



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 07:36 PM GMT

PDB ID : 3C16
Title : Complex of GS-Alpha with the Catalytic Domains of Mammalian Adenylyl Cyclase: Complex with Adenosine-5'-Triphosphate and Ca
Authors : Mou, T.-C.; Sprang, S.R.
Deposited on : 2008-01-22
Resolution : 2.87 Å (reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

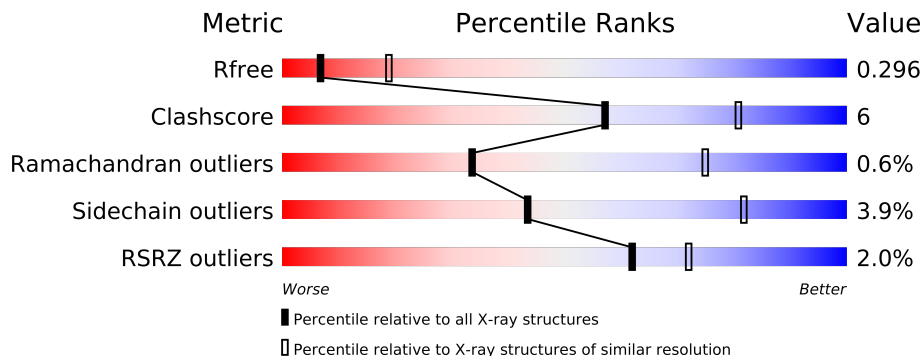
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.87 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}	66092	1360 (2.90-2.86)
Clashscore	79885	1696 (2.90-2.86)
Ramachandran outliers	78287	1647 (2.90-2.86)
Sidechain outliers	78261	1650 (2.90-2.86)
RSRZ outliers	66119	1362 (2.90-2.86)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	225	
2	B	212	
3	C	402	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
5	MG	C	398	-	X

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 5739 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Adenylate cyclase type 5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	190	Total	C	N	O	S	0	0	0
			1484	933	260	274	17			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	356	MET	-	EXPRESSION TAG	UNP P30803
A	357	HIS	-	EXPRESSION TAG	UNP P30803
A	358	HIS	-	EXPRESSION TAG	UNP P30803
A	359	HIS	-	EXPRESSION TAG	UNP P30803
A	360	HIS	-	EXPRESSION TAG	UNP P30803
A	361	HIS	-	EXPRESSION TAG	UNP P30803
A	362	HIS	-	EXPRESSION TAG	UNP P30803
A	476	MET	VAL	ENGINEERED	UNP P30803

- Molecule 2 is a protein called Adenylate cyclase type 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	186	Total	C	N	O	S	0	0	0
			1443	921	236	276	10			

- Molecule 3 is a protein called Guanine nucleotide-binding protein G(s) subunit alpha isoforms short.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	331	Total	C	N	O	S	0	0	0
			2710	1718	472	507	13			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	373	GLY	-	EXPRESSION TAG	UNP P04896

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	374	GLY	-	EXPRESSION TAG	UNP P04896
C	375	HIS	-	EXPRESSION TAG	UNP P04896
C	376	HIS	-	EXPRESSION TAG	UNP P04896
C	377	HIS	-	EXPRESSION TAG	UNP P04896
C	378	HIS	-	EXPRESSION TAG	UNP P04896
C	379	HIS	-	EXPRESSION TAG	UNP P04896
C	380	HIS	-	EXPRESSION TAG	UNP P04896

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	C	1	Total Cl 1 1	0	0

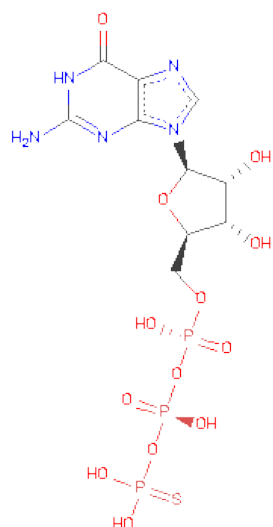
- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	C	1	Total Mg 1 1	0	0

- Molecule 6 is CALCIUM ION (three-letter code: CA) (formula: Ca).

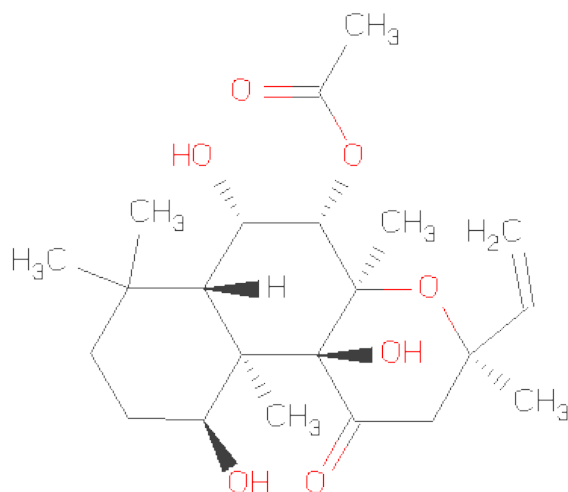
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	2	Total Ca 2 2	0	0

- Molecule 7 is 5'-GUANOSINE-DIPHOSPHATE-MONOTHIOPHOSPHATE (three-letter code: GSP) (formula: C₁₀H₁₆N₅O₁₃P₃S).



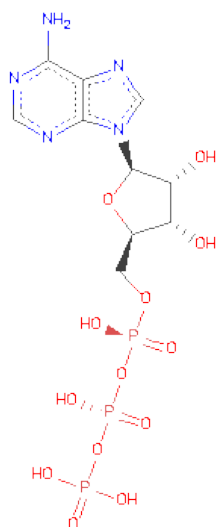
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	C	1	Total	C	N	O	P	S	
			32	10	5	13	3	1	
									0
									0

- Molecule 8 is FORSKOLIN (three-letter code: FOK) (formula: $C_{22}H_{34}O_7$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	O		
			29	22	7	0	0

- Molecule 9 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
9	A	1	31	10	5	13	3	0	0

- Molecule 10 is water.

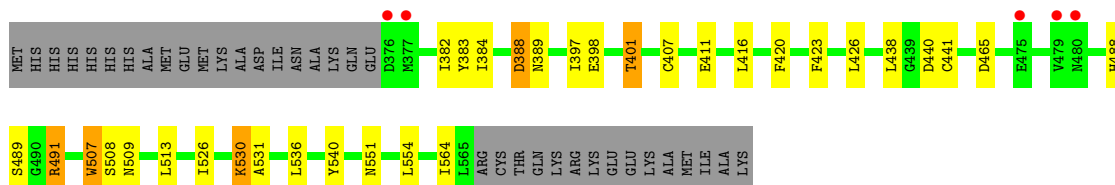
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	6	Total	O	0	0
			6	6		

3 Residue-property plots

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.

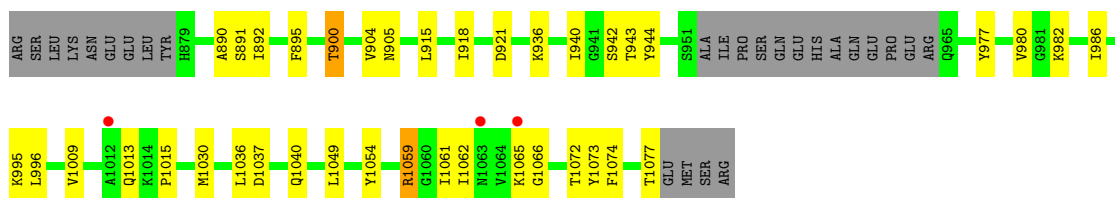
- Molecule 1: Adenylate cyclase type 5

Chain A: 



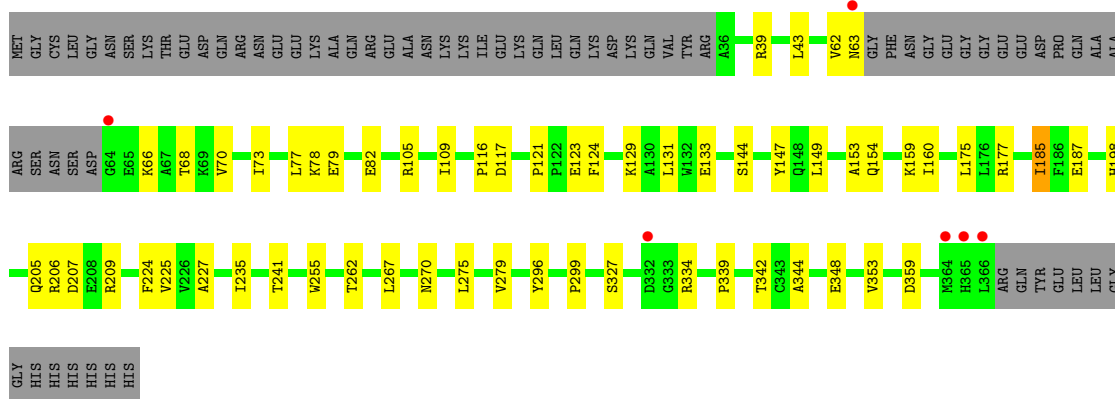
- Molecule 2: Adenylate cyclase type 2

Chain B: 



- Molecule 3: Guanine nucleotide-binding protein G(s) subunit alpha isoforms short

Chain C: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	118.33Å 133.53Å 70.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.87 24.64 – 2.87	Depositor EDS
% Data completeness (in resolution range)	85.9 (15.00-2.87) 86.0 (24.64-2.87)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.66 (at 2.89Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.249 , 0.299 0.246 , 0.296	Depositor DCC
R_{free} test set	1156 reflections (5.41%)	DCC
Wilson B-factor (Å ²)	43.0	Xtriage
Anisotropy	0.129	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 3.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Outliers	1 of 22537 reflections (0.004%)	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	5739	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.55% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GSP, MG, CL, CA, ATP, FOK

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	0/1512	0.49	0/2038
2	B	0.35	0/1468	0.50	0/1982
3	C	0.35	0/2767	0.50	0/3744
All	All	0.35	0/5747	0.50	0/7764

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1484	0	1454	18	0
2	B	1443	0	1441	23	0
3	C	2710	0	2657	31	0
4	C	1	0	0	1	0
5	C	1	0	0	0	0
6	A	2	0	0	0	0
7	C	32	0	12	2	0
8	A	29	0	34	3	0
9	A	31	0	12	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
10	A	6	0	0	0	0
All	All	5739	0	5610	73	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 6.

All (73) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:C:62:VAL:HG12	3:C:63:ASN:H	1.46	0.79
2:B:890:ALA:HB1	2:B:996:LEU:HD11	1.65	0.78
8:A:101:FOK:H201	8:A:101:FOK:H173	1.69	0.74
3:C:79:GLU:HA	3:C:82:GLU:HG2	1.70	0.73
8:A:101:FOK:H202	8:A:101:FOK:H193	1.73	0.70
3:C:62:VAL:HG12	3:C:63:ASN:N	2.10	0.67
2:B:891:SER:HA	2:B:943:THR:HA	1.80	0.64
2:B:1059:ARG:HH11	2:B:1062:ILE:HD11	1.66	0.61
3:C:129:LYS:O	3:C:133:GLU:HG2	2.01	0.60
1:A:397:ILE:HD12	1:A:440:ASP:HB2	1.84	0.59
2:B:918:ILE:HG12	2:B:986:ILE:HG13	1.85	0.58
1:A:426:LEU:HD22	1:A:465:ASP:HB3	1.85	0.58
2:B:1062:ILE:HD12	2:B:1072:THR:HG21	1.87	0.57
3:C:144:SER:HA	3:C:147:TYR:CE2	2.39	0.57
3:C:121:PRO:HB2	3:C:123:GLU:OE2	2.05	0.57
3:C:66:LYS:HE3	3:C:175:LEU:HD11	1.88	0.56
2:B:900:THR:HB	2:B:905:ASN:ND2	2.22	0.55
1:A:407:CYS:HB2	1:A:411:GLU:HB3	1.91	0.53
2:B:977:TYR:O	2:B:980:VAL:HB	2.09	0.52
2:B:900:THR:HB	2:B:905:ASN:HD21	1.73	0.52
1:A:526:ILE:HB	1:A:564:ILE:HD12	1.91	0.51
2:B:904:VAL:HG11	3:C:185:ILE:HB	1.93	0.51
3:C:187:GLU:OE2	3:C:198:HIS:ND1	2.44	0.51
3:C:344:ALA:HB3	7:C:395:GSP:N7	2.26	0.51
3:C:255:TRP:CE2	3:C:327:SER:HA	2.46	0.51
3:C:62:VAL:CG1	3:C:63:ASN:H	2.19	0.50
3:C:79:GLU:HA	3:C:82:GLU:CG	2.41	0.50
2:B:921:ASP:HB3	2:B:982:LYS:HD2	1.93	0.50
3:C:77:LEU:HD21	3:C:160:ILE:HG12	1.93	0.50
3:C:270:ASN:HA	3:C:342:THR:O	2.11	0.50
1:A:530:LYS:HD2	1:A:531:ALA:N	2.27	0.49
3:C:78:LYS:HA	3:C:124:PHE:HZ	1.77	0.49
3:C:70:VAL:HA	3:C:73:ILE:HD12	1.94	0.49
1:A:388:ASP:OD1	1:A:491:ARG:HB3	2.13	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:401:THR:HG21	2:B:1066:GLY:HA3	1.93	0.48
3:C:177:ARG:HA	7:C:395:GSP:O2'	2.12	0.48
2:B:891:SER:HB3	2:B:943:THR:HG23	1.95	0.48
2:B:995:LYS:HB2	2:B:1036:LEU:HD22	1.96	0.48
3:C:105:ARG:O	3:C:109:ILE:HG12	2.13	0.47
2:B:940:ILE:HB	2:B:943:THR:HB	1.96	0.47
8:A:101:FOK:H201	8:A:101:FOK:C17	2.42	0.46
2:B:892:ILE:HD12	2:B:942:SER:HB2	1.98	0.46
1:A:388:ASP:HB3	1:A:389:ASN:ND2	2.31	0.46
3:C:235:ILE:HG12	3:C:241:THR:O	2.15	0.45
2:B:1049:LEU:HB3	2:B:1054:TYR:HB2	1.99	0.45
3:C:144:SER:HB2	3:C:149:LEU:HD23	1.97	0.45
3:C:62:VAL:CG1	3:C:63:ASN:N	2.79	0.45
1:A:438:LEU:HB2	1:A:441:CYS:HB3	1.98	0.45
3:C:206:ARG:HG2	3:C:209:ARG:NH1	2.32	0.45
1:A:488:HIS:CD2	1:A:489:SER:O	2.70	0.45
3:C:275:LEU:O	3:C:279:VAL:HG23	2.17	0.44
1:A:488:HIS:CG	1:A:489:SER:N	2.86	0.43
3:C:43:LEU:O	3:C:224:PHE:HA	2.17	0.43
3:C:109:ILE:HG13	3:C:131:LEU:CD1	2.48	0.43
3:C:227:ALA:HB3	4:C:397:CL:CL	2.56	0.43
1:A:536:LEU:HD11	1:A:540:TYR:CD1	2.54	0.42
3:C:205:GLN:O	3:C:209:ARG:HG3	2.20	0.42
2:B:1059:ARG:HG3	2:B:1074:PHE:CZ	2.54	0.41
3:C:342:THR:HG22	3:C:353:VAL:HG21	2.03	0.41
1:A:383:TYR:C	1:A:384:ILE:HG13	2.39	0.41
2:B:995:LYS:HD2	2:B:1037:ASP:OD2	2.20	0.41
1:A:420:PHE:HA	1:A:423:PHE:HB2	2.03	0.41
2:B:936:LYS:HD2	2:B:944:TYR:OH	2.20	0.41
2:B:1009:VAL:HG11	2:B:1015:PRO:HB3	2.01	0.41
3:C:296:TYR:OH	3:C:299:PRO:HD3	2.21	0.41
1:A:509:ASN:O	1:A:513:LEU:N	2.50	0.41
2:B:895:PHE:CZ	2:B:915:LEU:HB2	2.56	0.41
2:B:1040:GLN:HA	2:B:1073:TYR:O	2.21	0.41
1:A:507:TRP:O	1:A:508:SER:HB3	2.20	0.40
1:A:423:PHE:HA	1:A:426:LEU:HD12	2.03	0.40
2:B:1030:MET:O	2:B:1040:GLN:NE2	2.50	0.40
3:C:267:LEU:HB3	3:C:339:PRO:HA	2.04	0.40
1:A:551:ASN:HB3	1:A:554:LEU:HD12	2.04	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	188/225 (84%)	171 (91%)	17 (9%)	0	100	100
2	B	182/212 (86%)	174 (96%)	6 (3%)	2 (1%)	21	58
3	C	327/402 (81%)	306 (94%)	19 (6%)	2 (1%)	33	75
All	All	697/839 (83%)	651 (93%)	42 (6%)	4 (1%)	33	75

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	116	PRO
2	B	1065	LYS
3	C	153	ALA
2	B	1013	GLN

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	159/189 (84%)	151 (95%)	8 (5%)	34	72
2	B	160/184 (87%)	156 (98%)	4 (2%)	60	90
3	C	298/357 (84%)	286 (96%)	12 (4%)	42	80
All	All	617/730 (84%)	593 (96%)	24 (4%)	43	82

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	382	ILE
1	A	388	ASP

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	398	GLU
1	A	401	THR
1	A	416	LEU
1	A	491	ARG
1	A	507	TRP
1	A	530	LYS
2	B	900	THR
2	B	1059	ARG
2	B	1061	ILE
2	B	1077	THR
3	C	39	ARG
3	C	68	THR
3	C	117	ASP
3	C	154	GLN
3	C	159	LYS
3	C	185	ILE
3	C	207	ASP
3	C	225	VAL
3	C	262	THR
3	C	334	ARG
3	C	348	GLU
3	C	359	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	385	GLN
1	A	537	ASN
2	B	880	GLN
2	B	1001	ASN
2	B	1050	GLN
2	B	1076	ASN
3	C	75	ASN
3	C	245	GLN
3	C	349	ASN
3	C	355	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 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	$\# Z > 2$
9	ATP	A	1	6	33,33,33	1.60	7 (21%)	52,52,52	1.84	9 (17%)
8	FOK	A	101	-	31,31,31	1.01	1 (3%)	54,54,54	1.50	8 (14%)
7	GSP	C	395	5	34,34,34	2.06	3 (8%)	51,54,54	2.65	12 (23%)

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
9	ATP	A	1	6	-	0/22/38/38	0/1/3/3
8	FOK	A	101	-	-	0/7/80/80	0/0/3/3
7	GSP	C	395	5	-	0/21/38/38	0/1/3/3

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	C	395	GSP	PG-S1G	-10.12	1.72	1.90
8	A	101	FOK	O4-C21	4.52	1.45	1.35
7	C	395	GSP	C2-N3	4.14	1.38	1.33
9	A	1	ATP	PB-O3A	3.82	1.66	1.59

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	A	1	ATP	PG-O1G	3.61	1.63	1.51
9	A	1	ATP	O4'-C1'	3.02	1.46	1.41
9	A	1	ATP	PA-O3A	2.84	1.65	1.59
7	C	395	GSP	PG-O3B	2.46	1.63	1.60
9	A	1	ATP	C4-N9	-2.29	1.34	1.37
9	A	1	ATP	PG-O2G	2.26	1.63	1.54
9	A	1	ATP	PB-O3B	2.10	1.63	1.59

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	C	395	GSP	C6-C5-N7	-12.70	132.43	134.14
9	A	1	ATP	N3-C2-N1	-7.99	122.03	128.71
7	C	395	GSP	O3G-PG-S1G	-6.56	105.81	112.73
7	C	395	GSP	O3B-PG-S1G	-5.34	112.16	114.53
8	A	101	FOK	C13-O1-C8	4.73	124.85	120.08
9	A	1	ATP	N3-C4-N9	4.66	133.85	125.43
8	A	101	FOK	C7-O4-C21	-4.25	111.49	117.80
7	C	395	GSP	C5-C4-N3	-3.97	120.19	125.94
7	C	395	GSP	PA-O3A-PB	-3.94	120.14	131.68
7	C	395	GSP	N3-C4-N9	3.79	132.46	126.91
7	C	395	GSP	C2-N3-C4	3.70	120.29	115.09
8	A	101	FOK	O4-C21-C22	3.47	117.65	111.12
9	A	1	ATP	C5-C4-N3	-3.37	118.36	125.70
7	C	395	GSP	PB-O3B-PG	-3.25	120.71	131.81
9	A	1	ATP	PB-O3B-PG	-3.08	122.64	131.68
8	A	101	FOK	O4-C7-C6	2.97	112.84	107.59
7	C	395	GSP	C3'-C2'-C1'	2.97	105.56	100.91
9	A	1	ATP	C4'-O4'-C1'	-2.96	106.53	109.75
9	A	1	ATP	C4-C5-N7	-2.81	107.12	109.52
9	A	1	ATP	C3'-C2'-C1'	2.74	105.20	100.91
9	A	1	ATP	PA-O3A-PB	-2.63	123.96	131.68
9	A	1	ATP	C2-N3-C4	2.46	121.02	114.01
8	A	101	FOK	C3-C4-C5	2.37	111.61	107.88
7	C	395	GSP	O2G-PG-S1G	-2.29	110.31	112.73
7	C	395	GSP	O4'-C1'-N9	2.21	110.50	108.44
8	A	101	FOK	O1-C8-C7	2.13	106.59	102.43
7	C	395	GSP	N7-C8-N9	-2.08	108.48	114.36
8	A	101	FOK	O7-C11-C12	-2.05	117.20	121.90
8	A	101	FOK	C10-C5-C6	2.04	115.71	111.45

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	190/225 (84%)	0.09	5 (2%)	53 63	18, 51, 82, 86	0
2	B	186/212 (87%)	-0.17	3 (1%)	68 78	16, 31, 71, 85	0
3	C	331/402 (82%)	-0.19	6 (1%)	65 75	17, 34, 66, 86	0
All	All	707/839 (84%)	-0.11	14 (1%)	62 72	16, 37, 73, 86	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	376	ASP	4.4
1	A	480	ASN	4.0
2	B	1012	ALA	4.0
1	A	479	VAL	3.9
1	A	475	GLU	3.6
2	B	1063	ASN	3.1
3	C	364	MET	2.9
3	C	332	ASP	2.7
1	A	377	MET	2.7
3	C	365	HIS	2.7
3	C	64	GLY	2.4
3	C	366	LEU	2.3
2	B	1065	LYS	2.3
3	C	63	ASN	2.1

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
5	MG	C	398	1/1	0.19	7.03	17,17,17,17	0
4	CL	C	397	1/1	0.21	1.38	30,30,30,30	0
8	FOK	A	101	29/29	0.19	0.95	16,33,49,50	0
7	GSP	C	395	32/32	0.16	0.19	17,22,32,41	0
9	ATP	A	1	31/31	0.17	-0.03	31,51,64,67	0
6	CA	A	29	1/1	0.05	-8.39	49,49,49,49	0
6	CA	A	30	1/1	0.05	-197.00	38,38,38,38	0

6.5 Other polymers ⓘ

There are no such residues in this entry.