	2.08	0.53
1:B:153:ASP:N	1:B:153:ASP:OD1	2.41	0.53
1:B:319:VAL:HG21	1:B:364:ASN:HB3	1.90	0.53
1:A:116:ALA:O	1:A:118:ILE:HG13	2.09	0.53
1:B:267:ASN:HB3	1:B:272:MET:HE1	1.89	0.52
1:B:262:ALA:HB2	1:B:269:VAL:CG1	2.39	0.52
1:A:288:ARG:HD2	1:A:331:HIS:HB2	1.90	0.52
1:B:21:MET:HE3	1:B:361:TYR:CD2	2.44	0.52
1:A:257:CYS:HG	1:A:260:SER:HG	1.58	0.52
1:A:288:ARG:HH11	1:A:288:ARG:HG2	1.73	0.52
1:B:187:GLN:OE1	1:B:187:GLN:HA	2.10	0.52
1:B:56:PRO:HB2	1:B:59:GLN:HG3	1.92	0.52
1:B:217:ALA:HB1	1:B:255:THR:OG1	2.10	0.52
1:A:35:THR:HG22	1:A:37:GLU:H	1.75	0.52
1:B:244:LEU:HB3	1:B:279:ARG:NH1	2.25	0.52
1:A:319:VAL:HG13	1:A:365:MET:CE	2.39	0.51
1:A:21:MET:CE	1:A:23:MET:HE1	2.40	0.51
1:B:275:TYR:OH	1:B:279:ARG:NH2	2.43	0.51
1:A:331:HIS:CD2	1:A:331:HIS:H	2.28	0.51
1:A:91:ASP:CB	1:A:98:ASN:HD21	2.23	0.51
1:A:103:ILE:O	1:A:107:LYS:HG2	2.11	0.51
1:A:224:PRO:HG2	1:A:264:ASP:HB2	1.93	0.51
1:B:285:MET:O	1:B:327:LEU:HA	2.10	0.51
1:B:127:PRO:HB3	1:B:185:TYR:CZ	2.45	0.51
1:A:376:ILE:HG21	1:B:305:GLN:O	2.12	0.50
1:A:73:VAL:CG1	1:A:78:LEU:HB2	2.42	0.50
1:B:137:HIS:O	1:B:139:PRO:HD3	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:149:PHE:HE1	1:A:154:LEU:HD22	1.76	0.50
1:A:362:GLU:OE2	1:B:304:SER:OG	2.18	0.50
1:A:277:LEU:CD2	1:A:282:ILE:HD12	2.42	0.49
1:B:148:ALA:HB2	1:B:298:GLN:HG3	1.93	0.49
1:B:21:MET:CE	1:B:361:TYR:CD2	2.95	0.49
1:B:299:GLU:OE2	1:B:331:HIS:HD2	1.96	0.49
1:A:231:PRO:C	1:A:232:ARG:HD2	2.33	0.49
1:A:303:LEU:HA	1:A:356:TYR:CZ	2.48	0.49
1:B:262:ALA:HB2	1:B:269:VAL:HG11	1.95	0.48
1:A:21:MET:CE	1:A:23:MET:HE2	2.43	0.48
1:A:63:LEU:H	1:A:112:ASN:HD22	1.62	0.48
1:B:256:MET:HB3	1:B:285:MET:SD	2.54	0.48
1:B:129:PHE:HB2	1:B:232:ARG:NH1	2.29	0.48
1:A:345:TYR:HE1	3:A:388:CS2:H61	1.78	0.47
1:A:287:THR:HG23	1:A:357:PHE:CZ	2.49	0.47
1:A:372:PRO:HG2	1:B:305:GLN:NE2	2.29	0.47
1:A:61:TRP:CZ3	1:A:83:ILE:HG21	2.49	0.47
1:B:24:SER:O	1:B:329:PRO:HD2	2.14	0.47
1:B:275:TYR:HD1	1:B:278:LYS:NZ	2.13	0.46
1:B:223:PRO:O	1:B:225:TYR:N	2.40	0.46
1:B:275:TYR:O	1:B:278:LYS:HB2	2.15	0.46
1:A:93:LYS:HE2	1:A:125:PHE:HA	1.97	0.46
1:B:342:THR:HA	1:B:343:PRO:HD2	1.66	0.46
1:B:82:VAL:HG11	1:B:284:PHE:CE1	2.51	0.46
1:B:49:VAL:CG1	1:B:84:GLU:HB2	2.45	0.46
1:A:108:THR:OG1	1:A:109:SER:N	2.49	0.46
1:A:97:PRO:C	1:A:99:ARG:H	2.19	0.46
1:B:312:ASN:HD21	1:B:364:ASN:HD21	1.62	0.46
1:B:23:MET:SD	1:B:327:LEU:HD22	2.56	0.46
1:A:219:HIS:CE1	3:A:388:CS2:O1B	2.68	0.46
1:A:79:GLU:HG3	1:A:79:GLU:O	2.16	0.46
1:B:263:SER:HB2	1:B:297:PHE:CD1	2.50	0.46
1:B:365:MET:O	1:B:370:LYS:HB2	2.15	0.46
1:A:63:LEU:N	1:A:112:ASN:HD22	2.13	0.46
1:B:182:ILE:HG23	1:B:230:LEU:HD21	1.98	0.46
1:A:329:PRO:HG3	1:A:354:ALA:HA	1.98	0.45
1:A:61:TRP:HB2	1:A:109:SER:HA	1.98	0.45
1:B:59:GLN:O	1:B:60:ALA:HB3	2.17	0.45
1:A:199:GLU:HG3	1:A:246:LEU:HD22	1.97	0.45
1:B:153:ASP:C	1:B:155:ALA:H	2.19	0.45
1:A:303:LEU:HD11	1:B:374:PHE:CE2	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:283:ASN:O	1:A:284:PHE:CB	2.65	0.45
1:B:267:ASN:HB3	1:B:272:MET:CE	2.46	0.45
1:A:133:ARG:HG2	1:A:133:ARG:HH11	1.83	0.44
1:B:207:PRO:O	1:B:211:GLU:HG3	2.17	0.44
1:B:365:MET:HG3	1:B:372:PRO:HB3	1.99	0.44
1:B:315:VAL:O	1:B:318:LEU:N	2.51	0.44
1:A:287:THR:HG23	1:A:287:THR:O	2.17	0.44
1:A:342:THR:O	1:A:343:PRO:C	2.55	0.44
1:A:333:ARG:HD3	1:A:349:ASP:OD2	2.18	0.44
1:B:349:ASP:HA	1:B:352:LEU:HG	1.98	0.44
1:A:24:SER:OG	1:A:326:SER:HB2	2.18	0.44
1:A:284:PHE:HD2	1:A:326:SER:O	2.00	0.44
1:B:190:SER:H	1:B:193:ASP:HB2	1.82	0.44
1:B:206:LEU:O	1:B:207:PRO:C	2.53	0.44
1:A:48:ILE:HG22	1:A:80:ILE:N	2.32	0.44
1:B:190:SER:N	1:B:193:ASP:HB2	2.33	0.44
1:B:238:GLU:H	1:B:238:GLU:CD	2.21	0.43
1:A:59:GLN:C	1:A:105:ASN:HB3	2.39	0.43
1:B:284:PHE:CE1	1:B:286:HIS:CE1	3.06	0.43
1:A:139:PRO:HA	1:A:145:THR:HG22	2.01	0.43
1:B:256:MET:HG3	1:B:258:VAL:HG13	2.00	0.43
1:A:104:GLU:O	1:A:107:LYS:HB2	2.18	0.43
1:B:275:TYR:HD1	1:B:278:LYS:HZ2	1.65	0.43
1:A:50:THR:O	1:A:83:ILE:HA	2.17	0.43
1:B:63:LEU:HA	1:B:66:ILE:HD12	2.00	0.43
1:A:106:TYR:O	1:A:110:ILE:HG13	2.19	0.43
1:A:251:HIS:ND1	1:A:251:HIS:N	2.66	0.43
1:A:328:ARG:NH2	1:A:345:TYR:HE2	2.16	0.43
1:B:109:SER:O	1:B:113:VAL:HG23	2.18	0.43
1:B:334:ARG:HG2	1:B:338:ASP:HB3	2.01	0.43
1:B:356:TYR:O	1:B:360:LEU:HG	2.19	0.43
1:A:362:GLU:HG3	1:A:372:PRO:HG3	2.01	0.43
1:B:336:TRP:CH2	1:B:352:LEU:HD21	2.53	0.42
1:A:110:ILE:HD12	1:A:205:ILE:HB	2.01	0.42
1:A:315:VAL:HG11	1:A:360:LEU:HB3	2.01	0.42
1:A:21:MET:HE3	1:A:23:MET:CE	2.49	0.42
1:B:137:HIS:HD2	1:B:334:ARG:HH21	1.68	0.42
1:A:179:LYS:HG2	1:A:179:LYS:O	2.19	0.42
1:A:378:ALA:O	1:B:141:PRO:HG3	2.19	0.42
1:B:44:GLY:HA2	1:B:374:PHE:HB3	2.01	0.42
1:B:54:ASP:N	1:B:54:ASP:OD1	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:93:LYS:HE2	1:B:125:PHE:HA	2.02	0.42
1:A:238:GLU:H	1:A:238:GLU:HG3	1.60	0.42
1:A:24:SER:OG	1:A:81:THR:HG21	2.20	0.42
1:A:256:MET:HB3	1:A:285:MET:SD	2.60	0.42
1:B:219:HIS:HE1	1:B:286:HIS:HD2	1.68	0.41
1:A:185:TYR:HA	1:A:189:ILE:HB	2.01	0.41
1:A:121:VAL:CG2	1:A:214:VAL:HG11	2.50	0.41
1:A:279:ARG:HE	1:A:279:ARG:HB3	1.40	0.41
1:A:328:ARG:NH2	1:A:330:ASP:OD1	2.54	0.41
1:B:336:TRP:HH2	1:B:352:LEU:HD21	1.85	0.41
1:B:310:ASP:OD1	1:B:313:ALA:CB	2.69	0.41
1:B:64:GLU:N	1:B:64:GLU:OE1	2.53	0.41
1:B:68:GLU:O	1:B:72:MET:HG3	2.19	0.41
1:B:257:CYS:HA	1:B:286:HIS:HB2	2.03	0.41
1:A:121:VAL:HG23	1:A:214:VAL:HG11	2.01	0.40
1:B:202:ILE:HA	1:B:205:ILE:HG12	2.02	0.40
1:B:284:PHE:HD2	1:B:326:SER:O	2.04	0.40
1:A:244:LEU:HG	1:A:281:ARG:HG3	2.02	0.40
1:A:52:VAL:CG1	1:A:54:ASP:OD1	2.69	0.40
1:B:270:LEU:HD21	1:B:313:ALA:HB3	2.02	0.40
1:B:384:LYS:HG3	1:B:385:GLU:H	1.86	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	345/386 (89%)	290 (84%)	43 (12%)	12 (4%)	3 14
1	B	345/386 (89%)	301 (87%)	33 (10%)	11 (3%)	4 16
All	All	690/772 (89%)	591 (86%)	76 (11%)	23 (3%)	4 15

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	57	VAL
1	A	284	PHE
1	A	331	HIS
1	A	343	PRO
1	B	343	PRO
1	A	28	TYR
1	A	85	SER
1	A	346	GLY
1	B	330	ASP
1	B	60	ALA
1	B	220	PRO
1	B	384	LYS
1	A	56	PRO
1	A	98	ASN
1	B	224	PRO
1	B	304	SER
1	B	348	TYR
1	A	361	TYR
1	B	338	ASP
1	B	363	ALA
1	A	68	GLU
1	B	85	SER
1	A	214	VAL

5.3.2 Protein sidechains

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

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	288/321 (90%)	263 (91%)	25 (9%)	10 30
1	B	288/321 (90%)	261 (91%)	27 (9%)	8 26
All	All	576/642 (90%)	524 (91%)	52 (9%)	9 29

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	54	ASP
1	A	59	GLN
1	A	63	LEU
1	A	66	ILE
1	A	78	LEU
1	A	80	ILE
1	A	107	LYS
1	A	152	SER
1	A	153	ASP
1	A	157	VAL
1	A	203	LYS
1	A	206	LEU
1	A	232	ARG
1	A	238	GLU
1	A	242	ARG
1	A	244	LEU
1	A	251	HIS
1	A	279	ARG
1	A	298	GLN
1	A	315	VAL
1	A	320	ASP
1	A	327	LEU
1	A	328	ARG
1	A	341	LYS
1	A	381	VAL
1	B	19	SER
1	B	25	PHE
1	B	34	VAL
1	B	59	GLN
1	B	79	GLU
1	B	86	ILE
1	B	98	ASN
1	B	134	SER
1	B	152	SER
1	B	153	ASP
1	B	187	GLN
1	B	238	GLU
1	B	244	LEU
1	B	254	ILE
1	B	268	ASP
1	B	274	GLU
1	B	279	ARG
1	B	280	ASN

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Mol	Chain	Res	Type
1	B	287	THR
1	B	288	ARG
1	B	315	VAL
1	B	331	HIS
1	B	334	ARG
1	B	341	LYS
1	B	342	THR
1	B	355	THR
1	B	371	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	98	ASN
1	A	105	ASN
1	A	112	ASN
1	A	280	ASN
1	A	283	ASN
1	A	298	GLN
1	A	305	GLN
1	A	312	ASN
1	B	59	GLN
1	B	98	ASN
1	B	105	ASN
1	B	137	HIS
1	B	219	HIS
1	B	237	GLN
1	B	251	HIS
1	B	267	ASN
1	B	283	ASN
1	B	286	HIS
1	B	298	GLN
1	B	305	GLN
1	B	312	ASN
1	B	324	GLN
1	B	331	HIS

5.3.3 RNA

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 carbohydrates 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	$\#\|Z\| > 2$
3	CS2	A	388	2	9,12,12	0.74	0	12,16,16	0.83	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
3	CS2	A	388	2	-	7/14/18/18	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	388	CS2	C1-C2-C3-O3
3	A	388	CS2	O2-C2-C3-O3
3	A	388	CS2	O2-C2-C3-C4
3	A	388	CS2	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
3	A	388	CS2	O5-C5-C6-O6
3	A	388	CS2	C3-C4-C5-C6
3	A	388	CS2	C1-C2-C3-C4

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	388	CS2	2	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	349/386 (90%)	-0.15	6 (1%)	70	69	3, 22, 42, 49	0
1	B	349/386 (90%)	-0.22	8 (2%)	60	58	2, 20, 42, 63	0
All	All	698/772 (90%)	-0.19	14 (2%)	65	63	2, 21, 42, 63	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	180	ALA	4.4
1	A	56	PRO	3.6
1	B	148	ALA	2.9
1	B	340	THR	2.8
1	B	293	GLY	2.7
1	A	19	SER	2.6
1	B	156	GLY	2.5
1	B	339	GLN	2.4
1	B	322	ASP	2.4
1	B	143	GLY	2.3
1	A	184	ASN	2.2
1	A	182	ILE	2.1
1	A	371	THR	2.1
1	B	296	GLY	2.1

6.2 Non-standard residues in protein, DNA, RNA chains

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates

There are no carbohydrates in this entry.

6.4 Ligands

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, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	CS2	A	388	13/13	0.93	0.17	41,44,46,47	0
2	MN	B	387	1/1	0.95	0.13	52,52,52,52	0
2	MN	A	387	1/1	0.95	0.06	31,31,31,31	0

6.5 Other polymers

There are no such residues in this entry.