



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:19 PM GMT

PDB ID : 1T3T
Title : Structure of Formylglycinamide synthetase
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Deposited on : 2004-04-27
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
<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

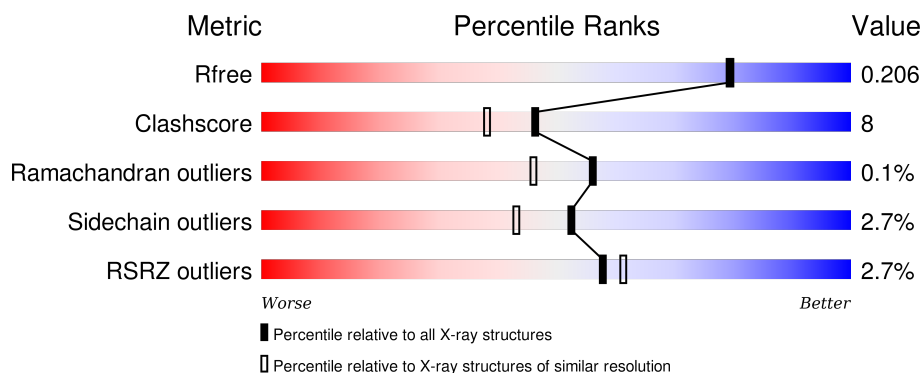
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	91344	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (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. 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	1303	<div> <div>3%</div> <div>82%</div> <div>15%</div> <div>..</div> </div>

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	MG	A	2006	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	MG	A	2007	-	-	-	X
4	ADP	A	2005	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 11024 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 Phosphoribosylformylglycinamidine synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1284	Total	C	N	O	S	0	0	0
			9885	6204	1761	1872	48			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	GLY	-	CLONING ARTIFACT	UNP P74881
A	-6	LEU	-	CLONING ARTIFACT	UNP P74881
A	-5	VAL	-	CLONING ARTIFACT	UNP P74881
A	-4	PRO	-	CLONING ARTIFACT	UNP P74881
A	-3	ARG	-	CLONING ARTIFACT	UNP P74881
A	-2	GLY	-	CLONING ARTIFACT	UNP P74881
A	-1	SER	-	CLONING ARTIFACT	UNP P74881
A	0	HIS	-	CLONING ARTIFACT	UNP P74881
A	1135	CYG	CYS	MODIFIED RESIDUE	UNP P74881

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	3	Total	Mg	0	0
			3	3		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

- Molecule 4 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

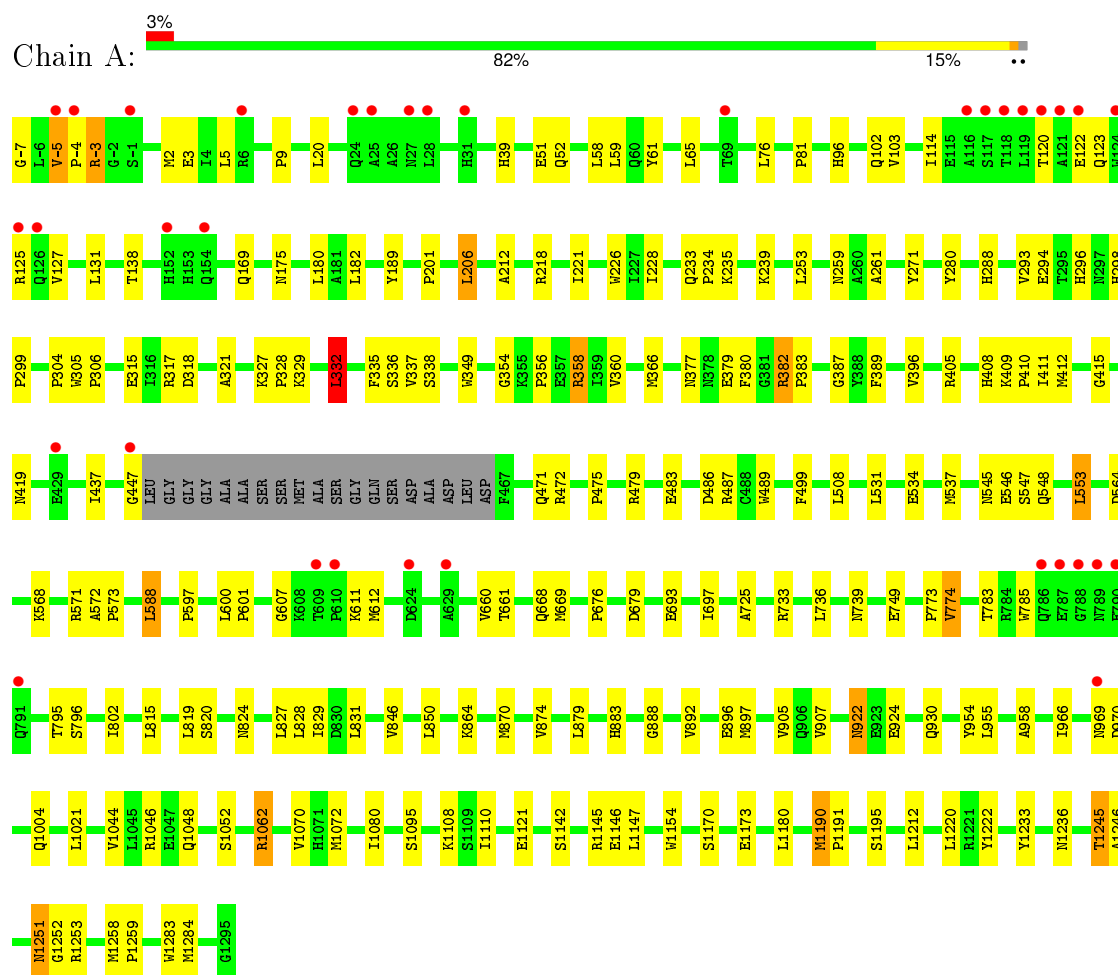
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1079	Total	O	0	0
			1079	1079		

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 errors 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: Phosphoribosylformylglycinamide synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	145.81Å 145.81Å 141.35Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	47.73 – 1.90 47.73 – 1.80	Depositor EDS
% Data completeness (in resolution range)	95.4 (47.73-1.90) 95.8 (47.73-1.80)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.57 (at 1.79Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.188 , 0.206 0.188 , 0.206	Depositor DCC
R_{free} test set	10822 reflections (8.48%)	DCC
Wilson B-factor (Å ²)	20.0	Xtriage
Anisotropy	0.069	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 47.9	EDS
Estimated twinning fraction	0.027 for h,-h-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 150565 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11024	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.29% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG, CYG, SO4, ADP

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	0.33	0/10082	0.62	3/13686 (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	332	LEU	CA-CB-CG	7.31	132.12	115.30
1	A	409	LYS	N-CA-C	-5.78	95.39	111.00
1	A	1236	ASN	N-CA-C	-5.29	96.72	111.00

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	9885	0	9662	149	0
2	A	3	0	0	0	0
3	A	30	0	0	0	0
4	A	27	0	12	1	0
5	A	1079	0	0	25	0
All	All	11024	0	9674	149	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 8.

All (149) 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:175:ASN:HD21	1:A:182:LEU:H	1.16	0.93
1:A:820:SER:H	1:A:930:GLN:HE22	1.16	0.91
1:A:824:ASN:HD21	1:A:958:ALA:H	1.15	0.88
1:A:1190:MET:HE2	1:A:1190:MET:H	1.46	0.80
1:A:358:ARG:HD3	1:A:358:ARG:H	1.55	0.71
1:A:396:VAL:CG1	1:A:850:LEU:HD22	2.22	0.70
1:A:733:ARG:HB3	5:A:2810:HOH:O	1.90	0.70
1:A:396:VAL:HG11	1:A:850:LEU:HD22	1.74	0.69
1:A:305:TRP:HB3	1:A:306:PRO:HD3	1.73	0.69
1:A:545:ASN:HD22	1:A:547:SER:H	1.40	0.69
1:A:922:ASN:HD22	1:A:924:GLU:H	1.41	0.68
1:A:175:ASN:HD22	1:A:180:LEU:HB2	1.59	0.67
1:A:321:ALA:HB3	5:A:2828:HOH:O	1.93	0.67
1:A:1251:ASN:ND2	1:A:1253:ARG:H	1.92	0.67
1:A:471:GLN:HE21	1:A:472:ARG:H	1.41	0.67
1:A:437:ILE:HB	1:A:553:LEU:HD12	1.78	0.66
1:A:1052:SER:HB3	5:A:2537:HOH:O	1.94	0.66
1:A:534:GLU:HG3	1:A:537:MET:HG3	1.78	0.65
1:A:669:MET:SD	1:A:676:PRO:HG3	2.36	0.65
1:A:5:LEU:HD22	1:A:59:LEU:HD12	1.77	0.65
1:A:175:ASN:ND2	1:A:182:LEU:H	1.94	0.64
1:A:796:SER:O	5:A:2760:HOH:O	2.15	0.64
1:A:337:VAL:HG13	1:A:366:MET:HE1	1.81	0.63
1:A:76:LEU:HD23	5:A:2982:HOH:O	1.98	0.63
1:A:114:ILE:HD11	1:A:131:LEU:HD11	1.82	0.62
1:A:487:ARG:NH2	5:A:2846:HOH:O	2.33	0.61
1:A:-5:VAL:HG12	5:A:2587:HOH:O	2.01	0.61
1:A:317:ARG:HH22	1:A:548:GLN:NE2	1.99	0.60
1:A:356:PRO:HG3	1:A:785:TRP:HB3	1.83	0.60
1:A:1222:TYR:OH	1:A:1245:THR:HG21	2.01	0.60
1:A:358:ARG:HD3	1:A:358:ARG:N	2.17	0.59
1:A:317:ARG:HH22	1:A:548:GLN:HE22	1.50	0.59
1:A:827:LEU:HD23	1:A:954:TYR:HA	1.83	0.59
1:A:221:ILE:HD13	1:A:612:MET:SD	2.43	0.58
1:A:228:ILE:HD13	1:A:233:GLN:HG3	1.86	0.58
1:A:259:ASN:HB2	5:A:2299:HOH:O	2.03	0.58
1:A:1146:GLU:HG3	1:A:1147:LEU:HD13	1.85	0.58
1:A:354:GLY:O	1:A:408:HIS:HE1	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:52:GLN:NE2	5:A:2964:HOH:O	2.37	0.57
1:A:1190:MET:HE2	1:A:1190:MET:N	2.16	0.56
1:A:259:ASN:ND2	5:A:2828:HOH:O	2.38	0.56
1:A:337:VAL:HG22	1:A:366:MET:HE2	1.87	0.56
1:A:261:ALA:HB2	1:A:774:VAL:HG13	1.87	0.56
1:A:335:PHE:HB3	1:A:366:MET:CE	2.35	0.56
1:A:96:HIS:HE1	1:A:103:VAL:O	1.89	0.56
1:A:1245:THR:HG22	1:A:1258:MET:HB2	1.88	0.56
1:A:336:SER:HB2	1:A:411:ILE:HB	1.88	0.56
1:A:1044:VAL:HG12	5:A:3079:HOH:O	2.05	0.55
1:A:1251:ASN:HD22	1:A:1251:ASN:C	2.08	0.55
1:A:1004:GLN:NE2	1:A:1233:TYR:H	2.04	0.55
1:A:39:HIS:HE1	1:A:61:TYR:OH	1.90	0.54
1:A:1070:VAL:HG22	1:A:1080:ILE:HD13	1.88	0.54
1:A:380:PHE:O	1:A:479:ARG:HD3	2.07	0.54
1:A:358:ARG:CD	1:A:358:ARG:H	2.15	0.53
1:A:379:GLU:HB3	1:A:475:PRO:HB2	1.90	0.53
1:A:405:ARG:NH2	5:A:2760:HOH:O	2.41	0.53
1:A:-3:ARG:HG3	1:A:2:MET:HB2	1.91	0.53
1:A:377:ASN:OD1	1:A:382:ARG:HD2	2.08	0.53
1:A:1251:ASN:HD22	1:A:1253:ARG:H	1.56	0.52
1:A:366:MET:HE3	1:A:389:PHE:CD1	2.46	0.51
1:A:120:THR:HG23	1:A:123:GLN:HE21	1.74	0.51
1:A:332:LEU:HG	1:A:697:ILE:HD11	1.93	0.51
1:A:905:VAL:HG13	1:A:907:VAL:HG13	1.91	0.51
1:A:668:GLN:HG2	4:A:2005:ADP:HI'	1.92	0.51
1:A:679:ASP:OD2	1:A:883:HIS:HD2	1.95	0.50
1:A:201:PRO:HB2	1:A:206:LEU:HD13	1.94	0.50
1:A:315:GLU:HG2	5:A:2833:HOH:O	2.12	0.49
1:A:58:LEU:HD23	1:A:1108:LYS:HD3	1.95	0.49
1:A:356:PRO:HG2	1:A:783:THR:CG2	2.43	0.49
1:A:1190:MET:CE	1:A:1190:MET:N	2.76	0.49
1:A:888:GLY:HA3	1:A:892:VAL:HG21	1.95	0.49
1:A:483:GLU:O	1:A:487:ARG:HG2	2.13	0.48
1:A:531:LEU:HD12	5:A:2814:HOH:O	2.13	0.48
1:A:1021:LEU:HD13	5:A:2907:HOH:O	2.12	0.48
1:A:829:ILE:HB	5:A:3018:HOH:O	2.13	0.48
1:A:-7:GLY:O	1:A:3:GLU:HA	2.14	0.48
1:A:486:ASP:HA	1:A:489:TRP:NE1	2.29	0.48
1:A:366:MET:HE2	1:A:389:PHE:HE1	1.79	0.47
1:A:564:ASP:O	1:A:568:LYS:HG2	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1251:ASN:HD22	1:A:1252:GLY:N	2.13	0.47
1:A:1220:LEU:HB2	1:A:1246:ALA:HB3	1.97	0.47
1:A:226:TRP:HB3	1:A:228:ILE:CD1	2.45	0.46
1:A:122:GLU:HG2	1:A:125:ARG:NH1	2.30	0.46
1:A:819:LEU:HG	1:A:897:MET:HB3	1.98	0.46
1:A:114:ILE:HD13	1:A:127:VAL:HG11	1.98	0.46
1:A:-5:VAL:HG13	1:A:-4:PRO:HD2	1.97	0.46
1:A:883:HIS:HE1	1:A:896:GLU:OE1	1.99	0.46
1:A:828:LEU:HB2	1:A:955:LEU:HD11	1.97	0.45
1:A:1072:MET:HE1	1:A:1110:ILE:HG12	1.99	0.45
1:A:293:VAL:HB	1:A:739:ASN:HD21	1.82	0.45
1:A:1212:LEU:HD22	5:A:2531:HOH:O	2.17	0.45
1:A:298:HIS:HB3	1:A:299:PRO:HD3	1.99	0.45
1:A:1170:SER:O	1:A:1191:PRO:HA	2.17	0.45
1:A:736:LEU:HG	1:A:802:ILE:HG23	1.99	0.45
1:A:499:PHE:HA	5:A:2968:HOH:O	2.17	0.45
1:A:545:ASN:ND2	1:A:547:SER:H	2.13	0.44
1:A:749:GLU:CD	5:A:2760:HOH:O	2.55	0.44
1:A:1245:THR:HG22	1:A:1259:PRO:HD3	2.00	0.44
1:A:1145:ARG:HD3	1:A:1154:TRP:HB2	1.99	0.44
1:A:20:LEU:HD11	5:A:2906:HOH:O	2.17	0.44
1:A:437:ILE:HB	1:A:553:LEU:CD1	2.46	0.44
1:A:212:ALA:HB2	5:A:2535:HOH:O	2.17	0.44
1:A:588:LEU:O	1:A:597:PRO:HD2	2.18	0.43
1:A:335:PHE:HB3	1:A:366:MET:HE3	2.00	0.43
1:A:338:SER:OG	1:A:408:HIS:HD2	2.02	0.43
1:A:81:PRO:HD3	1:A:138:THR:HG21	1.99	0.43
1:A:9:PRO:HB2	1:A:65:LEU:HB2	2.00	0.43
1:A:366:MET:CE	1:A:389:PHE:CE1	3.01	0.43
1:A:1222:TYR:CZ	1:A:1245:THR:HG21	2.53	0.43
1:A:387:GLY:HA2	1:A:697:ILE:HD12	2.01	0.43
1:A:1142:SER:O	1:A:1145:ARG:HG3	2.19	0.43
1:A:271:TYR:CZ	1:A:280:TYR:HB3	2.54	0.43
1:A:233:GLN:HA	1:A:234:PRO:HD3	1.80	0.42
1:A:189:TYR:OH	1:A:607:GLY:HA3	2.18	0.42
1:A:736:LEU:O	1:A:773:PRO:HD2	2.18	0.42
1:A:123:GLN:O	1:A:127:VAL:HG23	2.19	0.42
1:A:969:ASN:CG	1:A:970:ASP:H	2.21	0.42
1:A:228:ILE:HD13	1:A:233:GLN:CG	2.49	0.42
1:A:1121:GLU:HB2	1:A:1147:LEU:HG	2.02	0.42
1:A:1048:GLN:O	1:A:1095:SER:HA	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:349:TRP:CZ3	1:A:846:VAL:HG22	2.55	0.42
1:A:1062:ARG:HB3	1:A:1284:MET:CE	2.50	0.42
1:A:337:VAL:CG2	1:A:366:MET:HE2	2.48	0.41
1:A:611:LYS:NZ	1:A:611:LYS:HB2	2.35	0.41
1:A:1190:MET:HE1	1:A:1283:TRP:CZ2	2.54	0.41
1:A:304:PRO:HG3	1:A:360:VAL:HG21	2.02	0.41
1:A:318:ASP:HA	5:A:2828:HOH:O	2.20	0.41
1:A:1046:ARG:HB3	5:A:3079:HOH:O	2.18	0.41
1:A:815:LEU:HD13	1:A:879:LEU:HD12	2.02	0.41
1:A:486:ASP:HA	1:A:489:TRP:CD1	2.55	0.41
1:A:1190:MET:CE	1:A:1283:TRP:CZ2	3.03	0.41
1:A:2:MET:HE2	5:A:3033:HOH:O	2.20	0.41
1:A:235:LYS:HD3	1:A:239:LYS:HG3	2.03	0.41
1:A:572:ALA:HA	1:A:573:PRO:HD3	1.94	0.41
1:A:487:ARG:HA	1:A:487:ARG:HD3	1.81	0.41
1:A:383:PRO:HD3	1:A:660:VAL:HG13	2.03	0.41
1:A:51:GLU:CD	1:A:102:GLN:HE22	2.24	0.41
1:A:175:ASN:ND2	1:A:180:LEU:HB2	2.32	0.40
1:A:405:ARG:HA	1:A:795:THR:O	2.21	0.40
1:A:725:ALA:HB1	1:A:815:LEU:HB2	2.03	0.40
1:A:288:HIS:O	1:A:419:ASN:HA	2.21	0.40
1:A:329:LYS:HE2	1:A:693:GLU:OE1	2.21	0.40
1:A:447:GLY:CA	1:A:546:GLU:H	2.34	0.40
1:A:332:LEU:HD22	1:A:415:GLY:CA	2.51	0.40
1:A:870:MET:O	1:A:874:VAL:HG23	2.20	0.40
1:A:366:MET:HE2	1:A:389:PHE:CE1	2.56	0.40
1:A:327:LYS:HA	1:A:328:PRO:HD3	1.94	0.40
1:A:966:ILE:HG13	5:A:3046:HOH:O	2.21	0.40
1:A:600:LEU:HA	1:A:601:PRO:HD3	1.96	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	1279/1303 (98%)	1239 (97%)	39 (3%)	1 (0%)	56 46

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	661	THR

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	1026/1040 (99%)	998 (97%)	28 (3%)	52 43

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-5	VAL
1	A	-3	ARG
1	A	169	GLN
1	A	206	LEU
1	A	218	ARG
1	A	253	LEU
1	A	294	GLU
1	A	296	HIS
1	A	332	LEU
1	A	358	ARG
1	A	382	ARG
1	A	410	PRO
1	A	412	MET
1	A	508	LEU
1	A	553	LEU
1	A	571	ARG
1	A	588	LEU
1	A	774	VAL
1	A	831	LEU
1	A	864	LYS

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Mol	Chain	Res	Type
1	A	922	ASN
1	A	1062	ARG
1	A	1173	GLU
1	A	1180	LEU
1	A	1190	MET
1	A	1195	SER
1	A	1245	THR
1	A	1251	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	HIS
1	A	39	HIS
1	A	96	HIS
1	A	123	GLN
1	A	175	ASN
1	A	191	GLN
1	A	216	HIS
1	A	233	GLN
1	A	259	ASN
1	A	298	HIS
1	A	339	ASN
1	A	408	HIS
1	A	419	ASN
1	A	445	ASN
1	A	471	GLN
1	A	545	ASN
1	A	548	GLN
1	A	674	GLN
1	A	739	ASN
1	A	818	GLN
1	A	824	ASN
1	A	835	HIS
1	A	883	HIS
1	A	922	ASN
1	A	930	GLN
1	A	993	GLN
1	A	1004	GLN
1	A	1006	HIS
1	A	1018	ASN
1	A	1026	ASN

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Mol	Chain	Res	Type
1	A	1189	GLN
1	A	1251	ASN
1	A	1260	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	$\# Z > 2$
1	CYG	A	1135	1	10,14,15	0.61	0	6,17,19	6.02	4 (66%)

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
1	CYG	A	1135	1	-	0/10/16/18	0/0/0/0

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1135	CYG	OE2-CD1-CG1	-11.56	115.99	123.94
1	A	1135	CYG	CB1-CG1-CD1	-2.85	108.90	113.12
1	A	1135	CYG	O-C-CA	-2.12	119.98	125.49
1	A	1135	CYG	CG1-CD1-SG	8.44	120.95	113.36

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 3 are monoatomic - leaving 7 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	SO4	A	2001	-	4,4,4	0.26	0	6,6,6	0.12	0
3	SO4	A	2002	-	4,4,4	0.29	0	6,6,6	0.18	0
4	ADP	A	2005	2	22,29,29	1.00	0	27,45,45	1.77	3 (11%)
3	SO4	A	2009	-	4,4,4	0.25	0	6,6,6	0.12	0
3	SO4	A	2010	-	4,4,4	0.30	0	6,6,6	0.20	0
3	SO4	A	2011	-	4,4,4	0.26	0	6,6,6	0.12	0
3	SO4	A	2012	-	4,4,4	0.30	0	6,6,6	0.20	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	SO4	A	2001	-	-	0/0/0/0	0/0/0/0
3	SO4	A	2002	-	-	0/0/0/0	0/0/0/0
4	ADP	A	2005	2	-	0/12/32/32	0/3/3/3
3	SO4	A	2009	-	-	0/0/0/0	0/0/0/0
3	SO4	A	2010	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	SO4	A	2011	-	-	0/0/0/0	0/0/0/0
3	SO4	A	2012	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	2005	ADP	N3-C2-N1	-6.47	123.94	128.89
4	A	2005	ADP	C4-C5-N7	2.37	111.66	109.48
4	A	2005	ADP	C2-N1-C6	3.69	125.36	118.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	2005	ADP	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

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	1283/1303 (98%)	-0.06	35 (2%) 58 61	10, 20, 40, 67	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	118	THR	8.3
1	A	121	ALA	6.4
1	A	789	ASN	6.1
1	A	117	SER	5.4
1	A	790	GLU	4.7
1	A	120	THR	4.5
1	A	122	GLU	4.3
1	A	447	GLY	4.0
1	A	125	ARG	3.9
1	A	6	ARG	3.9
1	A	25	ALA	3.7
1	A	28	LEU	3.7
1	A	24	GLN	3.6
1	A	27	ASN	3.5
1	A	119	LEU	3.3
1	A	969	ASN	3.2
1	A	788	GLY	3.1
1	A	31	HIS	2.9
1	A	429	GLU	2.9
1	A	791	GLN	2.9
1	A	787	GLU	2.8
1	A	116	ALA	2.8
1	A	-1	SER	2.7
1	A	124	TRP	2.7
1	A	154	GLN	2.6
1	A	-5	VAL	2.5
1	A	69	THR	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	610	PRO	2.4
1	A	126	GLN	2.3
1	A	609	THR	2.3
1	A	152	HIS	2.3
1	A	786	GLN	2.3
1	A	624	ASP	2.2
1	A	-4	PRO	2.2
1	A	629	ALA	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [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(Å ²)	Q<0.9
1	CYG	A	1135	15/16	0.96	0.11	-	13,15,16,18	0

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(Å ²)	Q<0.9
4	ADP	A	2005	27/27	0.97	0.41	11.35	1,1,3,3	27
2	MG	A	2006	1/1	0.99	0.14	7.86	15,15,15,15	0
2	MG	A	2007	1/1	0.97	0.16	2.11	21,21,21,21	0
3	SO4	A	2009	5/5	0.97	0.10	1.28	38,39,40,41	0
3	SO4	A	2010	5/5	0.95	0.13	0.60	27,29,31,31	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	SO4	A	2001	5/5	0.95	0.10	-0.31	41,41,43,44	0
3	SO4	A	2002	5/5	0.97	0.07	-1.67	23,25,27,27	0
3	SO4	A	2012	5/5	0.98	0.09	-	32,32,33,33	0
2	MG	A	2008	1/1	0.98	0.18	-	19,19,19,19	0
3	SO4	A	2011	5/5	0.97	0.13	-	41,42,43,43	0

6.5 Other polymers [i](#)

There are no such residues in this entry.