



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:31 PM GMT

PDB ID : 1G16
Title : CRYSTAL STRUCTURE OF SEC4-GDP
Authors : Stroupe, C.; Brunger, A.T.
Deposited on : 2000-10-10
Resolution : 1.80 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

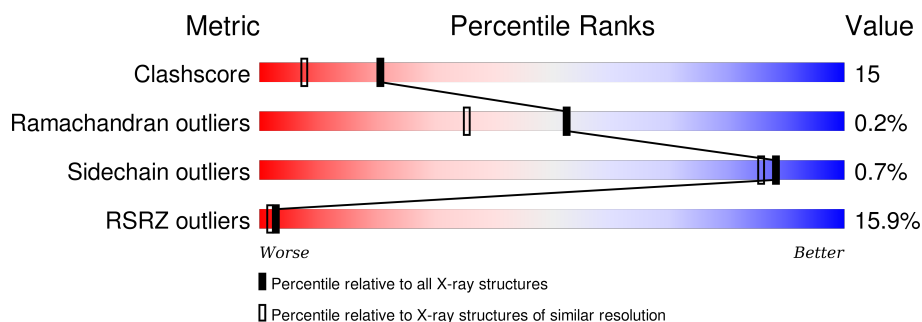
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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))
Clashscore	102246	5383 (1.80-1.80)
Ramachandran outliers	100387	5320 (1.80-1.80)
Sidechain outliers	100360	5319 (1.80-1.80)
RSRZ outliers	91569	4547 (1.80-1.80)

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	170	
1	B	170	
1	C	170	
1	D	170	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5593 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 RAS-RELATED PROTEIN SEC4.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	159	Total	C	N	O	S	Se	0	0	0
			1266	810	211	241	1	3			
1	B	161	Total	C	N	O	S	Se	0	0	0
			1280	817	213	246	1	3			
1	C	167	Total	C	N	O	S	Se	0	0	0
			1325	848	219	254	1	3			
1	D	160	Total	C	N	O	S	Se	0	0	0
			1274	814	212	244	1	3			

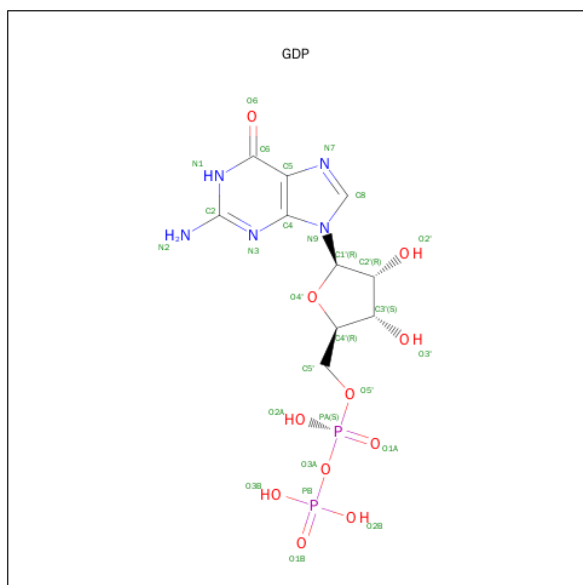
There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	21	MSE	MET	MODIFIED RESIDUE	UNP P07560
A	73	ILE	LEU	CONFLICT	UNP P07560
A	94	MSE	MET	MODIFIED RESIDUE	UNP P07560
A	102	ILE	VAL	CONFLICT	UNP P07560
A	137	MSE	MET	MODIFIED RESIDUE	UNP P07560
B	21	MSE	MET	MODIFIED RESIDUE	UNP P07560
B	73	ILE	LEU	CONFLICT	UNP P07560
B	94	MSE	MET	MODIFIED RESIDUE	UNP P07560
B	102	ILE	VAL	CONFLICT	UNP P07560
B	137	MSE	MET	MODIFIED RESIDUE	UNP P07560
C	21	MSE	MET	MODIFIED RESIDUE	UNP P07560
C	73	ILE	LEU	CONFLICT	UNP P07560
C	94	MSE	MET	MODIFIED RESIDUE	UNP P07560
C	102	ILE	VAL	CONFLICT	UNP P07560
C	137	MSE	MET	MODIFIED RESIDUE	UNP P07560
D	21	MSE	MET	MODIFIED RESIDUE	UNP P07560
D	73	ILE	LEU	CONFLICT	UNP P07560
D	94	MSE	MET	MODIFIED RESIDUE	UNP P07560
D	102	ILE	VAL	CONFLICT	UNP P07560
D	137	MSE	MET	MODIFIED RESIDUE	UNP P07560

- Molecule 2 is COBALT (II) ION (three-letter code: CO) (formula: Co).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Co	0	0
			1	1		
2	A	2	Total	Co	0	0
			2	2		
2	D	3	Total	Co	0	0
			3	3		
2	C	2	Total	Co	0	0
			2	2		

- Molecule 3 is GUANOSINE-5'-DIPHOSPHATE (three-letter code: GDP) (formula: C₁₀H₁₅N₅O₁₁P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			28	10	5	11	2		
3	B	1	Total	C	N	O	P	0	0
			28	10	5	11	2		
3	C	1	Total	C	N	O	P	0	0
			28	10	5	11	2		
3	D	1	Total	C	N	O	P	0	0
			28	10	5	11	2		

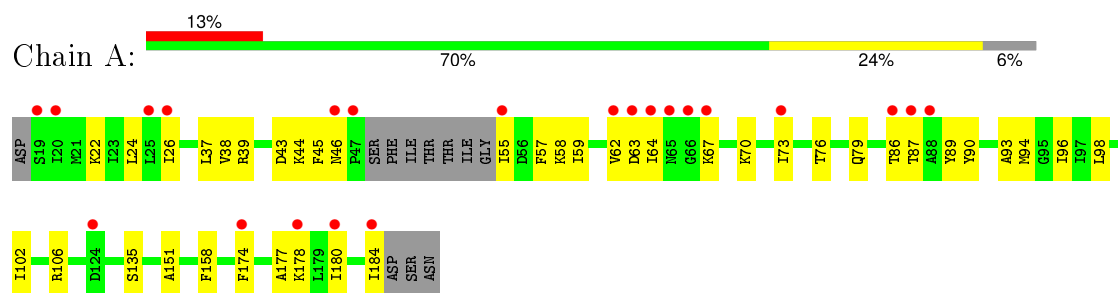
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	105	Total 105	O 105	0	0
4	B	86	Total 86	O 86	0	0
4	C	70	Total 70	O 70	0	0
4	D	67	Total 67	O 67	0	0

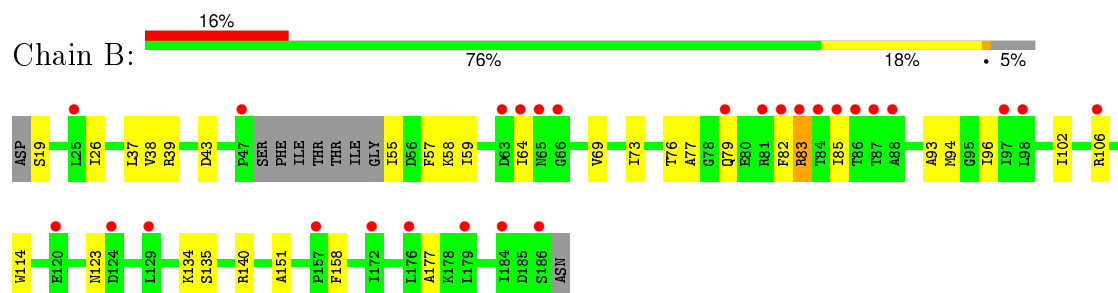
3 Residue-property plots

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.

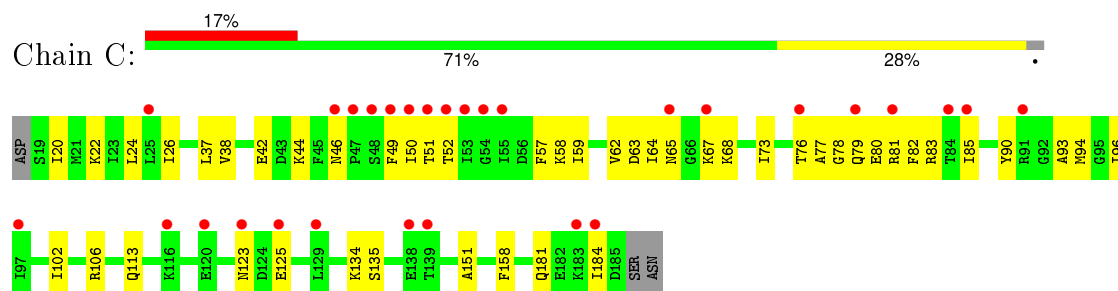
• Molecule 1: RAS-RELATED PROTEIN SEC4



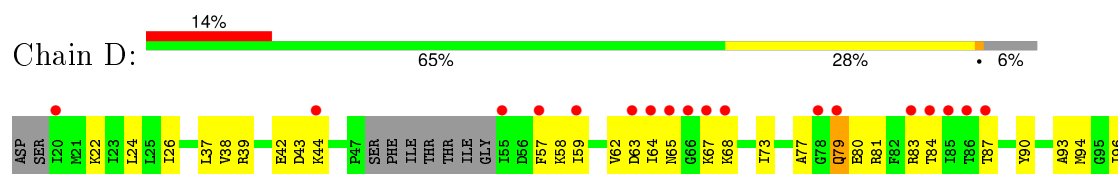
• Molecule 1: RAS-RELATED PROTEIN SEC4



• Molecule 1: RAS-RELATED PROTEIN SEC4



• Molecule 1: RAS-RELATED PROTEIN SEC4





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	56.01Å 56.38Å 59.37Å 95.25° 101.68° 116.20°	Depositor
Resolution (Å)	28.93 – 1.80 28.93 – 1.80	Depositor EDS
% Data completeness (in resolution range)	88.2 (28.93-1.80) 87.1 (28.93-1.80)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.07 (at 1.80Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.276 , 0.296 0.282 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	22.1	Xtriage
Anisotropy	0.180	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 52.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 109369 reflections	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	5593	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 44.05 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 1.6134e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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: GDP, CO

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	0.35	0/1280	0.60	0/1718
1	B	0.36	0/1294	0.57	0/1737
1	C	0.35	0/1341	0.58	0/1803
1	D	0.35	0/1288	0.56	0/1729
All	All	0.35	0/5203	0.58	0/6987

There are no bond length outliers.

There are no bond angle outliers.

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	1266	0	1290	35	0
1	B	1280	0	1299	33	0
1	C	1325	0	1348	51	0
1	D	1274	0	1294	40	0
2	A	2	0	0	0	0
2	B	1	0	0	0	0
2	C	2	0	0	0	0
2	D	3	0	0	0	0
3	A	28	0	12	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	28	0	12	1	0
3	C	28	0	12	0	0
3	D	28	0	12	0	0
4	A	105	0	0	2	0
4	B	86	0	0	3	0
4	C	70	0	0	0	0
4	D	67	0	0	5	0
All	All	5593	0	5279	153	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 15.

All (153) 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:B:85:ILE:H	1:B:85:ILE:HD12	1.21	1.05
1:C:181:GLN:O	1:C:184:ILE:HG22	1.64	0.96
1:C:46:ASN:HB2	1:C:49:PHE:HB2	1.50	0.91
1:C:79:GLN:HB2	1:D:171:GLU:OE2	1.78	0.82
1:A:37:LEU:HD13	1:A:73:ILE:HG21	1.64	0.80
1:C:181:GLN:NE2	1:C:184:ILE:HG21	1.96	0.80
1:D:81:ARG:O	1:D:84:THR:HG22	1.82	0.79
1:C:123:ASN:HD22	1:C:125:GLU:HG2	1.47	0.78
1:B:64:ILE:HD11	1:B:177:ALA:HB1	1.67	0.76
1:A:102:ILE:CD1	1:A:135:SER:HA	2.17	0.75
1:D:102:ILE:CD1	1:D:135:SER:HA	2.19	0.72
1:B:102:ILE:CD1	1:B:135:SER:HA	2.20	0.71
1:D:37:LEU:HD13	1:D:73:ILE:HG21	1.72	0.70
1:D:93:ALA:O	1:D:94:MSE:HE2	1.91	0.70
1:D:87:THR:HG22	1:D:121:HIS:HD2	1.55	0.70
1:C:102:ILE:CD1	1:C:135:SER:HA	2.21	0.69
1:C:37:LEU:HD13	1:C:73:ILE:HG21	1.73	0.69
1:C:79:GLN:HB2	1:D:171:GLU:CD	2.13	0.68
1:B:85:ILE:H	1:B:85:ILE:CD1	1.98	0.68
1:C:123:ASN:HD22	1:C:125:GLU:CG	2.06	0.67
1:B:37:LEU:HD13	1:B:73:ILE:HG21	1.77	0.67
1:A:102:ILE:HD13	1:A:135:SER:HA	1.76	0.66
1:D:102:ILE:HD13	1:D:135:SER:HA	1.77	0.66
1:C:94:MSE:HE1	1:C:123:ASN:ND2	2.11	0.66
1:C:79:GLN:HE22	1:D:168:ASN:ND2	1.92	0.66
1:C:82:PHE:O	1:C:85:ILE:HG12	1.96	0.66

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:LYS:NZ	1:A:46:ASN:HB2	2.11	0.66
1:C:102:ILE:HD13	1:C:135:SER:HA	1.79	0.65
1:C:64:ILE:O	1:C:67:LYS:HG2	1.97	0.65
1:B:93:ALA:O	1:B:94:MSE:HE2	1.98	0.64
1:C:83:ARG:NH1	1:D:178:LYS:HD3	2.13	0.64
1:B:64:ILE:HD12	1:B:69:VAL:HG21	1.80	0.62
1:D:42:GLU:O	1:D:44:LYS:HG3	1.99	0.62
1:D:138:GLU:HG2	4:D:660:HOH:O	1.99	0.62
1:B:102:ILE:HD13	1:B:135:SER:HA	1.80	0.62
1:B:79:GLN:HG2	1:B:82:PHE:HE2	1.64	0.61
1:B:85:ILE:HD11	1:B:114:TRP:HZ3	1.63	0.61
1:A:102:ILE:HD11	1:A:135:SER:HA	1.83	0.61
1:A:44:LYS:HZ1	1:A:46:ASN:HB2	1.66	0.61
1:D:170:ASN:ND2	4:D:424:HOH:O	2.32	0.61
1:C:42:GLU:O	1:C:44:LYS:HG3	2.03	0.59
1:B:82:PHE:CE1	1:B:83:ARG:HD2	2.38	0.58
1:D:102:ILE:HD11	1:D:135:SER:HA	1.85	0.57
1:A:39:ARG:O	1:A:43:ASP:HA	2.04	0.57
1:B:55:ILE:HG13	1:B:55:ILE:O	2.05	0.56
1:B:102:ILE:HD11	1:B:135:SER:HA	1.86	0.56
1:A:93:ALA:O	1:A:94:MSE:HE2	2.06	0.56
1:C:93:ALA:O	1:C:94:MSE:HE2	2.05	0.56
1:B:79:GLN:HG2	1:B:82:PHE:CE2	2.40	0.56
1:B:39:ARG:O	1:B:43:ASP:HA	2.05	0.56
1:D:65:ASN:ND2	1:D:181:GLN:HG2	2.22	0.55
1:D:62:VAL:HG12	1:D:64:ILE:HG23	1.88	0.55
1:B:19:SER:C	4:B:549:HOH:O	2.45	0.55
1:D:39:ARG:O	1:D:43:ASP:HA	2.07	0.54
1:C:102:ILE:HD11	1:C:135:SER:HA	1.88	0.54
1:C:62:VAL:HG23	1:C:64:ILE:HD13	1.89	0.53
1:B:85:ILE:HD12	1:B:85:ILE:N	2.06	0.53
1:B:76:THR:HG23	4:B:594:HOH:O	2.08	0.53
1:C:42:GLU:HB3	1:C:44:LYS:HE3	1.91	0.53
1:B:58:LYS:C	1:B:59:ILE:HD12	2.30	0.52
1:B:140:ARG:HD2	1:C:20:ILE:HD12	1.92	0.52
1:A:55:ILE:N	4:A:518:HOH:O	2.42	0.52
1:C:63:ASP:O	1:C:64:ILE:HD12	2.09	0.51
1:B:94:MSE:HE1	1:B:123:ASN:CB	2.41	0.51
1:A:24:LEU:HG	1:A:89:TYR:CE2	2.46	0.51
1:A:44:LYS:NZ	1:A:46:ASN:CB	2.74	0.51
1:B:64:ILE:CD1	1:B:177:ALA:HB1	2.38	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:37:LEU:HD13	1:A:73:ILE:CG2	2.38	0.50
1:C:63:ASP:OD1	1:C:68:LYS:HD3	2.12	0.50
1:A:58:LYS:C	1:A:59:ILE:HD12	2.32	0.49
1:B:77:ALA:O	1:B:79:GLN:HG3	2.12	0.49
1:D:174:PHE:O	1:D:178:LYS:HG3	2.13	0.49
1:C:58:LYS:C	1:C:59:ILE:HD12	2.33	0.49
1:A:26:ILE:HD12	1:A:96:ILE:CG2	2.43	0.49
1:C:64:ILE:O	1:C:65:ASN:HB2	2.13	0.48
1:A:62:VAL:CG1	1:A:63:ASP:N	2.76	0.48
1:A:44:LYS:HD3	1:A:45:PHE:N	2.28	0.48
1:A:180:ILE:O	1:A:184:ILE:HG13	2.14	0.48
1:A:98:LEU:N	1:A:98:LEU:HD12	2.28	0.48
1:C:64:ILE:O	1:C:67:LYS:HE2	2.14	0.47
1:C:63:ASP:C	1:C:64:ILE:HD12	2.34	0.47
1:B:82:PHE:HB3	4:B:566:HOH:O	2.15	0.47
1:A:86:THR:O	1:A:89:TYR:N	2.29	0.47
1:C:49:PHE:O	1:C:50:ILE:HD13	2.14	0.47
1:A:64:ILE:HD11	1:A:177:ALA:HB1	1.97	0.47
1:D:62:VAL:HG13	1:D:174:PHE:CE2	2.50	0.47
1:C:26:ILE:CD1	1:C:90:TYR:HE2	2.28	0.47
1:C:181:GLN:HE21	1:C:184:ILE:HG21	1.76	0.47
1:D:58:LYS:C	1:D:59:ILE:HD12	2.36	0.47
1:C:51:THR:HG22	1:C:52:THR:N	2.30	0.46
1:D:87:THR:HG22	1:D:121:HIS:CD2	2.43	0.46
1:A:22:LYS:HE2	1:A:24:LEU:HD11	1.97	0.46
1:D:79:GLN:CD	1:D:79:GLN:H	2.19	0.46
1:C:38:VAL:HG22	1:C:58:LYS:NZ	2.30	0.46
1:B:26:ILE:HD12	1:B:96:ILE:CG2	2.45	0.46
1:B:94:MSE:HE1	1:B:123:ASN:ND2	2.30	0.46
1:A:57:PHE:CE1	1:A:59:ILE:HD11	2.51	0.46
1:C:26:ILE:HD11	1:C:90:TYR:CE2	2.51	0.45
1:D:180:ILE:O	1:D:184:ILE:HG13	2.16	0.45
1:A:62:VAL:HG13	1:A:174:PHE:CE2	2.51	0.45
1:A:106:ARG:HB2	1:A:106:ARG:NH1	2.32	0.45
1:C:78:GLY:O	1:C:81:ARG:HB3	2.16	0.45
1:C:26:ILE:HD12	1:C:96:ILE:CG2	2.46	0.45
1:C:94:MSE:SE	1:C:125:GLU:HG3	2.66	0.45
1:D:79:GLN:NE2	1:D:79:GLN:H	2.14	0.45
1:C:76:THR:HG22	1:C:77:ALA:O	2.16	0.45
1:B:151:ALA:CB	1:B:158:PHE:HB2	2.47	0.45
1:D:80:GLU:O	1:D:83:ARG:HB3	2.17	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:67:LYS:HA	1:A:67:LYS:HD3	1.78	0.45
1:C:57:PHE:CE1	1:C:59:ILE:HD11	2.51	0.45
1:C:22:LYS:HE2	1:C:24:LEU:HD11	2.00	0.44
1:B:57:PHE:CE1	1:B:59:ILE:HD11	2.53	0.44
1:A:79:GLN:NE2	4:A:583:HOH:O	2.51	0.44
1:A:59:ILE:N	1:A:59:ILE:HD12	2.33	0.44
1:A:62:VAL:HG12	1:A:63:ASP:N	2.32	0.44
1:C:26:ILE:HD11	1:C:90:TYR:HE2	1.83	0.44
1:A:44:LYS:HZ1	1:A:46:ASN:CB	2.29	0.43
1:D:22:LYS:HE2	1:D:24:LEU:HD11	2.00	0.43
1:C:151:ALA:CB	1:C:158:PHE:HB2	2.47	0.43
1:A:38:VAL:HG22	1:A:58:LYS:NZ	2.33	0.43
1:D:67:LYS:HG2	1:D:68:LYS:N	2.34	0.43
1:D:151:ALA:CB	1:D:158:PHE:HB2	2.49	0.43
1:D:26:ILE:HD12	1:D:96:ILE:CG2	2.49	0.43
1:C:80:GLU:HG2	1:D:175:THR:OG1	2.19	0.43
1:D:57:PHE:CE1	1:D:59:ILE:HD11	2.54	0.42
1:C:123:ASN:ND2	1:C:125:GLU:CG	2.80	0.42
1:C:76:THR:CG2	1:C:77:ALA:N	2.82	0.42
1:D:26:ILE:HD11	1:D:90:TYR:CE2	2.54	0.42
1:A:151:ALA:CB	1:A:158:PHE:HB2	2.49	0.42
1:C:79:GLN:HG2	1:C:81:ARG:HH22	1.84	0.42
1:C:102:ILE:HD11	1:C:134:LYS:O	2.20	0.42
1:C:64:ILE:O	1:C:64:ILE:HG22	2.20	0.42
1:A:26:ILE:CD1	1:A:90:TYR:HE1	2.32	0.42
1:C:181:GLN:NE2	1:C:184:ILE:CG2	2.77	0.42
1:D:62:VAL:CG1	1:D:63:ASP:N	2.82	0.42
1:D:38:VAL:HG22	1:D:58:LYS:NZ	2.35	0.42
1:A:76:THR:O	1:A:79:GLN:HG3	2.19	0.42
1:D:178:LYS:O	1:D:182:GLU:HG3	2.20	0.42
1:B:59:ILE:N	1:B:59:ILE:HD12	2.35	0.42
1:C:26:ILE:HD12	1:C:96:ILE:HG23	2.02	0.42
1:C:82:PHE:CD1	1:C:85:ILE:HD11	2.54	0.42
1:B:38:VAL:HG22	1:B:58:LYS:NZ	2.35	0.42
1:A:174:PHE:O	1:A:178:LYS:HG3	2.20	0.41
1:A:70:LYS:HE3	1:A:70:LYS:HB2	1.82	0.41
1:D:77:ALA:HA	4:D:410:HOH:O	2.20	0.41
1:C:106:ARG:NH1	1:C:106:ARG:HB2	2.36	0.41
1:B:94:MSE:HE1	1:B:123:ASN:CG	2.42	0.40
1:D:59:ILE:HG22	4:D:681:HOH:O	2.21	0.40
1:B:134:LYS:HG2	3:B:202:GDP:C6	2.56	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:137:MSE:HE2	4:D:562:HOH:O	2.21	0.40
1:D:59:ILE:HD12	1:D:59:ILE:N	2.36	0.40
1:B:106:ARG:HB2	1:B:106:ARG:NH1	2.36	0.40
1:D:106:ARG:NH1	1:D:106:ARG:HB2	2.37	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	155/170 (91%)	152 (98%)	2 (1%)	1 (1%)	30	14
1	B	157/170 (92%)	152 (97%)	5 (3%)	0	100	100
1	C	165/170 (97%)	161 (98%)	4 (2%)	0	100	100
1	D	156/170 (92%)	152 (97%)	4 (3%)	0	100	100
All	All	633/680 (93%)	617 (98%)	15 (2%)	1 (0%)	52	35

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	87	THR

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	139/146 (95%)	139 (100%)	0	100	100
1	B	141/146 (97%)	140 (99%)	1 (1%)	88	86
1	C	146/146 (100%)	145 (99%)	1 (1%)	88	86
1	D	140/146 (96%)	138 (99%)	2 (1%)	74	65
All	All	566/584 (97%)	562 (99%)	4 (1%)	88	86

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	83	ARG
1	C	113	GLN
1	D	79	GLN
1	D	124	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	165	ASN
1	B	65	ASN
1	B	79	GLN
1	B	165	ASN
1	C	123	ASN
1	C	165	ASN
1	C	181	GLN
1	D	65	ASN
1	D	79	GLN
1	D	123	ASN
1	D	165	ASN
1	D	168	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 8 are monoatomic - leaving 4 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	GDP	A	201	2	23,30,30	2.71	8 (34%)	30,47,47	3.06	10 (33%)
3	GDP	B	202	2	23,30,30	2.96	9 (39%)	30,47,47	3.03	10 (33%)
3	GDP	C	203	2	23,30,30	2.94	8 (34%)	30,47,47	3.01	10 (33%)
3	GDP	D	204	2	23,30,30	2.94	9 (39%)	30,47,47	3.06	9 (30%)

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	GDP	A	201	2	-	0/12/32/32	0/3/3/3
3	GDP	B	202	2	-	0/12/32/32	0/3/3/3
3	GDP	C	203	2	-	0/12/32/32	0/3/3/3
3	GDP	D	204	2	-	0/12/32/32	0/3/3/3

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	202	GDP	PB-O2B	-4.03	1.40	1.54
3	D	204	GDP	PB-O2B	-3.96	1.40	1.54
3	C	203	GDP	PB-O2B	-3.20	1.43	1.54
3	A	201	GDP	PB-O2B	-2.48	1.45	1.54
3	C	203	GDP	C3'-C4'	2.03	1.58	1.53
3	B	202	GDP	O3'-C3'	2.05	1.47	1.43

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	204	GDP	O3'-C3'	2.14	1.48	1.43
3	C	203	GDP	C5-C4	2.15	1.45	1.40
3	A	201	GDP	O4'-C4'	2.21	1.50	1.45
3	A	201	GDP	O3'-C3'	2.25	1.48	1.43
3	B	202	GDP	C5-C4	2.30	1.45	1.40
3	A	201	GDP	C5-C4	2.32	1.45	1.40
3	C	203	GDP	O4'-C4'	2.46	1.50	1.45
3	B	202	GDP	PB-O3B	2.48	1.63	1.54
3	D	204	GDP	C5-C4	2.65	1.46	1.40
3	D	204	GDP	O4'-C4'	2.66	1.51	1.45
3	D	204	GDP	PB-O3B	2.68	1.64	1.54
3	B	202	GDP	O4'-C4'	2.91	1.51	1.45
3	C	203	GDP	C8-N7	3.22	1.40	1.34
3	D	204	GDP	C8-N7	3.24	1.40	1.34
3	A	201	GDP	C8-N7	3.32	1.41	1.34
3	B	202	GDP	C8-N7	3.67	1.41	1.34
3	A	201	GDP	O6-C6	3.96	1.34	1.24
3	B	202	GDP	O6-C6	4.14	1.34	1.24
3	D	204	GDP	O6-C6	4.17	1.34	1.24
3	C	203	GDP	O6-C6	4.18	1.34	1.24
3	A	201	GDP	O4'-C1'	6.45	1.49	1.41
3	D	204	GDP	O4'-C1'	6.49	1.49	1.41
3	B	202	GDP	O4'-C1'	6.60	1.49	1.41
3	C	203	GDP	O4'-C1'	6.85	1.49	1.41
3	A	201	GDP	C2-N1	7.83	1.49	1.35
3	B	202	GDP	C2-N1	8.54	1.50	1.35
3	D	204	GDP	C2-N1	8.65	1.50	1.35
3	C	203	GDP	C2-N1	8.84	1.51	1.35

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	201	GDP	C6-C5-C4	-9.63	109.39	120.90
3	C	203	GDP	C6-C5-C4	-9.49	109.56	120.90
3	D	204	GDP	C6-C5-C4	-9.36	109.71	120.90
3	B	202	GDP	C6-C5-C4	-9.34	109.73	120.90
3	A	201	GDP	C4-C5-N7	-5.87	104.08	109.48
3	C	203	GDP	C4-C5-N7	-5.85	104.09	109.48
3	B	202	GDP	C4-C5-N7	-5.50	104.42	109.48
3	D	204	GDP	C4-C5-N7	-5.45	104.46	109.48
3	D	204	GDP	N2-C2-N1	-5.35	108.34	117.20
3	B	202	GDP	N2-C2-N1	-5.23	108.53	117.20

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	201	GDP	N2-C2-N1	-5.20	108.59	117.20
3	D	204	GDP	N3-C2-N1	-5.03	119.78	127.44
3	C	203	GDP	N2-C2-N1	-4.96	108.98	117.20
3	B	202	GDP	N3-C2-N1	-4.93	119.93	127.44
3	C	203	GDP	N3-C2-N1	-4.89	120.00	127.44
3	A	201	GDP	N3-C2-N1	-4.84	120.07	127.44
3	A	201	GDP	C1'-N9-C4	-2.22	123.59	126.94
3	B	202	GDP	C1'-N9-C4	-2.21	123.61	126.94
3	C	203	GDP	C1'-N9-C4	-2.21	123.61	126.94
3	B	202	GDP	O2B-PB-O1B	2.15	117.49	110.58
3	A	201	GDP	O2B-PB-O1B	2.23	117.76	110.58
3	C	203	GDP	C2'-C3'-C4'	2.25	107.24	102.61
3	A	201	GDP	C2'-C3'-C4'	2.38	107.50	102.61
3	D	204	GDP	C2'-C3'-C4'	2.38	107.51	102.61
3	A	201	GDP	O2'-C2'-C3'	2.39	119.61	111.83
3	D	204	GDP	O2B-PB-O1B	2.43	118.40	110.58
3	C	203	GDP	O2B-PB-O1B	2.44	118.44	110.58
3	B	202	GDP	O2'-C2'-C3'	2.54	120.08	111.83
3	D	204	GDP	O2'-C2'-C3'	2.58	120.22	111.83
3	B	202	GDP	C2'-C3'-C4'	2.64	108.03	102.61
3	C	203	GDP	O2'-C2'-C3'	2.95	121.43	111.83
3	C	203	GDP	C4'-O4'-C1'	3.01	113.03	109.72
3	B	202	GDP	C4'-O4'-C1'	3.55	113.61	109.72
3	D	204	GDP	C4'-O4'-C1'	3.55	113.62	109.72
3	A	201	GDP	C4'-O4'-C1'	3.69	113.77	109.72
3	C	203	GDP	N2-C2-N3	6.89	131.02	117.80
3	A	201	GDP	N2-C2-N3	7.06	131.35	117.80
3	B	202	GDP	N2-C2-N3	7.16	131.54	117.80
3	D	204	GDP	N2-C2-N3	7.34	131.88	117.80

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
3	B	202	GDP	1	0

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	156/170 (91%)	0.96	22 (14%) 4 3	16, 29, 48, 56	0
1	B	158/170 (92%)	0.97	27 (17%) 2 1	16, 30, 49, 53	0
1	C	164/170 (96%)	1.12	29 (17%) 2 1	17, 32, 50, 54	0
1	D	157/170 (92%)	0.98	23 (14%) 3 2	15, 31, 49, 55	0
All	All	635/680 (93%)	1.01	101 (15%) 3 2	15, 30, 49, 56	0

All (101) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	51	THR	9.7
1	C	184	ILE	9.6
1	C	47	PRO	6.0
1	A	47	PRO	5.6
1	C	53	ILE	5.6
1	C	49	PHE	5.6
1	D	78	GLY	5.4
1	A	64	ILE	5.4
1	D	55	ILE	5.2
1	A	20	ILE	5.1
1	C	79	GLN	5.0
1	D	65	ASN	4.9
1	B	66	GLY	4.8
1	A	62	VAL	4.6
1	A	65	ASN	4.6
1	A	87	THR	4.4
1	B	84	THR	4.3
1	B	82	PHE	4.3
1	A	67	LYS	4.3
1	D	83	ARG	4.2
1	B	87	THR	4.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	46	ASN	4.0
1	D	85	ILE	3.9
1	A	86	THR	3.9
1	D	57	PHE	3.8
1	A	66	GLY	3.8
1	C	55	ILE	3.8
1	D	68	LYS	3.8
1	D	79	GLN	3.7
1	A	55	ILE	3.7
1	A	124	ASP	3.6
1	C	183	LYS	3.6
1	C	81	ARG	3.5
1	B	86	THR	3.5
1	B	65	ASN	3.5
1	A	46	ASN	3.4
1	B	47	PRO	3.4
1	C	139	THR	3.4
1	D	84	THR	3.3
1	C	138	GLU	3.3
1	C	50	ILE	3.2
1	C	48	SER	3.2
1	B	184	ILE	3.2
1	C	54	GLY	3.1
1	A	19	SER	3.1
1	B	129	LEU	3.1
1	C	123	ASN	3.1
1	D	87	THR	3.1
1	D	67	LYS	3.0
1	C	76	THR	3.0
1	D	64	ILE	3.0
1	C	65	ASN	3.0
1	D	63	ASP	2.9
1	A	25	LEU	2.9
1	D	184	ILE	2.9
1	C	85	ILE	2.8
1	C	84	THR	2.8
1	D	20	ILE	2.7
1	B	85	ILE	2.6
1	A	174	PHE	2.6
1	B	172	ILE	2.6
1	B	88	ALA	2.6
1	D	123	ASN	2.6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	64	ILE	2.6
1	B	186	SER	2.6
1	A	184	ILE	2.5
1	D	86	THR	2.5
1	B	97	ILE	2.5
1	D	59	ILE	2.5
1	C	25	LEU	2.5
1	C	97	ILE	2.5
1	A	178	LYS	2.5
1	B	81	ARG	2.4
1	A	63	ASP	2.4
1	A	26	ILE	2.4
1	C	120	GLU	2.4
1	A	73	ILE	2.4
1	C	129	LEU	2.4
1	D	66	GLY	2.4
1	B	120	GLU	2.4
1	C	125	GLU	2.4
1	D	175	THR	2.4
1	A	180	ILE	2.4
1	B	124	ASP	2.3
1	B	176	LEU	2.3
1	B	179	LEU	2.2
1	C	67	LYS	2.2
1	A	88	ALA	2.2
1	C	116	LYS	2.2
1	C	91	ARG	2.2
1	B	106	ARG	2.2
1	D	44	LYS	2.2
1	D	138	GLU	2.2
1	B	25	LEU	2.2
1	B	83	ARG	2.2
1	C	52	THR	2.1
1	B	98	LEU	2.1
1	B	63	ASP	2.1
1	B	157	PRO	2.1
1	B	79	GLN	2.1
1	D	131	VAL	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 [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
3	GDP	B	202	28/28	0.90	0.14	-0.08	29,32,37,43	0
3	GDP	D	204	28/28	0.94	0.12	-0.17	22,26,34,36	0
3	GDP	A	201	28/28	0.96	0.10	-0.54	16,19,25,27	0
3	GDP	C	203	28/28	0.97	0.08	-0.84	17,26,32,36	0
2	CO	C	302	1/1	0.97	0.04	-1.17	12,12,12,12	0
2	CO	D	307	1/1	0.95	0.08	-1.41	32,32,32,32	0
2	CO	D	308	1/1	0.90	0.05	-2.01	40,40,40,40	0
2	CO	B	304	1/1	0.93	0.04	-3.45	24,24,24,24	0
2	CO	A	301	1/1	0.99	0.02	-4.86	10,10,10,10	0
2	CO	D	303	1/1	0.99	0.07	-	19,19,19,19	0
2	CO	C	306	1/1	0.98	0.02	-	32,32,32,32	0
2	CO	A	305	1/1	0.96	0.03	-	30,30,30,30	0

6.5 Other polymers [i](#)

There are no such residues in this entry.