



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:49 AM GMT

PDB ID : 3G4D
Title : Crystal Structure of (+)-delta-Cadinene Synthase from *Gossypium arboreum* and Evolutionary Divergence of Metal Binding Motifs for Catalysis
Authors : Gennadios,H.A.; Di Costanzo,L.; Miller, D.J.; Allemann, R.K.; Christianson,D.W.
Deposited on : 2009-02-03
Resolution : 2.40 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
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

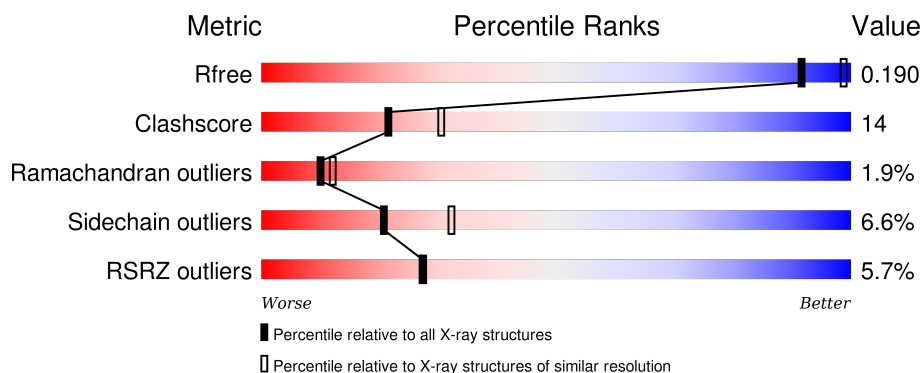
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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}	91344	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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. 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	554	
1	B	554	

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
2	BME	A	1274	-	-	-	X
3	GOL	B	2431	-	X	-	X
3	GOL	B	2432	-	X	-	X
3	GOL	B	2433	-	X	-	-
3	GOL	B	2434	-	X	-	-

2 Entry composition [i](#)

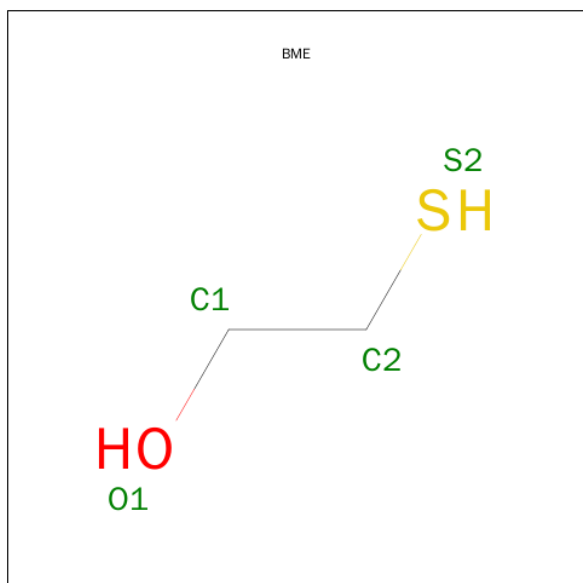
There are 4 unique types of molecules in this entry. The entry contains 8813 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 (+)-delta-cadinene synthase isozyme XC1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	516	Total	C	N	O	S	15	0	0
			4217	2694	702	799	22			
1	B	512	Total	C	N	O	S	8	0	0
			4178	2670	699	787	22			

- Molecule 2 is BETA-MERCAPTOETHANOL (three-letter code: BME) (formula: C_2H_6OS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	S	0	0
			4	2	1	1		
2	A	1	Total	C	O	S	0	0
			4	2	1	1		
2	A	1	Total	C	O	S	0	0
			4	2	1	1		

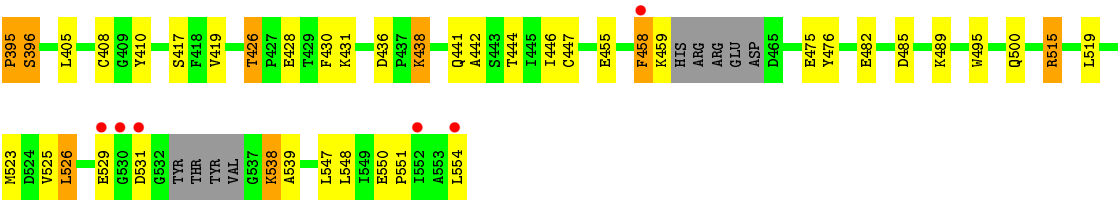
- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	278	Total	O	0	0
			278	278		
4	B	104	Total	O	0	0
			104	104		



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants a, b, c, α , β , γ	158.19Å 158.19Å 158.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.70 – 2.40 47.70 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.3 (47.70-2.40) 99.3 (47.70-2.40)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.05 (at 2.39Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.192 , 0.238 0.187 , 0.190	Depositor DCC
R_{free} test set	2605 reflections (5.09%)	DCC
Wilson B-factor (Å ²)	50.7	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 64.4	EDS
Estimated twinning fraction	0.026 for l,-k,h	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 51194 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8813	wwPDB-VP
Average B, all atoms (Å ²)	65.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.73% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: GOL, BME

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	0.31	0/4313	0.48	2/5835 (0.0%)
1	B	0.25	0/4271	0.43	0/5772
All	All	0.28	0/8584	0.45	2/11607 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	106	ASP	CB-CG-OD2	5.18	122.97	118.30
1	A	211	ARG	NE-CZ-NH1	5.02	122.81	120.30

There are no chirality outliers.

There are no planarity outliers.

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	4217	0	4103	105	0
1	B	4178	0	4088	131	0
2	A	12	0	18	5	0
3	B	24	0	24	0	0
4	A	278	0	0	7	0
4	B	104	0	0	7	0
All	All	8813	0	8233	234	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 14.

All (234) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:320:ILE:HG13	1:A:321:PRO:HD3	1.37	1.06
1:A:28:GLN:HB3	1:A:29:PRO:HA	1.41	1.02
1:B:222:GLN:HB3	1:B:223:ASP:CA	1.90	1.01
1:B:222:GLN:HB2	1:B:232:LEU:HD22	1.44	0.97
1:B:222:GLN:HB3	1:B:223:ASP:HA	1.49	0.95
1:A:392:ASN:HB2	1:B:140:GLN:HG3	1.50	0.93
1:A:28:GLN:HB3	1:A:29:PRO:CA	2.01	0.91
1:B:222:GLN:CB	1:B:223:ASP:HA	2.01	0.91
1:B:211:ARG:HG2	1:B:211:ARG:HH11	1.38	0.89
1:A:438:LYS:H	1:A:438:LYS:HE3	1.39	0.87
1:B:222:GLN:HB3	1:B:223:ASP:C	1.94	0.86
1:B:320:ILE:HG13	1:B:321:PRO:HD3	1.58	0.85
1:B:438:LYS:H	1:B:438:LYS:HE3	1.42	0.85
1:B:525:VAL:HG12	1:B:526:LEU:HD13	1.58	0.84
1:A:27:PHE:HD2	1:A:267:PRO:HA	1.43	0.83
1:A:276:GLY:HA2	1:A:279:TRP:CE3	2.13	0.83
1:A:211:ARG:HG2	1:A:211:ARG:HH11	1.44	0.81
1:A:438:LYS:HE3	1:A:438:LYS:N	1.94	0.81
1:B:211:ARG:CG	1:B:211:ARG:HH11	1.93	0.81
1:B:455:GLU:HG3	1:B:531:ASP:OD2	1.80	0.81
1:B:359:VAL:HG13	1:B:364:ARG:HB2	1.62	0.81
1:B:438:LYS:N	1:B:438:LYS:HE3	1.98	0.78
1:B:222:GLN:CB	1:B:223:ASP:CA	2.60	0.78
1:A:426:THR:HG23	1:A:428:GLU:H	1.48	0.77
1:A:27:PHE:HB2	1:A:28:GLN:HE21	1.50	0.77
1:A:320:ILE:HG13	1:A:321:PRO:CD	2.17	0.73
1:B:222:GLN:H	1:B:232:LEU:HD13	1.53	0.72
1:A:28:GLN:CB	1:A:29:PRO:HA	2.19	0.71
1:A:211:ARG:HG2	1:A:211:ARG:NH1	2.04	0.71
1:A:27:PHE:HB2	1:A:28:GLN:NE2	2.05	0.71
1:B:438:LYS:H	1:B:438:LYS:CE	2.05	0.70
1:A:359:VAL:HG13	1:A:364:ARG:HB2	1.73	0.69
1:B:458:PHE:O	1:B:459:LYS:HG3	1.93	0.69
1:B:114:THR:HG23	1:B:132:VAL:HG12	1.74	0.69
1:A:438:LYS:CE	1:A:438:LYS:H	2.06	0.68
1:B:135:LYS:HG2	4:B:645:HOH:O	1.93	0.67
1:A:211:ARG:CG	1:A:211:ARG:HH11	2.06	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:108:GLU:HB2	1:A:135:LYS:NZ	2.10	0.67
1:B:284:TYR:HB3	1:B:515:ARG:NH2	2.09	0.67
1:B:211:ARG:HG2	1:B:211:ARG:NH1	2.03	0.67
1:A:40:PRO:O	1:A:41:ASP:HB2	1.95	0.67
1:B:475:GLU:HG3	1:B:476:TYR:CD2	2.29	0.66
1:A:28:GLN:HG3	1:A:271:ASP:HB2	1.76	0.66
1:A:106:ASP:C	1:A:108:GLU:HA	2.16	0.66
1:A:69:ASN:ND2	1:A:72:GLN:H	1.94	0.66
1:A:328:ARG:NH2	1:A:333:CYS:SG	2.68	0.65
1:A:364:ARG:HD2	1:A:367:ARG:CZ	2.27	0.65
1:B:428:GLU:HA	1:B:431:LYS:HD3	1.79	0.65
1:B:164:ARG:HD2	1:B:172:ASP:OD1	1.97	0.65
1:B:426:THR:HG22	1:B:428:GLU:H	1.62	0.64
1:A:254:CYS:SG	4:A:623:HOH:O	2.55	0.64
1:B:60:ARG:O	1:B:64:VAL:HG23	1.98	0.64
1:B:226:SER:O	1:B:227:HIS:HB2	1.96	0.64
1:A:39:CYS:SG	2:A:1275:BME:S2	2.90	0.63
1:A:106:ASP:O	1:A:108:GLU:HA	1.97	0.63
1:B:364:ARG:HD3	4:B:635:HOH:O	1.99	0.63
1:A:276:GLY:HA2	1:A:279:TRP:HE3	1.63	0.62
1:A:27:PHE:CD2	1:A:267:PRO:HA	2.29	0.62
1:B:426:THR:CG2	1:B:428:GLU:H	2.13	0.62
1:A:40:PRO:HG3	1:A:247:ARG:HB3	1.82	0.62
1:A:426:THR:HG23	1:A:428:GLU:N	2.15	0.61
1:A:110:ASP:HB2	1:A:113:THR:H	1.64	0.61
1:B:320:ILE:HG13	1:B:321:PRO:CD	2.31	0.60
1:B:211:ARG:NH2	1:B:242:LEU:HD12	2.16	0.60
1:A:538:LYS:HD2	1:A:538:LYS:H	1.67	0.60
1:B:110:ASP:HB3	1:B:113:THR:H	1.66	0.60
1:A:107:ALA:HA	1:A:108:GLU:HB3	1.84	0.60
1:B:192:PRO:HG3	1:B:224:ILE:CG2	2.32	0.60
1:B:278:PHE:HE1	1:B:547:LEU:HD12	1.66	0.59
1:A:328:ARG:CZ	1:A:333:CYS:SG	2.91	0.59
1:A:535:TYR:N	1:A:536:VAL:HA	2.16	0.59
1:B:232:LEU:HD11	1:B:236:LYS:HE2	1.85	0.59
1:A:541:LYS:HE3	1:B:482:GLU:OE1	2.01	0.59
1:A:422:GLY:CA	2:A:1274:BME:H11	2.33	0.58
1:A:199:HIS:HA	1:A:202:LYS:NZ	2.19	0.57
1:B:165:VAL:CG1	1:B:166:HIS:N	2.68	0.57
1:A:274:VAL:HG12	1:A:544:ILE:HD13	1.87	0.57
1:B:180:HIS:CE1	1:B:184:LEU:HD11	2.39	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:169:ASP:O	1:A:173:GLU:HG3	2.04	0.56
1:B:222:GLN:NE2	1:B:227:HIS:O	2.39	0.56
1:B:364:ARG:HD2	1:B:367:ARG:CZ	2.35	0.56
1:A:417:SER:OG	1:A:515:ARG:HG2	2.04	0.56
1:B:221:TYR:C	1:B:221:TYR:CD1	2.79	0.56
1:B:396:SER:HB3	4:B:648:HOH:O	2.05	0.55
1:A:280:ILE:HD12	1:A:296:LEU:HG	1.87	0.55
1:B:426:THR:HG22	1:B:428:GLU:N	2.21	0.55
1:A:294:LYS:HG2	4:A:817:HOH:O	2.07	0.55
1:A:69:ASN:C	1:A:69:ASN:HD22	2.09	0.55
1:A:39:CYS:HG	2:A:1275:BME:HS2	1.33	0.55
1:A:108:GLU:HB2	1:A:135:LYS:HZ1	1.71	0.55
1:A:198:SER:O	1:A:202:LYS:HG3	2.07	0.55
1:A:106:ASP:HB3	4:A:761:HOH:O	2.07	0.55
1:A:169:ASP:OD1	1:A:169:ASP:N	2.38	0.55
1:B:225:GLU:O	1:B:225:GLU:HG2	2.06	0.55
1:B:538:LYS:HE3	1:B:538:LYS:HA	1.89	0.55
1:B:137:LYS:HE3	1:B:170:ILE:HG13	1.89	0.54
1:A:242:LEU:HD21	1:A:287:PRO:HG3	1.90	0.54
1:A:303:ALA:HB2	1:A:375:MET:HE1	1.89	0.54
1:B:225:GLU:O	1:B:225:GLU:CG	2.56	0.54
1:B:69:ASN:HD22	1:B:69:ASN:N	2.05	0.54
1:A:110:ASP:HB3	4:A:788:HOH:O	2.06	0.53
1:B:303:ALA:HB2	1:B:375:MET:HE1	1.91	0.53
1:A:69:ASN:HD22	1:A:72:GLN:H	1.56	0.53
1:B:59:VAL:O	1:B:63:ILE:HG13	2.09	0.53
1:B:395:PRO:O	1:B:396:SER:HB3	2.09	0.53
1:B:182:LEU:O	1:B:186:VAL:HG23	2.09	0.53
1:B:247:ARG:HD3	1:B:548:LEU:O	2.09	0.52
1:B:44:ILE:HD11	1:B:554:LEU:HD11	1.92	0.52
1:B:191:HIS:CG	1:B:192:PRO:HA	2.45	0.52
1:A:422:GLY:HA3	2:A:1274:BME:H11	1.92	0.52
1:A:334:ILE:O	1:A:342:LYS:HE2	2.09	0.52
1:B:222:GLN:HB2	1:B:223:ASP:HA	1.86	0.52
1:B:315:THR:OG1	1:B:318:GLU:HG3	2.09	0.52
1:A:185:ALA:O	1:A:189:LEU:HG	2.11	0.51
1:A:180:HIS:CE1	1:A:184:LEU:HD11	2.46	0.51
1:A:252:GLU:OE1	1:A:294:LYS:HE3	2.10	0.51
1:A:330:ASP:OD1	1:A:331:ILE:N	2.44	0.51
1:B:221:TYR:C	1:B:221:TYR:HD1	2.15	0.50
1:B:165:VAL:HG13	4:B:597:HOH:O	2.10	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:40:PRO:HG3	1:B:247:ARG:HB3	1.91	0.50
1:A:40:PRO:O	1:A:41:ASP:CB	2.59	0.50
1:B:295:MET:HE3	1:B:355:MET:HA	1.94	0.50
1:A:458:PHE:HB2	1:A:529:GLU:O	2.12	0.50
1:B:526:LEU:HD12	1:B:539:ALA:HB1	1.93	0.50
1:B:346:LYS:HG3	1:B:347:ALA:N	2.26	0.49
1:B:458:PHE:CD2	1:B:459:LYS:N	2.80	0.49
1:B:74:LEU:HD21	1:B:100:ILE:HG21	1.95	0.49
1:A:39:CYS:CB	2:A:1275:BME:HS2	2.25	0.49
1:B:279:TRP:HE3	1:B:300:ILE:HD11	1.78	0.49
1:B:118:PHE:O	1:B:122:ARG:HB2	2.13	0.49
1:B:256:TRP:O	1:B:260:LEU:HD23	2.12	0.49
1:B:287:PRO:HD2	4:B:619:HOH:O	2.12	0.49
1:A:191:HIS:CG	1:A:192:PRO:HA	2.48	0.48
1:A:109:ASN:HB3	1:A:113:THR:HB	1.94	0.48
1:B:114:THR:HG23	1:B:132:VAL:CG1	2.40	0.48
1:B:119:ARG:HG3	1:B:163:LEU:HD13	1.95	0.48
1:A:222:GLN:C	1:A:223:ASP:O	2.51	0.48
1:B:122:ARG:HD2	4:B:577:HOH:O	2.13	0.47
1:A:225:GLU:O	1:A:226:SER:CB	2.62	0.47
1:B:133:PHE:HB2	1:B:170:ILE:HD13	1.96	0.47
1:A:199:HIS:HA	1:A:202:LYS:HZ2	1.79	0.47
1:B:207:ARG:HG2	1:B:495:TRP:CH2	2.49	0.47
1:B:165:VAL:HG11	1:B:500:GLN:OE1	2.14	0.46
1:B:221:TYR:O	1:B:221:TYR:HD1	1.98	0.46
1:A:182:LEU:O	1:A:186:VAL:HG23	2.15	0.46
1:B:302:MET:HG3	1:B:351:VAL:HG21	1.97	0.46
1:B:302:MET:O	1:B:306:VAL:HG23	2.16	0.46
1:B:359:VAL:CG1	1:B:364:ARG:HB2	2.39	0.46
1:B:393:TYR:CE2	1:B:395:PRO:HA	2.50	0.46
1:A:447:CYS:HA	1:A:523:MET:HG3	1.98	0.46
1:B:417:SER:OG	1:B:515:ARG:HG2	2.15	0.46
1:B:447:CYS:HA	1:B:523:MET:HG3	1.97	0.46
1:A:398:GLU:HA	1:A:401:LYS:HB3	1.98	0.46
1:B:42:LYS:HD2	1:B:44:ILE:H	1.81	0.46
1:B:355:MET:HE3	1:B:419:VAL:HG11	1.97	0.46
1:B:41:ASP:O	1:B:42:LYS:HB3	2.16	0.46
1:B:129:SER:O	1:B:132:VAL:HG23	2.15	0.46
1:B:328:ARG:CZ	1:B:333:CYS:SG	3.04	0.46
1:B:52:HIS:HB2	1:B:237:ILE:HG21	1.97	0.45
1:B:436:ASP:OD2	1:B:441:GLN:HG3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:257:TRP:CD2	1:B:273:VAL:HG21	2.51	0.45
1:B:253:ILE:CG2	1:B:294:LYS:HG2	2.46	0.45
1:B:475:GLU:HG3	1:B:476:TYR:HD2	1.79	0.45
1:B:426:THR:CG2	1:B:428:GLU:OE1	2.65	0.45
1:B:458:PHE:CG	1:B:459:LYS:N	2.84	0.45
1:A:266:LEU:HA	1:A:267:PRO:HD2	1.80	0.45
1:B:191:HIS:CD2	1:B:224:ILE:HD13	2.52	0.45
1:B:111:LEU:HD22	1:B:147:VAL:HG13	1.99	0.45
1:A:257:TRP:CG	1:A:273:VAL:HG21	2.52	0.44
1:A:318:GLU:O	1:A:321:PRO:HD2	2.17	0.44
1:A:252:GLU:HB2	1:A:294:LYS:HE3	1.98	0.44
1:B:331:ILE:HG13	1:B:331:ILE:H	1.58	0.44
1:B:56:LYS:HG3	1:B:234:PHE:CZ	2.52	0.44
1:A:104:ASN:ND2	4:A:761:HOH:O	2.50	0.44
1:B:28:GLN:HB3	1:B:29:PRO:HD3	2.00	0.44
1:A:108:GLU:HB2	1:A:135:LYS:HZ2	1.79	0.44
1:B:138:ASP:HB3	1:B:144:LYS:HG2	2.00	0.44
1:B:485:ASP:O	1:B:489:LYS:HG3	2.18	0.44
1:A:45:ASP:HB3	1:A:48:THR:CG2	2.47	0.44
1:B:253:ILE:HD11	1:B:273:VAL:HG13	2.00	0.44
1:A:191:HIS:NE2	1:A:223:ASP:OD1	2.47	0.44
1:A:152:ARG:NH2	4:A:735:HOH:O	2.37	0.44
1:B:222:GLN:C	1:B:224:ILE:H	2.20	0.43
1:B:442:ALA:O	1:B:446:ILE:HG13	2.18	0.43
1:B:352:TYR:O	1:B:356:VAL:HG23	2.18	0.43
1:B:475:GLU:HG3	1:B:476:TYR:N	2.33	0.43
1:B:101:TYR:HD1	1:B:121:LEU:HD13	1.82	0.43
1:B:165:VAL:HG13	1:B:166:HIS:N	2.33	0.43
1:A:107:ALA:N	1:A:108:GLU:HA	2.29	0.43
1:B:295:MET:CE	1:B:355:MET:HA	2.48	0.43
1:B:410:TYR:CE2	1:B:519:LEU:HB3	2.53	0.43
1:B:278:PHE:CE1	1:B:547:LEU:HD12	2.52	0.43
1:B:44:ILE:HD13	1:B:237:ILE:HG23	1.99	0.42
1:B:70:SER:O	1:B:74:LEU:HG	2.19	0.42
1:B:100:ILE:HG22	1:B:100:ILE:O	2.20	0.42
1:A:437:PRO:HD2	1:A:440:ILE:HB	2.00	0.42
1:A:256:TRP:CE2	1:A:260:LEU:HD21	2.53	0.42
1:A:28:GLN:NE2	1:A:263:GLN:HA	2.35	0.42
1:A:27:PHE:HB3	1:A:266:LEU:O	2.19	0.42
1:A:191:HIS:ND1	1:A:192:PRO:HA	2.34	0.42
1:A:244:PHE:O	1:A:248:LYS:HG3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:318:GLU:C	1:A:321:PRO:HD2	2.39	0.42
1:B:395:PRO:HB2	1:B:396:SER:H	1.54	0.42
1:A:405:LEU:N	1:A:406:PRO:CD	2.82	0.42
1:B:526:LEU:HD12	1:B:526:LEU:HA	1.87	0.42
1:B:69:ASN:OD1	1:B:71:THR:HG22	2.19	0.42
1:A:256:TRP:O	1:A:260:LEU:HD22	2.20	0.42
1:B:405:LEU:HD22	1:B:444:THR:HG22	2.02	0.42
1:B:165:VAL:HG13	1:B:166:HIS:H	1.85	0.42
1:B:355:MET:CE	1:B:419:VAL:HG11	2.50	0.42
1:A:398:GLU:HA	1:A:401:LYS:CB	2.49	0.42
1:A:394:LYS:HA	1:A:395:PRO:HD3	1.89	0.41
1:A:134:ASN:HD22	1:A:134:ASN:HA	1.69	0.41
1:B:140:GLN:H	1:B:140:GLN:NE2	2.19	0.41
1:B:161:SER:O	1:B:164:ARG:HB3	2.21	0.41
1:A:39:CYS:HA	1:A:40:PRO:HD3	1.80	0.41
1:B:550:GLU:HA	1:B:551:PRO:HD2	1.77	0.41
1:A:364:ARG:HD3	4:A:683:HOH:O	2.20	0.41
1:B:257:TRP:CG	1:B:273:VAL:HG21	2.56	0.41
1:B:353:GLU:HG3	4:B:620:HOH:O	2.21	0.41
1:B:385:GLU:O	1:B:389:THR:HG23	2.21	0.41
1:A:217:TYR:HA	1:A:220:VAL:HG13	2.03	0.41
1:B:370:TYR:CD2	1:B:430:PHE:HB3	2.55	0.41
1:A:550:GLU:HA	1:A:551:PRO:HD2	1.94	0.41
1:A:68:ALA:HB3	1:A:72:GLN:CD	2.42	0.41
1:B:114:THR:CG2	1:B:132:VAL:HG12	2.46	0.40
1:B:328:ARG:NH2	1:B:333:CYS:SG	2.93	0.40
1:A:192:PRO:HG3	1:A:223:ASP:CB	2.50	0.40
1:A:27:PHE:CG	1:A:263:GLN:O	2.75	0.40
1:B:320:ILE:N	1:B:321:PRO:CD	2.85	0.40
1:A:109:ASN:HB3	1:A:113:THR:CG2	2.52	0.40
1:A:56:LYS:HB2	1:A:234:PHE:CE1	2.56	0.40
1:A:223:ASP:O	1:A:224:ILE:C	2.55	0.40
1:A:119:ARG:HG3	1:A:163:LEU:HD13	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	508/554 (92%)	489 (96%)	12 (2%)	7 (1%)	14	19
1	B	502/554 (91%)	467 (93%)	23 (5%)	12 (2%)	7	7
All	All	1010/1108 (91%)	956 (95%)	35 (4%)	19 (2%)	10	12

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	223	ASP
1	A	457	LYS
1	B	395	PRO
1	B	396	SER
1	B	458	PHE
1	A	28	GLN
1	A	110	ASP
1	B	110	ASP
1	B	225	GLU
1	B	226	SER
1	B	529	GLU
1	A	26	ASP
1	A	466	CYS
1	B	42	LYS
1	B	392	ASN
1	A	224	ILE
1	B	40	PRO
1	B	227	HIS
1	B	228	ASN

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	457/492 (93%)	427 (93%)	30 (7%)	21	32
1	B	453/492 (92%)	423 (93%)	30 (7%)	21	32
All	All	910/984 (92%)	850 (93%)	60 (7%)	21	32

All (60) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	GLN
1	A	35	LEU
1	A	37	LEU
1	A	48	THR
1	A	69	ASN
1	A	110	ASP
1	A	164	ARG
1	A	169	ASP
1	A	206	ARG
1	A	211	ARG
1	A	220	VAL
1	A	222	GLN
1	A	225	GLU
1	A	231	LEU
1	A	258	LYS
1	A	260	LEU
1	A	284	TYR
1	A	291	LEU
1	A	320	ILE
1	A	359	VAL
1	A	378	LEU
1	A	383	LEU
1	A	392	ASN
1	A	426	THR
1	A	431	LYS
1	A	438	LYS
1	A	457	LYS
1	A	515	ARG
1	A	526	LEU
1	A	535	TYR
1	B	36	PHE
1	B	37	LEU
1	B	42	LYS

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Mol	Chain	Res	Type
1	B	48	THR
1	B	69	ASN
1	B	139	GLU
1	B	165	VAL
1	B	206	ARG
1	B	211	ARG
1	B	220	VAL
1	B	221	TYR
1	B	223	ASP
1	B	225	GLU
1	B	231	LEU
1	B	253	ILE
1	B	264	ARG
1	B	284	TYR
1	B	291	LEU
1	B	320	ILE
1	B	357	GLN
1	B	378	LEU
1	B	383	LEU
1	B	392	ASN
1	B	394	LYS
1	B	408	CYS
1	B	426	THR
1	B	438	LYS
1	B	515	ARG
1	B	526	LEU
1	B	538	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	28	GLN
1	A	69	ASN
1	A	134	ASN
1	A	180	HIS
1	A	181	HIS
1	A	222	GLN
1	A	441	GLN
1	B	28	GLN
1	B	69	ASN
1	B	72	GLN
1	B	134	ASN

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Mol	Chain	Res	Type
1	B	180	HIS
1	B	191	HIS
1	B	357	GLN
1	B	392	ASN
1	B	441	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

7 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	BME	A	1273	-	3,3,3	0.26	0	2,2,2	0.38	0
2	BME	A	1274	-	3,3,3	0.38	0	2,2,2	0.42	0
2	BME	A	1275	-	3,3,3	0.22	0	2,2,2	0.40	0
3	GOL	B	2431	-	5,5,5	4.89	3 (60%)	5,5,5	2.31	3 (60%)
3	GOL	B	2432	-	5,5,5	5.01	3 (60%)	5,5,5	2.22	4 (80%)
3	GOL	B	2433	-	5,5,5	4.87	3 (60%)	5,5,5	2.31	4 (80%)
3	GOL	B	2434	-	5,5,5	4.96	3 (60%)	5,5,5	2.23	5 (100%)

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
2	BME	A	1273	-	-	0/1/1/1	0/0/0/0
2	BME	A	1274	-	-	0/1/1/1	0/0/0/0
2	BME	A	1275	-	-	0/1/1/1	0/0/0/0
3	GOL	B	2431	-	-	0/4/4/4	0/0/0/0
3	GOL	B	2432	-	-	0/4/4/4	0/0/0/0
3	GOL	B	2433	-	-	0/4/4/4	0/0/0/0
3	GOL	B	2434	-	-	0/4/4/4	0/0/0/0

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	2432	GOL	O2-C2	-8.30	1.18	1.43
3	B	2434	GOL	O2-C2	-8.24	1.18	1.43
3	B	2433	GOL	O2-C2	-8.18	1.19	1.43
3	B	2431	GOL	O2-C2	-8.06	1.19	1.43
3	B	2432	GOL	C3-C2	-5.36	1.31	1.52
3	B	2434	GOL	C3-C2	-5.29	1.32	1.52
3	B	2432	GOL	C1-C2	-5.26	1.32	1.52
3	B	2431	GOL	C3-C2	-5.25	1.32	1.52
3	B	2434	GOL	C1-C2	-5.17	1.32	1.52
3	B	2431	GOL	C1-C2	-5.16	1.32	1.52
3	B	2433	GOL	C3-C2	-5.10	1.32	1.52
3	B	2433	GOL	C1-C2	-5.07	1.32	1.52

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	2434	GOL	O1-C1-C2	2.01	119.94	110.18
3	B	2434	GOL	O3-C3-C2	2.05	120.15	110.18
3	B	2432	GOL	O1-C1-C2	2.06	120.15	110.18
3	B	2432	GOL	C3-C2-C1	2.14	119.50	111.12
3	B	2434	GOL	C3-C2-C1	2.19	119.72	111.12
3	B	2433	GOL	O2-C2-C1	2.29	119.14	108.65
3	B	2431	GOL	O1-C1-C2	2.31	121.39	110.18
3	B	2433	GOL	O3-C3-C2	2.31	121.41	110.18
3	B	2433	GOL	O2-C2-C3	2.40	119.64	108.65
3	B	2434	GOL	O2-C2-C3	2.42	119.73	108.65
3	B	2432	GOL	O2-C2-C3	2.42	119.73	108.65

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	2431	GOL	O2-C2-C3	2.42	119.75	108.65
3	B	2434	GOL	O2-C2-C1	2.45	119.86	108.65
3	B	2433	GOL	C3-C2-C1	2.57	121.21	111.12
3	B	2432	GOL	O2-C2-C1	2.60	120.59	108.65
3	B	2431	GOL	O2-C2-C1	2.74	121.23	108.65

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1274	BME	2	0
2	A	1275	BME	3	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	515/554 (92%)	-0.07	11 (2%) 67 66	31, 50, 85, 126	2 (0%)
1	B	511/554 (92%)	0.30	47 (9%) 11 11	42, 74, 120, 146	0
All	All	1026/1108 (92%)	0.11	58 (5%) 27 27	31, 60, 112, 146	2 (0%)

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	268	TYR	9.1
1	A	458	PHE	8.8
1	A	107	ALA	7.6
1	B	44	ILE	6.6
1	A	25	ALA	6.6
1	B	269	ALA	6.2
1	B	29	PRO	6.1
1	B	43	ASN	5.3
1	B	267	PRO	5.2
1	B	531	ASP	5.1
1	A	535	TYR	4.6
1	B	142	ASN	4.5
1	B	138	ASP	4.2
1	B	264	ARG	4.2
1	B	141	GLY	4.2
1	B	263	GLN	4.1
1	B	313	TYR	4.1
1	B	140	GLN	3.6
1	A	27	PHE	3.6
1	B	32	TRP	3.6
1	B	232	LEU	3.6
1	B	50	LYS	3.5
1	B	237	ILE	3.5
1	A	139	GLU	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	143	PHE	3.4
1	B	137	LYS	3.3
1	A	536	VAL	3.3
1	B	530	GLY	3.2
1	B	270	ARG	3.2
1	B	147	VAL	3.2
1	B	47	GLU	3.0
1	B	28	GLN	3.0
1	A	137	LYS	2.9
1	B	136	PHE	2.9
1	B	221	TYR	2.8
1	B	48	THR	2.8
1	B	223	ASP	2.8
1	B	458	PHE	2.8
1	B	143	PHE	2.8
1	B	266	LEU	2.8
1	A	140	GLN	2.7
1	B	233	GLU	2.7
1	B	257	TRP	2.6
1	B	53	GLN	2.6
1	B	554	LEU	2.5
1	B	63	ILE	2.4
1	B	51	ARG	2.4
1	B	392	ASN	2.4
1	B	529	GLU	2.4
1	B	103	ASN	2.3
1	B	265	LYS	2.3
1	A	279	TRP	2.3
1	B	552	ILE	2.2
1	B	139	GLU	2.2
1	B	65	ALA	2.2
1	B	133	PHE	2.2
1	B	330	ASP	2.2
1	B	229	LYS	2.0

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 [i](#)

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
3	GOL	B	2431	6/6	0.78	0.21	3.48	69,76,88,91	0
2	BME	A	1274	4/4	0.71	0.28	2.92	75,76,78,88	0
3	GOL	B	2432	6/6	0.96	0.18	2.64	67,76,81,82	0
3	GOL	B	2433	6/6	0.65	0.21	0.67	70,90,98,98	0
2	BME	A	1273	4/4	0.94	0.16	0.55	73,75,75,76	0
3	GOL	B	2434	6/6	0.96	0.14	-0.43	80,84,92,98	0
2	BME	A	1275	4/4	0.92	0.14	-	69,71,73,100	0

6.5 Other polymers [i](#)

There are no such residues in this entry.