



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 9, 2020 – 01:00 AM BST

PDB ID : 4YTQ
Title : Crystal structure of D-tagatose 3-epimerase C66S from *Pseudomonas cichorii*
in complex with 1-deoxy D-tagatose
Authors : Yoshida, H.; Yoshihara, A.; Ishii, T.; Izumori, K.; Kamitori, S.
Deposited on : 2015-03-18
Resolution : 1.90 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 : 2.13.1
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) : 2.13.1

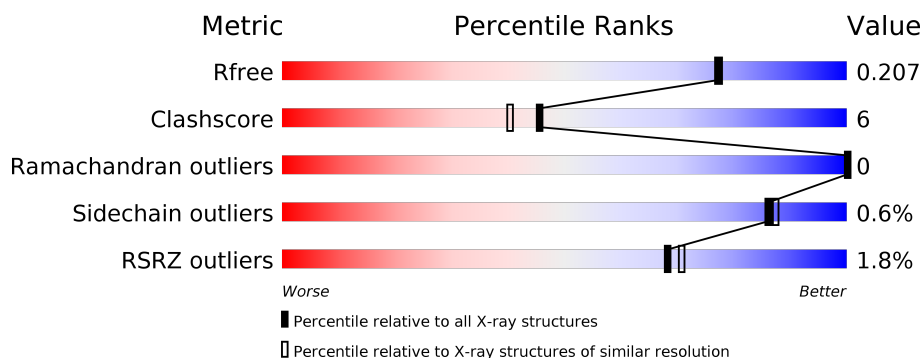
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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 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 $\leq 5\%$. 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	300	<div> <div>2%</div> <div> <div></div> <div>86%</div> <div>11%</div> <div>••</div> </div> </div>
1	B	300	<div> <div>2%</div> <div> <div></div> <div>85%</div> <div>12%</div> <div>•</div> </div> </div>
1	C	300	<div> <div>2%</div> <div> <div></div> <div>85%</div> <div>13%</div> <div>••</div> </div> </div>
1	D	300	<div> <div>2%</div> <div> <div></div> <div>78%</div> <div>18%</div> <div>••</div> </div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10508 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 D-tagatose 3-epimerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	294	Total	C	N	O	S	0	0	0
			2314	1465	399	432	18			
1	B	292	Total	C	N	O	S	0	0	0
			2297	1456	394	429	18			
1	C	297	Total	C	N	O	S	0	0	0
			2344	1483	408	435	18			
1	D	290	Total	C	N	O	S	0	0	0
			2287	1451	392	426	18			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	66	SER	CYS	engineered mutation	UNP O50580
A	291	GLY	-	expression tag	UNP O50580
A	292	SER	-	expression tag	UNP O50580
A	293	ARG	-	expression tag	UNP O50580
A	294	SER	-	expression tag	UNP O50580
A	295	HIS	-	expression tag	UNP O50580
A	296	HIS	-	expression tag	UNP O50580
A	297	HIS	-	expression tag	UNP O50580
A	298	HIS	-	expression tag	UNP O50580
A	299	HIS	-	expression tag	UNP O50580
A	300	HIS	-	expression tag	UNP O50580
B	66	SER	CYS	engineered mutation	UNP O50580
B	291	GLY	-	expression tag	UNP O50580
B	292	SER	-	expression tag	UNP O50580
B	293	ARG	-	expression tag	UNP O50580
B	294	SER	-	expression tag	UNP O50580
B	295	HIS	-	expression tag	UNP O50580
B	296	HIS	-	expression tag	UNP O50580
B	297	HIS	-	expression tag	UNP O50580
B	298	HIS	-	expression tag	UNP O50580
B	299	HIS	-	expression tag	UNP O50580

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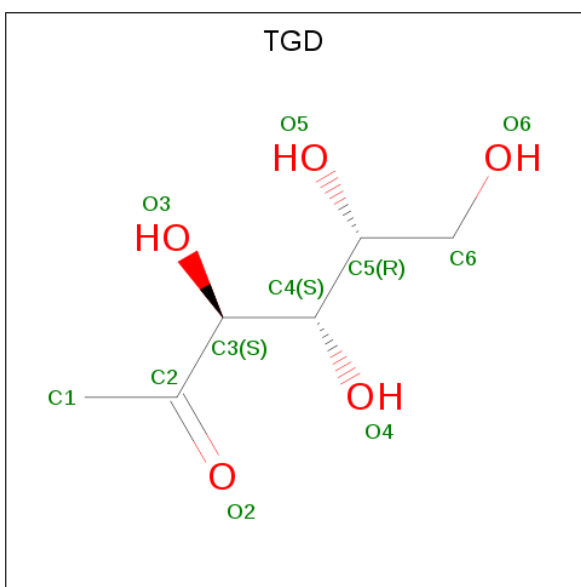
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Chain	Residue	Modelled	Actual	Comment	Reference
B	300	HIS	-	expression tag	UNP O50580
C	66	SER	CYS	engineered mutation	UNP O50580
C	291	GLY	-	expression tag	UNP O50580
C	292	SER	-	expression tag	UNP O50580
C	293	ARG	-	expression tag	UNP O50580
C	294	SER	-	expression tag	UNP O50580
C	295	HIS	-	expression tag	UNP O50580
C	296	HIS	-	expression tag	UNP O50580
C	297	HIS	-	expression tag	UNP O50580
C	298	HIS	-	expression tag	UNP O50580
C	299	HIS	-	expression tag	UNP O50580
C	300	HIS	-	expression tag	UNP O50580
D	66	SER	CYS	engineered mutation	UNP O50580
D	291	GLY	-	expression tag	UNP O50580
D	292	SER	-	expression tag	UNP O50580
D	293	ARG	-	expression tag	UNP O50580
D	294	SER	-	expression tag	UNP O50580
D	295	HIS	-	expression tag	UNP O50580
D	296	HIS	-	expression tag	UNP O50580
D	297	HIS	-	expression tag	UNP O50580
D	298	HIS	-	expression tag	UNP O50580
D	299	HIS	-	expression tag	UNP O50580
D	300	HIS	-	expression tag	UNP O50580

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

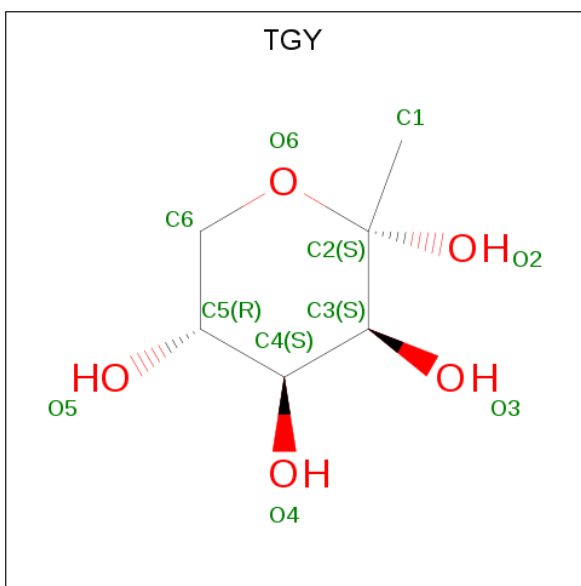
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Mn 1 1	0	0
2	A	1	Total Mn 1 1	0	0
2	D	1	Total Mn 1 1	0	0
2	C	1	Total Mn 1 1	0	0

- Molecule 3 is 1-deoxy-D-tagatose (three-letter code: TGD) (formula: C₆H₁₂O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	1
			22	12	10		

- Molecule 4 is 1-deoxy-alpha-D-tagatopyranose (three-letter code: TGY) (formula: $C_6H_{12}O_5$).

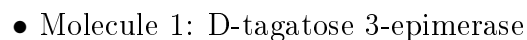
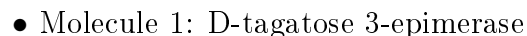
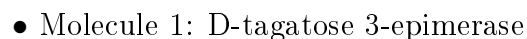


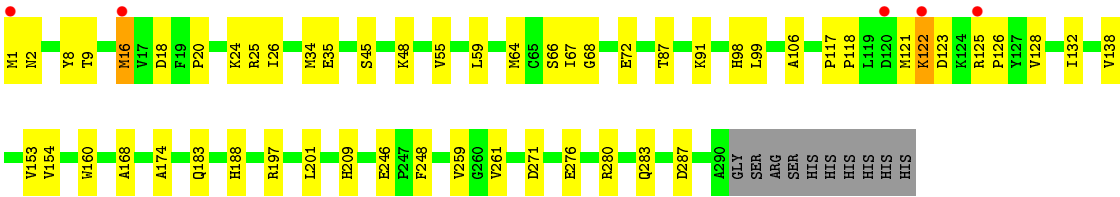
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	C	1	Total	C	O	0	0
			11	6	5		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	324	Total 324	O 324	0	0
5	B	324	Total 324	O 324	0	0
5	C	339	Total 339	O 339	0	0
5	D	242	Total 242	O 242	0	0

- Molecule 1: D-tagatose 3-epimerase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	52.33 Å 126.31 Å 99.50 Å 90.00° 101.77° 90.00°	Depositor
Resolution (Å)	19.32 – 1.90 19.32 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.5 (19.32-1.90) 99.6 (19.32-1.90)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.13 (at 1.90 Å)	Xtriage
Refinement program	CNS 1.3	Depositor
R, R_{free}	0.174 , 0.213 0.168 , 0.207	Depositor DCC
R_{free} test set	9910 reflections (10.02%)	wwPDB-VP
Wilson B-factor (Å ²)	18.3	Xtriage
Anisotropy	0.308	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 56.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	10508	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.05% 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: MN, TGY, TGD

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	0.32	0/2365	0.59	1/3190 (0.0%)
1	B	0.32	0/2348	0.58	1/3168 (0.0%)
1	C	0.31	0/2398	0.58	1/3235 (0.0%)
1	D	0.30	0/2338	0.57	0/3155
All	All	0.31	0/9449	0.58	3/12748 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	68	GLY	N-CA-C	-5.47	99.41	113.10
1	B	68	GLY	N-CA-C	-5.38	99.66	113.10
1	C	68	GLY	N-CA-C	-5.11	100.32	113.10

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	2314	0	2271	27	0
1	B	2297	0	2253	28	0
1	C	2344	0	2292	25	0
1	D	2287	0	2245	36	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	22	0	0	3	0
4	C	11	0	0	0	0
5	A	324	0	0	2	0
5	B	324	0	0	1	0
5	C	339	0	0	1	0
5	D	242	0	0	4	0
All	All	10508	0	9061	117	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 (117) 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:D:55:VAL:O	1:D:59:LEU:HD13	1.83	0.78
1:C:169:ILE:HD13	1:C:182:VAL:HG21	1.70	0.72
1:C:133:GLU:OE2	1:C:137:ARG:HD2	1.91	0.70
1:A:106:ALA:HB1	1:A:183:GLN:OE1	1.92	0.69
1:B:169:ILE:HD13	1:B:182:VAL:HG21	1.76	0.67
1:D:106:ALA:HB1	1:D:183:GLN:OE1	1.96	0.65
1:D:276:GLU:HG3	1:D:280:ARG:NH1	2.12	0.65
1:C:106:ALA:HB1	1:C:183:GLN:OE1	1.97	0.64
1:D:125:ARG:HB2	1:D:126:PRO:HD3	1.80	0.62
1:B:106:ALA:HB1	1:B:183:GLN:OE1	2.00	0.60
1:D:24:LYS:HA	1:D:59:LEU:HD21	1.84	0.59
1:A:292:SER:O	1:A:293:ARG:HD3	2.02	0.59
1:D:72:GLU:HG2	5:D:655:HOH:O	2.02	0.59
1:D:87:THR:O	1:D:91:LYS:HG3	2.04	0.58
1:B:67:ILE:HG13	1:B:68:GLY:N	2.19	0.57
1:A:108:LEU:HG	1:A:113:TRP:HB2	1.88	0.55
1:C:165:ALA:O	1:C:169:ILE:HG12	2.05	0.55
1:B:259:VAL:HG23	1:B:261:VAL:HG23	1.88	0.55
1:D:16:MET:HB2	5:D:576:HOH:O	2.06	0.55
1:A:270:THR:OG1	1:A:273:GLU:HG3	2.07	0.55
1:A:197:ARG:CZ	1:A:201:LEU:HD11	2.36	0.55
1:C:246:GLU:HG2	1:C:248:PHE:CE2	2.41	0.55
1:D:122:LYS:H	1:D:122:LYS:HD3	1.70	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:69:LEU:HB2	1:C:111:CYS:O	2.07	0.54
1:D:153:VAL:HG22	1:D:168:ALA:HB2	1.89	0.54
1:D:123:ASP:OD2	1:D:125:ARG:HG3	2.08	0.54
1:D:45:SER:OG	1:D:48:LYS:HG2	2.08	0.54
1:D:118:PRO:HD2	1:D:121:MET:HB2	1.90	0.54
1:A:259:VAL:HG23	1:A:261:VAL:HG23	1.89	0.53
1:D:99:LEU:C	1:D:99:LEU:HD23	2.29	0.53
1:A:125:ARG:HB3	1:A:126:PRO:HD3	1.90	0.53
1:A:153:VAL:HG22	1:A:168:ALA:HB2	1.92	0.52
1:C:117:PRO:HG3	1:C:160:TRP:CG	2.44	0.52
1:A:99:LEU:C	1:A:99:LEU:HD23	2.29	0.52
1:B:99:LEU:HD23	1:B:99:LEU:C	2.30	0.52
1:D:117:PRO:HG3	1:D:160:TRP:CG	2.45	0.52
1:D:67:ILE:HG13	1:D:68:GLY:N	2.25	0.52
1:C:99:LEU:HD23	1:C:99:LEU:C	2.29	0.52
1:C:67:ILE:HG13	1:C:68:GLY:N	2.25	0.52
1:C:197:ARG:CZ	1:C:201:LEU:HD11	2.40	0.52
1:B:125:ARG:HB3	1:B:126:PRO:HD3	1.92	0.51
1:B:249:MET:CE	1:B:277:ARG:NH2	2.73	0.51
1:D:246:GLU:HG2	1:D:248:PHE:CE2	2.45	0.51
1:C:153:VAL:HG22	1:C:168:ALA:HB2	1.92	0.51
1:D:117:PRO:HG3	1:D:160:TRP:CD2	2.48	0.49
1:C:259:VAL:HG23	1:C:261:VAL:HG23	1.95	0.49
1:A:117:PRO:HG3	1:A:160:TRP:CG	2.48	0.49
1:D:1:MET:HG2	1:D:2:ASN:N	2.28	0.48
1:B:67:ILE:HG13	1:B:68:GLY:H	1.78	0.48
1:B:246:GLU:HG2	1:B:248:PHE:CE2	2.48	0.48
1:B:117:PRO:HG3	1:B:160:TRP:CG	2.48	0.48
1:B:117:PRO:HG3	1:B:160:TRP:CD2	2.49	0.48
1:D:91:LYS:HE3	1:D:138:VAL:HG12	1.96	0.47
1:A:246:GLU:HG2	1:A:248:PHE:CE2	2.49	0.47
1:A:66:SER:HB2	3:A:402[B]:TGD:O6	2.14	0.47
1:B:240:ASP:O	1:B:240:ASP:OD1	2.32	0.47
1:C:132:ILE:HD13	1:C:174:ALA:HB3	1.97	0.47
1:B:249:MET:HE3	1:B:277:ARG:NH2	2.29	0.47
1:A:66:SER:HA	1:A:106:ALA:O	2.15	0.46
1:B:18:ASP:OD2	1:B:20:PRO:HD2	2.15	0.46
1:B:2:ASN:ND2	5:B:501:HOH:O	2.40	0.46
1:D:98:HIS:HD2	5:D:534:HOH:O	1.98	0.46
1:B:50:ARG:HH11	1:B:50:ARG:HG2	1.80	0.46
1:C:132:ILE:HD13	1:C:174:ALA:CB	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:255:VAL:O	1:C:259:VAL:HG22	2.16	0.46
1:D:66:SER:HA	1:D:106:ALA:O	2.16	0.46
1:A:154:VAL:HA	1:A:188:HIS:CG	2.51	0.46
1:A:3:LYS:HE3	5:A:770:HOH:O	2.16	0.46
1:D:132:ILE:HD13	1:D:174:ALA:CB	2.46	0.46
1:D:197:ARG:CZ	1:D:201:LEU:HD11	2.46	0.46
1:B:66:SER:HA	1:B:106:ALA:O	2.15	0.46
1:B:75:PHE:CZ	1:B:90:VAL:HG21	2.51	0.46
1:A:67:ILE:HG13	1:A:68:GLY:N	2.31	0.46
1:C:8:TYR:CE1	1:C:9:THR:HG23	2.51	0.45
1:B:165:ALA:O	1:B:169:ILE:HG12	2.17	0.45
1:B:153:VAL:HG22	1:B:168:ALA:HB2	1.98	0.45
1:C:152:GLU:HA	1:C:183:GLN:HB3	1.99	0.45
1:C:20:PRO:O	1:C:24:LYS:HG3	2.17	0.45
1:A:117:PRO:HG3	1:A:160:TRP:CD2	2.51	0.44
1:C:66:SER:HA	1:C:106:ALA:O	2.17	0.44
1:B:16:MET:CE	1:B:44:LEU:HD21	2.47	0.44
1:A:188:HIS:HE2	3:A:402[A]:TGD:C6	2.30	0.44
1:A:108:LEU:N	1:A:108:LEU:HD22	2.32	0.44
1:C:117:PRO:HG3	1:C:160:TRP:CD2	2.52	0.44
1:B:69:LEU:HB2	1:B:111:CYS:O	2.17	0.44
1:B:8:TYR:CD1	1:B:36:ILE:HA	2.52	0.44
1:D:154:VAL:HA	1:D:188:HIS:CG	2.52	0.44
1:A:197:ARG:O	1:A:201:LEU:HG	2.18	0.44
1:D:259:VAL:HG23	1:D:261:VAL:HG23	2.00	0.44
3:A:402[A]:TGD:C1	3:A:402[A]:TGD:O4	2.66	0.43
1:D:26:ILE:HD12	1:D:34:MET:SD	2.58	0.43
1:B:249:MET:HE1	1:B:277:ARG:NH2	2.34	0.43
1:C:125:ARG:HB2	1:C:126:PRO:HD3	2.00	0.43
1:D:18:ASP:OD2	1:D:20:PRO:HD2	2.19	0.43
1:C:249:MET:HA	1:C:265:MET:HB2	2.02	0.42
1:D:128:VAL:O	1:D:132:ILE:HG13	2.19	0.42
1:D:8:TYR:CE1	1:D:9:THR:HG23	2.54	0.42
1:A:1:MET:HG3	1:A:2:ASN:N	2.35	0.42
1:C:35:GLU:HG3	1:C:64:MET:HG3	2.01	0.42
1:D:2:ASN:ND2	5:D:502:HOH:O	2.50	0.42
1:A:7:PHE:HB3	5:A:694:HOH:O	2.18	0.42
1:D:35:GLU:HG3	1:D:64:MET:HG3	2.00	0.42
1:A:154:VAL:CG2	1:A:159:GLN:HB2	2.50	0.42
1:A:99:LEU:HD23	1:A:99:LEU:O	2.20	0.42
1:D:276:GLU:HG3	1:D:280:ARG:HH12	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:249:MET:HA	1:B:265:MET:HB2	2.02	0.41
1:C:1:MET:HG2	5:C:578:HOH:O	2.21	0.41
1:B:50:ARG:HG2	1:B:50:ARG:NH1	2.36	0.41
1:C:266:SER:O	1:C:267:ASN:HB2	2.21	0.41
1:A:219:PRO:HA	1:A:220:PRO:HD3	1.96	0.41
1:D:25:ARG:NH2	1:D:271:ASP:HB3	2.35	0.41
1:D:283:GLN:NE2	1:D:287:ASP:OD2	2.52	0.41
1:B:152:GLU:HA	1:B:183:GLN:HB3	2.02	0.40
1:B:160:TRP:N	1:B:160:TRP:CD1	2.89	0.40
1:D:183:GLN:HG3	1:D:209:HIS:HB3	2.03	0.40
1:A:160:TRP:CD1	1:A:160:TRP:N	2.89	0.40
1:A:272:GLU:OE2	1:A:276:GLU:OE2	2.39	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	292/300 (97%)	287 (98%)	5 (2%)	0	100	100
1	B	290/300 (97%)	285 (98%)	5 (2%)	0	100	100
1	C	295/300 (98%)	288 (98%)	7 (2%)	0	100	100
1	D	288/300 (96%)	282 (98%)	6 (2%)	0	100	100
All	All	1165/1200 (97%)	1142 (98%)	23 (2%)	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	243/249 (98%)	241 (99%)	2 (1%)	81	82
1	B	241/249 (97%)	241 (100%)	0	100	100
1	C	246/249 (99%)	244 (99%)	2 (1%)	81	82
1	D	240/249 (96%)	238 (99%)	2 (1%)	81	82
All	All	970/996 (97%)	964 (99%)	6 (1%)	86	87

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	TYR
1	A	293	ARG
1	C	8	TYR
1	C	297	HIS
1	D	16	MET
1	D	122	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	267	ASN
1	C	267	ASN
1	D	98	HIS
1	D	267	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates

There are no monosaccharides in this entry.

5.6 Ligand geometry

Of 7 ligands modelled in this entry, 4 are monoatomic - leaving 3 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	TGD	A	402[B]	2	10,10,10	0.43	0	11,13,13	0.91	1 (9%)
3	TGD	A	402[A]	2	10,10,10	0.47	0	11,13,13	0.76	0
4	TGY	C	402	-	11,11,11	0.70	0	16,17,17	0.97	1 (6%)

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	TGD	A	402[B]	2	-	4/14/14/14	-
3	TGD	A	402[A]	2	-	0/14/14/14	-
4	TGY	C	402	-	-	-	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	402	TGY	O6-C2-C3	2.78	110.16	105.96
3	A	402[B]	TGD	C1-C2-C3	2.30	119.97	118.39

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	402[B]	TGD	C3-C4-C5-O5
3	A	402[B]	TGD	O4-C4-C5-O5
3	A	402[B]	TGD	C3-C4-C5-C6
3	A	402[B]	TGD	O4-C4-C5-C6

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	402[B]	TGD	1	0
3	A	402[A]	TGD	2	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 ⓘ

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	294/300 (98%)	-0.27	5 (1%) 70 72	9, 16, 27, 60	0
1	B	292/300 (97%)	-0.31	6 (2%) 63 66	12, 16, 27, 51	0
1	C	297/300 (99%)	-0.34	5 (1%) 70 72	11, 16, 27, 42	0
1	D	290/300 (96%)	-0.02	5 (1%) 70 72	13, 21, 33, 55	0
All	All	1173/1200 (97%)	-0.24	21 (1%) 68 71	9, 17, 30, 60	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1	MET	8.3
1	D	1	MET	7.9
1	A	294	SER	6.8
1	A	1	MET	6.2
1	B	292	SER	5.1
1	C	296	HIS	4.3
1	D	122	LYS	3.9
1	C	120	ASP	3.9
1	D	125	ARG	3.8
1	C	1	MET	3.3
1	A	119	LEU	2.9
1	C	122	LYS	2.9
1	D	120	ASP	2.7
1	A	122	LYS	2.7
1	A	2	ASN	2.7
1	B	122	LYS	2.6
1	C	297	HIS	2.4
1	B	119	LEU	2.2
1	B	240	ASP	2.2
1	B	120	ASP	2.1
1	D	16	MET	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	TGD	A	402[B]	11/11	0.71	0.28	41,44,47,47	11
3	TGD	A	402[A]	11/11	0.71	0.28	53,56,57,58	11
4	TGY	C	402	11/11	0.74	0.23	35,37,39,39	0
2	MN	A	401	1/1	0.99	0.03	16,16,16,16	0
2	MN	B	401	1/1	0.99	0.03	23,23,23,23	0
2	MN	D	401	1/1	1.00	0.02	20,20,20,20	0
2	MN	C	401	1/1	1.00	0.05	22,22,22,22	0

6.5 Other polymers [i](#)

There are no such residues in this entry.