



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:51 AM GMT

PDB ID : 2OI7
Title : E. coli GlmU- Complex with UDP-GlcNAc, desulpho-CoA and GlcNAc-1-PO4
Authors : Olsen, L.R.; Vetting, M.W.; Roderick, S.L.
Deposited on : 2007-01-10
Resolution : 2.54 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

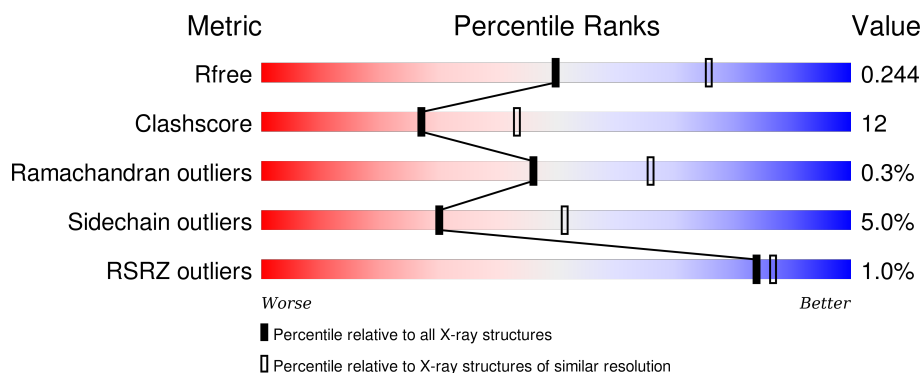
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 2.54 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	4549 (2.58-2.50)
Clashscore	102246	5292 (2.58-2.50)
Ramachandran outliers	100387	5194 (2.58-2.50)
Sidechain outliers	100360	5196 (2.58-2.50)
RSRZ outliers	91569	4561 (2.58-2.50)

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	456	 71% 25% 2% 2% 2%
1	B	456	 73% 23% 2% 2% 2%

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	GN1	B	5001	-	-	-	X
6	DCA	A	2000	-	-	-	X
6	DCA	B	2001	-	-	-	X

2 Entry composition [i](#)

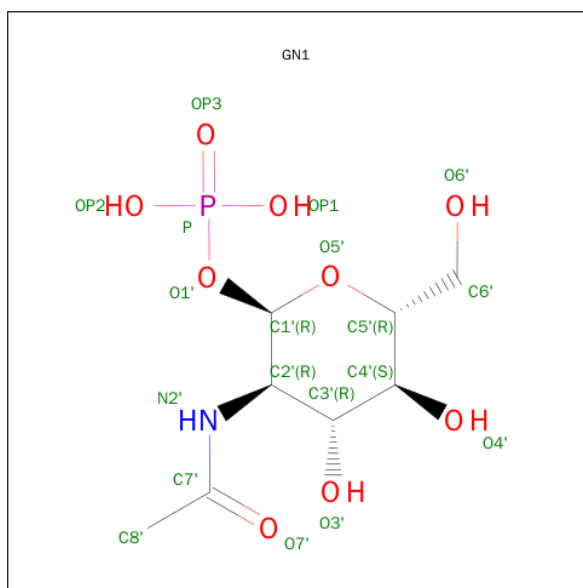
There are 8 unique types of molecules in this entry. The entry contains 7160 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Bifunctional protein glmU.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	449	Total	C	N	O	S	0	0	0
			3372	2103	607	650	12			
1	B	450	Total	C	N	O	S	0	0	0
			3392	2115	611	654	12			

- Molecule 2 is SUGAR (2-(ACETYLAMINO)-2-DEOXY-1-O-PHOSPHONO-ALPHA-D-GLUCOPYRANOSE) (three-letter code: GN1) (formula: $C_8H_{16}NO_9P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			19	8	1	9	1		
2	B	1	Total	C	N	O	P	0	0
			19	8	1	9	1		

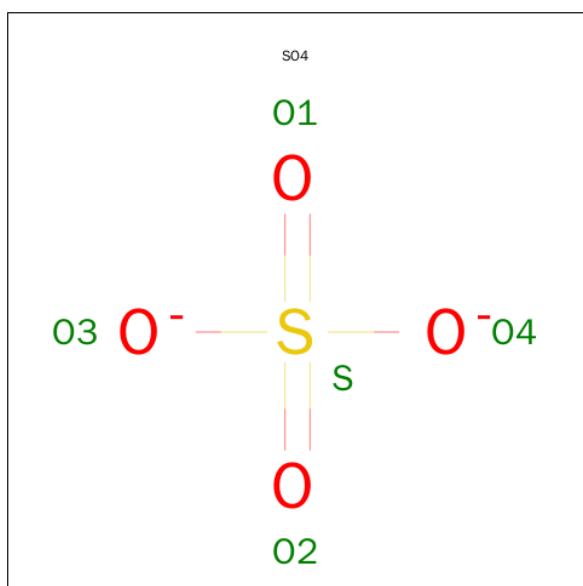
- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Mg	0	0
			1	1		

- Molecule 4 is COBALT (II) ION (three-letter code: CO) (formula: Co).

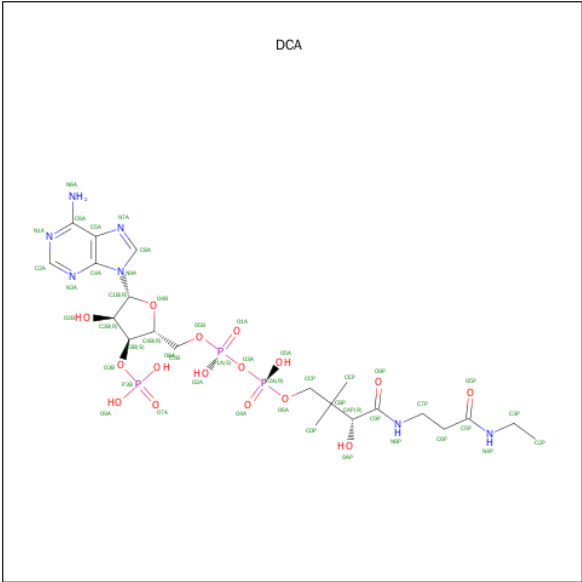
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	2	Total	Co	0	0
			2	2		
4	A	1	Total	Co	0	0
			1	1		

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



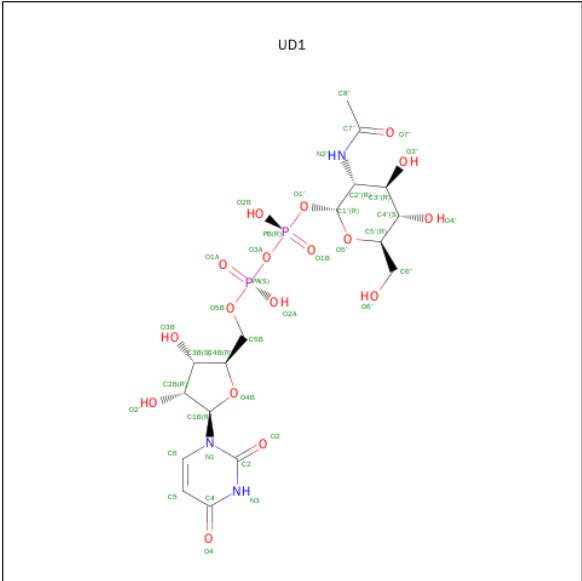
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 6 is DESULFO-COENZYME A (three-letter code: DCA) (formula: C₂₁H₃₆N₇O₁₆P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	N	O	P	0	0
			47	21	7	16	3		
6	B	1	Total	C	N	O	P	0	0
			47	21	7	16	3		

- Molecule 7 is URIDINE-DIPHOSPHATE-N-ACETYLGLUCOSAMINE (three-letter code: UD1) (formula: C₁₇H₂₇N₃O₁₇P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	A	1	Total	C	N	O	P	0	0
			39	17	3	17	2		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	B	1	Total	C	N	O	P	0	0
			39	17	3	17	2		

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	89	Total	O	0	0
			89	89		
8	B	88	Total	O	0	0
			88	88		

4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	102.97Å 102.97Å 644.08Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	27.45 – 2.54 29.88 – 2.54	Depositor EDS
% Data completeness (in resolution range)	94.3 (27.45-2.54) 94.3 (29.88-2.54)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.87 (at 2.54Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.209 , 0.247 0.208 , 0.244	Depositor DCC
R_{free} test set	2057 reflections (4.92%)	DCC
Wilson B-factor (Å ²)	25.0	Xtriage
Anisotropy	0.178	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 29.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	6 of 44213 reflections (0.014%)	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7160	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 8.05% 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, 0.375 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, GN1, CO, DCA, UD1, SO4

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	0.69	0/3420	0.73	1/4640 (0.0%)
1	B	0.70	1/3440 (0.0%)	0.72	2/4664 (0.0%)
All	All	0.70	1/6860 (0.0%)	0.73	3/9304 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	385	CYS	CB-SG	-5.35	1.73	1.81

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	385	CYS	CA-CB-SG	-5.45	104.19	114.00
1	A	182	ARG	NE-CZ-NH1	-5.07	117.77	120.30
1	B	68	ASP	CB-CG-OD1	-5.06	113.75	118.30

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	3372	0	3369	89	0
1	B	3392	0	3401	76	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	19	0	14	0	0
2	B	19	0	14	1	0
3	B	1	0	0	0	0
4	A	1	0	0	0	0
4	B	2	0	0	0	0
5	B	5	0	0	0	0
6	A	47	0	32	1	0
6	B	47	0	32	3	0
7	A	39	0	25	0	0
7	B	39	0	25	1	0
8	A	89	0	0	6	0
8	B	88	0	0	4	0
All	All	7160	0	6912	167	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 12.

All (167) 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:270:ARG:H	1:A:288:HIS:HD2	1.04	1.00
1:B:270:ARG:H	1:B:288:HIS:HD2	1.13	0.94
1:B:243:GLN:HG3	8:B:8065:HOH:O	1.72	0.89
1:B:94:ALA:HB3	1:B:97:GLU:HG2	1.54	0.87
1:B:19:MET:HG3	1:B:25:LYS:HG3	1.61	0.82
1:A:270:ARG:H	1:A:288:HIS:CD2	1.96	0.80
1:A:197:TYR:HB3	1:A:199:THR:HG23	1.63	0.80
1:B:270:ARG:H	1:B:288:HIS:CD2	2.00	0.78
1:A:9:VAL:HG22	1:A:52:HIS:HB2	1.67	0.75
1:A:444:THR:HG21	1:B:444:THR:HG21	1.70	0.74
1:A:94:ALA:HB3	1:A:97:GLU:HG2	1.71	0.70
1:B:386:ASN:O	1:B:392:LYS:HA	1.90	0.70
1:B:182:ARG:HH22	1:B:208:GLU:CD	1.97	0.68
1:B:345:GLY:HA3	1:B:362:GLY:O	1.94	0.68
1:A:76:GLN:HE21	1:A:84:HIS:HD2	1.43	0.66
1:B:203:ALA:O	1:B:207:GLN:HG2	1.95	0.66
1:A:255:ARG:HA	1:A:255:ARG:NE	2.10	0.65
1:A:354:ARG:HD2	8:A:8108:HOH:O	1.97	0.65
1:A:221:SER:HA	1:A:224:GLU:HG3	1.79	0.64
1:B:363:HIS:CD2	1:B:380:ALA:HB2	2.33	0.64
1:A:413:VAL:HG21	1:A:427:VAL:HG12	1.81	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:142:ILE:HD13	1:A:152:ILE:HG12	1.80	0.63
1:A:21:SER:HB2	1:A:229:ARG:HH12	1.63	0.62
1:B:151:GLY:HA2	1:B:202:ILE:CG2	2.30	0.62
1:A:345:GLY:HA3	1:A:362:GLY:O	1.99	0.62
1:B:248:LEU:HD11	1:B:257:PRO:HG3	1.82	0.61
1:B:68:ASP:HB3	8:B:8129:HOH:O	2.01	0.61
1:A:94:ALA:HB3	1:A:97:GLU:CG	2.30	0.60
1:A:156:LYS:HG2	8:A:8162:HOH:O	2.02	0.60
1:B:151:GLY:HA2	1:B:202:ILE:HG22	1.83	0.60
1:A:59:GLY:HA2	1:A:62:LEU:HD12	1.84	0.59
1:B:300:ASN:O	1:B:317:ASP:HA	2.01	0.59
1:B:11:LEU:HD12	1:B:103:TYR:CE2	2.37	0.59
1:B:299:LYS:HD3	1:B:316:GLU:OE2	2.03	0.59
1:B:131:VAL:HG21	1:B:220:LEU:HD13	1.84	0.59
1:B:178:ALA:O	1:B:182:ARG:HG3	2.02	0.59
1:A:449:TRP:HZ3	1:A:451:ARG:HG2	1.68	0.58
1:A:248:LEU:HD11	1:A:257:PRO:HG3	1.85	0.58
1:A:347:PHE:HB2	1:A:364:LEU:HD13	1.86	0.58
1:A:26:VAL:HB	1:A:36:VAL:HB	1.84	0.58
1:A:212:ILE:HD12	1:A:212:ILE:N	2.19	0.58
1:B:82:THR:OG1	1:B:198:ILE:HG22	2.04	0.57
1:B:307:CYS:SG	1:B:324:CYS:HB2	2.44	0.57
1:B:393:PHE:CG	1:B:412:PRO:HD3	2.40	0.57
1:A:151:GLY:HA2	1:A:202:ILE:HG22	1.86	0.57
1:B:385:CYS:HB2	1:B:410:VAL:HA	1.85	0.57
1:B:26:VAL:HB	1:B:36:VAL:HB	1.85	0.57
1:A:270:ARG:N	1:A:288:HIS:HD2	1.88	0.56
1:A:386:ASN:O	1:A:392:LYS:HA	2.06	0.56
1:A:219:ARG:O	1:A:222:GLU:HB2	2.06	0.56
1:B:94:ALA:HB3	1:B:97:GLU:CG	2.33	0.56
1:A:68:ASP:O	1:A:71:LEU:HB2	2.05	0.55
1:A:24:PRO:HD2	1:A:27:LEU:HD12	1.88	0.55
1:A:161:GLU:O	1:A:164:GLN:HB2	2.07	0.55
1:A:246:LYS:HG3	8:A:8170:HOH:O	2.07	0.54
1:B:12:ALA:HB1	1:B:26:VAL:HG21	1.89	0.54
1:A:141:ARG:HG3	1:A:165:ILE:HB	1.89	0.53
1:A:286:LEU:CD2	1:A:303:ILE:HD12	2.38	0.53
1:A:193:GLN:HE21	1:A:193:GLN:HA	1.74	0.53
1:A:363:HIS:N	1:A:363:HIS:CD2	2.76	0.53
1:B:67:LYS:HD2	1:B:67:LYS:O	2.09	0.53
1:B:145:GLU:O	1:B:146:ASN:HB2	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:300:ASN:O	1:A:317:ASP:HA	2.09	0.52
1:A:116:ARG:NH2	1:A:216:HIS:O	2.43	0.52
1:A:351:LYS:O	1:A:368:GLY:HA2	2.10	0.52
1:A:408:GLN:NE2	1:A:426:THR:OG1	2.35	0.52
1:B:132:LYS:O	1:B:133:LEU:HD23	2.10	0.52
1:A:151:GLY:HA2	1:A:202:ILE:CG2	2.40	0.51
1:B:59:GLY:O	1:B:63:LYS:HG3	2.11	0.51
1:A:54:VAL:HA	1:A:74:VAL:O	2.11	0.51
1:A:68:ASP:H	1:A:71:LEU:HD12	1.76	0.51
1:A:176:ASN:HB3	1:A:179:ASP:OD2	2.11	0.50
1:A:385:CYS:HB2	1:A:410:VAL:HA	1.93	0.50
1:B:68:ASP:HB2	1:B:71:LEU:HD12	1.94	0.50
1:A:387:TYR:CE2	1:A:389:GLY:HA2	2.46	0.50
1:B:221:SER:HB3	1:B:238:VAL:HG11	1.93	0.50
1:B:219:ARG:O	1:B:222:GLU:HG2	2.12	0.49
1:A:150:THR:HG22	1:A:206:TYR:CZ	2.47	0.49
1:A:197:TYR:HB3	1:A:199:THR:CG2	2.39	0.49
1:B:112:GLU:HB2	1:B:218:GLN:NE2	2.27	0.49
1:A:133:LEU:HD12	1:A:168:ILE:HG12	1.95	0.49
1:A:19:MET:O	1:A:229:ARG:NH1	2.46	0.48
1:B:335:GLY:HA3	1:B:352:LYS:HA	1.95	0.48
1:B:6:MET:HG2	1:B:118:ARG:CZ	2.43	0.48
1:B:283:ASN:O	1:B:300:ASN:HA	2.13	0.48
2:B:5001:GN1:C7'	6:B:2001:DCA:H21	2.44	0.48
1:B:40:ILE:HD13	1:B:71:LEU:HD11	1.95	0.48
1:B:63:LYS:HA	1:B:73:TRP:CH2	2.49	0.48
1:B:6:MET:HE3	1:B:118:ARG:HD2	1.96	0.48
1:B:121:LYS:HD2	1:B:175:ALA:HA	1.95	0.48
1:A:129:LEU:HD12	1:A:215:VAL:O	2.14	0.48
1:A:328:PRO:O	1:A:346:ASN:HA	2.14	0.47
1:A:277:ASN:O	1:A:295:GLY:HA2	2.13	0.47
1:A:150:THR:HG22	1:A:206:TYR:CE1	2.49	0.47
1:B:365:THR:HG23	1:B:382:THR:O	2.15	0.47
1:A:182:ARG:HH11	1:A:182:ARG:HG3	1.80	0.47
1:A:131:VAL:HG12	1:A:217:PRO:HD2	1.97	0.46
1:B:394:LYS:HG2	1:B:395:THR:N	2.30	0.46
1:A:281:GLU:O	1:A:299:LYS:HA	2.15	0.46
1:B:121:LYS:HG3	1:B:174:ILE:HG23	1.97	0.46
1:B:33:LYS:HD2	1:B:38:HIS:CD2	2.51	0.46
1:B:141:ARG:HG2	1:B:162:GLN:O	2.15	0.46
1:B:69:ASP:OD1	1:B:69:ASP:C	2.55	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:131:VAL:CG2	1:B:220:LEU:HD13	2.46	0.45
1:A:334:PRO:HD3	8:A:8111:HOH:O	2.16	0.45
1:A:76:GLN:HE21	1:A:84:HIS:CD2	2.29	0.45
1:B:319:ASN:HB3	1:B:337:GLU:HG3	1.98	0.45
7:B:4001:UD1:H5'1	8:B:8008:HOH:O	2.15	0.45
1:B:21:SER:HB2	1:B:229:ARG:NH2	2.32	0.45
1:A:108:LEU:HB2	8:A:8062:HOH:O	2.16	0.45
1:A:393:PHE:CD2	1:A:412:PRO:HD3	2.51	0.45
1:B:24:PRO:HD3	1:B:61:LEU:HD12	1.98	0.45
1:B:68:ASP:HB3	1:B:71:LEU:HG	1.99	0.44
1:B:408:GLN:HE21	1:B:408:GLN:HB3	1.46	0.44
1:A:248:LEU:HD11	1:A:257:PRO:CG	2.47	0.44
1:B:69:ASP:C	1:B:71:LEU:H	2.21	0.44
1:A:333:ARG:HB3	1:A:334:PRO:HD2	2.00	0.44
1:B:122:PRO:HG2	1:B:125:GLY:HA3	2.00	0.44
1:B:54:VAL:HA	1:B:74:VAL:O	2.17	0.44
1:B:25:LYS:HE3	8:B:8057:HOH:O	2.17	0.44
1:A:133:LEU:CD1	1:A:168:ILE:HG12	2.48	0.44
1:A:177:GLY:O	1:A:181:LYS:HG3	2.17	0.44
1:B:176:ASN:OD1	1:B:178:ALA:HB3	2.18	0.43
6:A:2000:DCA:H31	8:A:8091:HOH:O	2.17	0.43
1:B:265:THR:HB	1:B:283:ASN:OD1	2.19	0.43
1:B:405:SER:OG	6:B:2001:DCA:H32	2.19	0.43
1:B:176:ASN:HB3	1:B:179:ASP:OD2	2.19	0.43
1:A:253:MET:HE3	1:A:255:ARG:NH2	2.34	0.43
1:A:129:LEU:HD23	1:A:172:ILE:HD12	2.00	0.43
1:A:122:PRO:HG2	1:A:211:GLU:HG3	2.01	0.43
1:A:152:ILE:HB	1:A:199:THR:OG1	2.19	0.42
1:A:141:ARG:HD3	1:A:141:ARG:HA	1.85	0.42
1:B:208:GLU:HB3	1:B:210:ARG:HG2	2.00	0.42
1:B:438:ILE:HG12	1:B:439:SER:N	2.35	0.42
1:A:286:LEU:HD22	1:A:290:VAL:HG11	2.02	0.42
1:B:9:VAL:HG13	1:B:101:MET:HA	2.01	0.42
1:A:363:HIS:CD2	1:A:380:ALA:HB2	2.55	0.42
1:A:283:ASN:O	1:A:300:ASN:HA	2.20	0.42
1:A:255:ARG:HB2	1:A:275:ASP:HA	2.02	0.42
1:B:112:GLU:HB2	1:B:218:GLN:HE22	1.84	0.42
1:A:333:ARG:HB3	1:A:334:PRO:CD	2.50	0.42
1:A:76:GLN:NE2	1:A:85:ALA:HB2	2.35	0.41
1:A:180:MET:HE3	1:A:184:LEU:HG	2.01	0.41
1:A:299:LYS:O	1:A:300:ASN:C	2.59	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:333:ARG:HB3	1:B:334:PRO:HD2	2.01	0.41
1:B:299:LYS:O	1:B:300:ASN:C	2.59	0.41
1:A:141:ARG:NH1	1:A:163:ARG:O	2.52	0.41
1:A:346:ASN:C	1:A:346:ASN:HD22	2.22	0.41
1:A:429:ARG:HB3	1:A:429:ARG:CZ	2.51	0.41
1:A:286:LEU:HD21	1:A:303:ILE:HD12	2.01	0.41
1:B:30:LEU:O	1:B:38:HIS:HE1	2.03	0.41
1:A:90:ALA:HB3	1:A:91:PRO:HD3	2.01	0.41
1:A:136:PRO:HG3	1:A:166:GLN:HA	2.02	0.41
1:A:363:HIS:O	1:A:380:ALA:HA	2.21	0.41
1:A:283:ASN:H	1:A:300:ASN:ND2	2.18	0.41
1:B:405:SER:HG	6:B:2001:DCA:H32	1.85	0.41
1:A:121:LYS:HA	1:A:122:PRO:HD3	1.91	0.41
1:A:94:ALA:O	1:A:97:GLU:HG2	2.21	0.41
1:B:362:GLY:C	1:B:363:HIS:CG	2.95	0.41
1:A:393:PHE:CG	1:A:412:PRO:HD3	2.56	0.41
1:A:238:VAL:O	1:A:242:GLU:HG3	2.21	0.41
1:B:90:ALA:HB3	1:B:91:PRO:HD3	2.01	0.41
1:A:323:ALA:O	1:A:341:GLY:HA2	2.21	0.41
1:B:73:TRP:CD1	1:B:73:TRP:N	2.89	0.40
1:B:321:ALA:HB3	1:B:340:GLU:N	2.36	0.40
1:B:121:LYS:HE3	1:B:122:PRO:O	2.20	0.40
1:B:4:ASN:HD22	1:B:4:ASN:C	2.24	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	447/456 (98%)	425 (95%)	21 (5%)	1 (0%)	52	73
1	B	448/456 (98%)	428 (96%)	18 (4%)	2 (0%)	39	60

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	895/912 (98%)	853 (95%)	39 (4%)	3 (0%)	46 66

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	70	ASN
1	A	192	ALA
1	B	56	GLY

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	349/362 (96%)	331 (95%)	18 (5%)	29 49
1	B	353/362 (98%)	336 (95%)	17 (5%)	31 53
All	All	702/724 (97%)	667 (95%)	35 (5%)	30 51

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	53	LEU
1	A	57	HIS
1	A	123	GLN
1	A	141	ARG
1	A	164	GLN
1	A	182	ARG
1	A	193	GLN
1	A	199	THR
1	A	219	ARG
1	A	229	ARG
1	A	234	ARG
1	A	255	ARG
1	A	331	ARG
1	A	346	ASN
1	A	363	HIS

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Mol	Chain	Res	Type
1	A	405	SER
1	A	408	GLN
1	A	447	GLU
1	B	4	ASN
1	B	6	MET
1	B	50	HIS
1	B	67	LYS
1	B	69	ASP
1	B	123	GLN
1	B	132	LYS
1	B	219	ARG
1	B	241	SER
1	B	255	ARG
1	B	331	ARG
1	B	346	ASN
1	B	363	HIS
1	B	391	ASN
1	B	408	GLN
1	B	446	LYS
1	B	449	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	79	GLN
1	A	84	HIS
1	A	193	GLN
1	A	288	HIS
1	A	300	ASN
1	A	363	HIS
1	B	4	ASN
1	B	28	HIS
1	B	38	HIS
1	B	72	ASN
1	B	166	GLN
1	B	193	GLN
1	B	218	GLN
1	B	243	GLN
1	B	288	HIS
1	B	363	HIS
1	B	377	ASN
1	B	408	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 11 ligands modelled in this entry, 4 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$
6	DCA	A	2000	-	39,49,49	2.32	13 (33%)	50,74,74	1.70	6 (12%)
7	UD1	A	4000	-	32,41,41	1.73	1 (3%)	46,62,62	1.94	1 (2%)
2	GN1	A	5000	-	18,19,19	2.18	7 (38%)	24,28,28	0.91	0
6	DCA	B	2001	-	39,49,49	2.40	17 (43%)	50,74,74	1.64	6 (12%)
7	UD1	B	4001	3	32,41,41	2.00	3 (9%)	46,62,62	2.01	2 (4%)
2	GN1	B	5001	-	18,19,19	2.13	7 (38%)	24,28,28	0.96	1 (4%)
5	SO4	B	7000	-	4,4,4	1.04	0	6,6,6	0.35	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
6	DCA	A	2000	-	-	0/43/63/63	0/3/3/3
7	UD1	A	4000	-	-	0/22/63/63	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GN1	A	5000	-	-	0/10/31/31	0/1/1/1
6	DCA	B	2001	-	-	0/43/63/63	0/3/3/3
7	UD1	B	4001	3	-	0/22/63/63	0/3/3/3
2	GN1	B	5001	-	-	0/10/31/31	0/1/1/1
5	SO4	B	7000	-	-	0/0/0/0	0/0/0/0

All (48) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	B	4001	UD1	C6-C5	-2.14	1.33	1.38
2	B	5001	GN1	P-OP2	-2.11	1.47	1.54
7	B	4001	UD1	PB-O2B	-2.05	1.46	1.54
6	A	2000	DCA	C3P-N4P	2.08	1.54	1.46
6	A	2000	DCA	P2A-O4A	2.10	1.58	1.51
6	B	2001	DCA	O4B-C1B	2.12	1.43	1.41
6	B	2001	DCA	P1A-O5B	2.12	1.68	1.59
6	A	2000	DCA	P1A-O5B	2.20	1.69	1.59
6	A	2000	DCA	P3B-O7A	2.33	1.58	1.51
6	B	2001	DCA	C5A-N7A	2.33	1.47	1.39
2	A	5000	GN1	C1'-C2'	2.39	1.57	1.53
6	A	2000	DCA	O5P-C5P	2.39	1.28	1.23
6	A	2000	DCA	P2A-O6A	2.43	1.70	1.59
6	B	2001	DCA	O6A-CCP	2.44	1.52	1.43
6	B	2001	DCA	C2B-C3B	2.47	1.58	1.53
6	B	2001	DCA	P2A-O4A	2.48	1.60	1.51
6	B	2001	DCA	P3B-O3B	2.52	1.67	1.60
2	B	5001	GN1	C1'-C2'	2.56	1.57	1.53
2	B	5001	GN1	O5'-C5'	2.65	1.51	1.44
6	B	2001	DCA	C3P-N4P	2.65	1.56	1.46
6	B	2001	DCA	P1A-O1A	2.67	1.60	1.51
2	A	5000	GN1	O5'-C5'	2.73	1.51	1.44
6	B	2001	DCA	P2A-O6A	2.81	1.71	1.59
2	B	5001	GN1	O4'-C4'	2.87	1.49	1.43
2	A	5000	GN1	O5'-C1'	2.88	1.49	1.41
6	B	2001	DCA	CEP-CBP	2.90	1.60	1.53
2	A	5000	GN1	C3'-C2'	2.93	1.59	1.53
2	A	5000	GN1	C2'-N2'	2.94	1.50	1.45
6	B	2001	DCA	C8A-N7A	3.01	1.40	1.34
6	A	2000	DCA	O2B-C2B	3.05	1.50	1.43
6	A	2000	DCA	O6A-CCP	3.06	1.54	1.43
6	A	2000	DCA	C5A-N7A	3.08	1.49	1.39
2	B	5001	GN1	O5'-C1'	3.10	1.49	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	5000	GN1	O3'-C3'	3.24	1.50	1.43
6	A	2000	DCA	C8A-N7A	3.41	1.41	1.34
6	B	2001	DCA	P3B-O7A	3.50	1.62	1.51
6	B	2001	DCA	O2B-C2B	3.57	1.51	1.43
2	B	5001	GN1	C4'-C5'	3.97	1.61	1.53
2	B	5001	GN1	C2'-N2'	4.05	1.52	1.45
6	A	2000	DCA	C2A-N1A	4.30	1.42	1.33
6	B	2001	DCA	C6P-C5P	4.51	1.60	1.51
6	B	2001	DCA	C2A-N1A	4.75	1.42	1.33
6	A	2000	DCA	C6P-C5P	4.77	1.60	1.51
2	A	5000	GN1	C4'-C5'	5.03	1.63	1.53
6	B	2001	DCA	C4A-N3A	6.58	1.45	1.35
6	A	2000	DCA	C4A-N3A	7.27	1.46	1.35
7	A	4000	UD1	C6-N1	8.39	1.47	1.35
7	B	4001	UD1	C6-N1	9.87	1.49	1.35

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	2000	DCA	C1B-N9A-C4A	-8.17	114.61	126.94
6	B	2001	DCA	C1B-N9A-C4A	-7.86	115.08	126.94
6	A	2000	DCA	N3A-C2A-N1A	-3.25	126.41	128.89
6	A	2000	DCA	C6P-C5P-N4P	-2.74	111.69	116.46
6	B	2001	DCA	O3A-P2A-O6A	-2.08	97.41	102.94
2	B	5001	GN1	O6'-C6'-C5'	-2.02	104.66	111.33
6	A	2000	DCA	C3B-C2B-C1B	2.03	104.85	99.98
6	A	2000	DCA	O6A-CCP-CBP	2.10	113.93	110.55
7	B	4001	UD1	C4B-O4B-C1B	2.13	112.06	109.72
6	B	2001	DCA	O4B-C1B-N9A	2.35	113.01	108.10
6	B	2001	DCA	C3P-N4P-C5P	2.40	126.67	122.31
6	B	2001	DCA	O6A-CCP-CBP	2.45	114.49	110.55
6	A	2000	DCA	CEP-CBP-CCP	2.94	112.32	108.50
6	B	2001	DCA	CEP-CBP-CCP	3.97	113.65	108.50
7	A	4000	UD1	C4-N3-C2	11.67	125.70	114.14
7	B	4001	UD1	C4-N3-C2	12.25	126.27	114.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	2000	DCA	1	0
6	B	2001	DCA	3	0
7	B	4001	UD1	1	0
2	B	5001	GN1	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	449/456 (98%)	-0.30	0 100 100	12, 21, 31, 37	0
1	B	450/456 (98%)	-0.23	9 (2%) 68 73	13, 22, 31, 42	0
All	All	899/912 (98%)	-0.26	9 (1%) 84 87	12, 22, 31, 42	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	69	ASP	3.9
1	B	4	ASN	2.6
1	B	70	ASN	2.5
1	B	146	ASN	2.5
1	B	64	GLN	2.3
1	B	164	GLN	2.2
1	B	161	GLU	2.2
1	B	67	LYS	2.1
1	B	3	ASN	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 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
6	DCA	A	2000	47/47	0.69	0.45	7.27	29,40,57,58	0
6	DCA	B	2001	47/47	0.65	0.43	6.35	33,41,57,60	0
2	GN1	B	5001	19/19	0.93	0.28	3.84	27,35,40,41	0
2	GN1	A	5000	19/19	0.93	0.20	1.11	34,36,38,39	0
7	UD1	B	4001	39/39	0.98	0.19	0.97	14,19,23,25	0
7	UD1	A	4000	39/39	0.97	0.14	-0.57	17,23,31,31	0
5	SO4	B	7000	5/5	0.95	0.11	-0.61	43,45,46,47	0
4	CO	B	6001	1/1	0.94	0.16	-	23,23,23,23	1
4	CO	A	6003	1/1	0.98	0.19	-	23,23,23,23	1
4	CO	B	6002	1/1	0.96	0.12	-	23,23,23,23	1
3	MG	B	6000	1/1	0.91	0.17	-	26,26,26,26	0

6.5 Other polymers [i](#)

There are no such residues in this entry.