



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 05:59 PM UTC

PDB ID : 3DM5 / pdb_00003dm5
Title : Structures of SRP54 and SRP19, the two proteins assembling the ribonucleic core of the Signal Recognition Particle from the archaeon *Pyrococcus furiosus*.
Authors : Egea, P.F.; Napetschnig, J.; Walter, P.; Stroud, R.M.
Deposited on : 2008-06-30
Resolution : 2.51 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

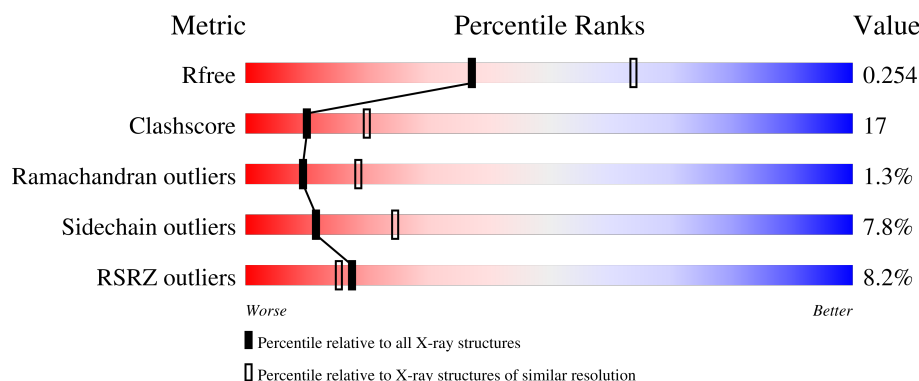
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.51 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}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	443	
1	B	443	

2 Entry composition [i](#)

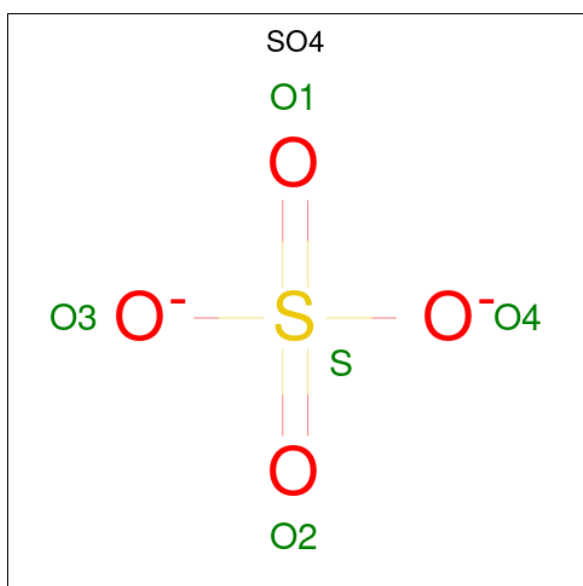
There are 5 unique types of molecules in this entry. The entry contains 6649 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Signal recognition 54 kDa protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	416	Total	C	N	O	S	0	0	0
			3192	2028	552	601	11			
1	B	413	Total	C	N	O	S	11	1	0
			3196	2034	556	595	11			

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



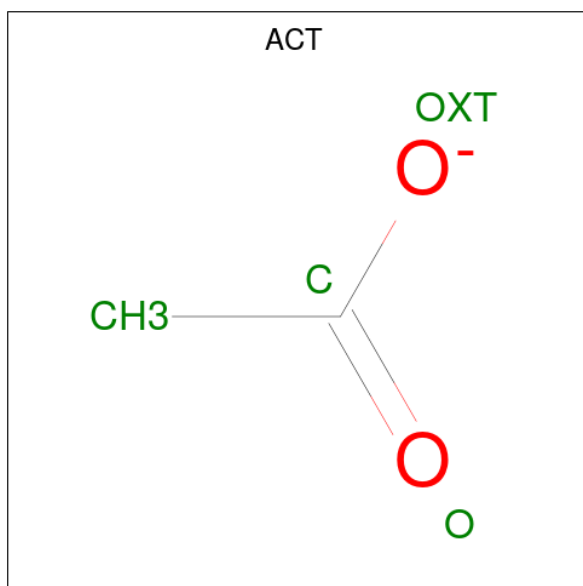
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		

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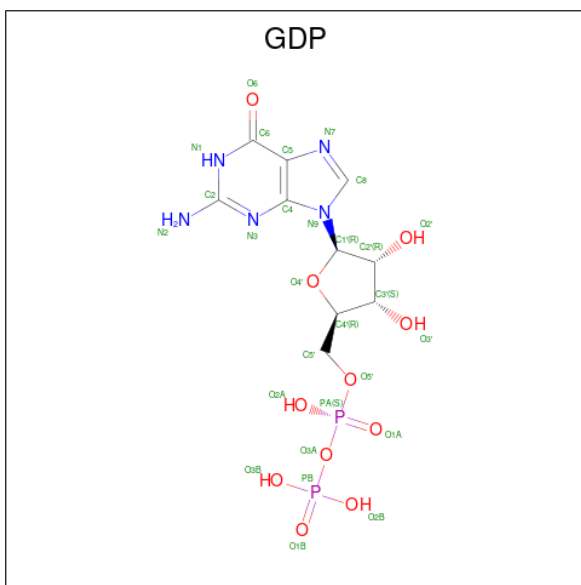
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



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

- Molecule 4 is GUANOSINE-5'-DIPHOSPHATE (CCD ID: GDP) (formula: $C_{10}H_{15}N_5O_{11}P_2$).



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

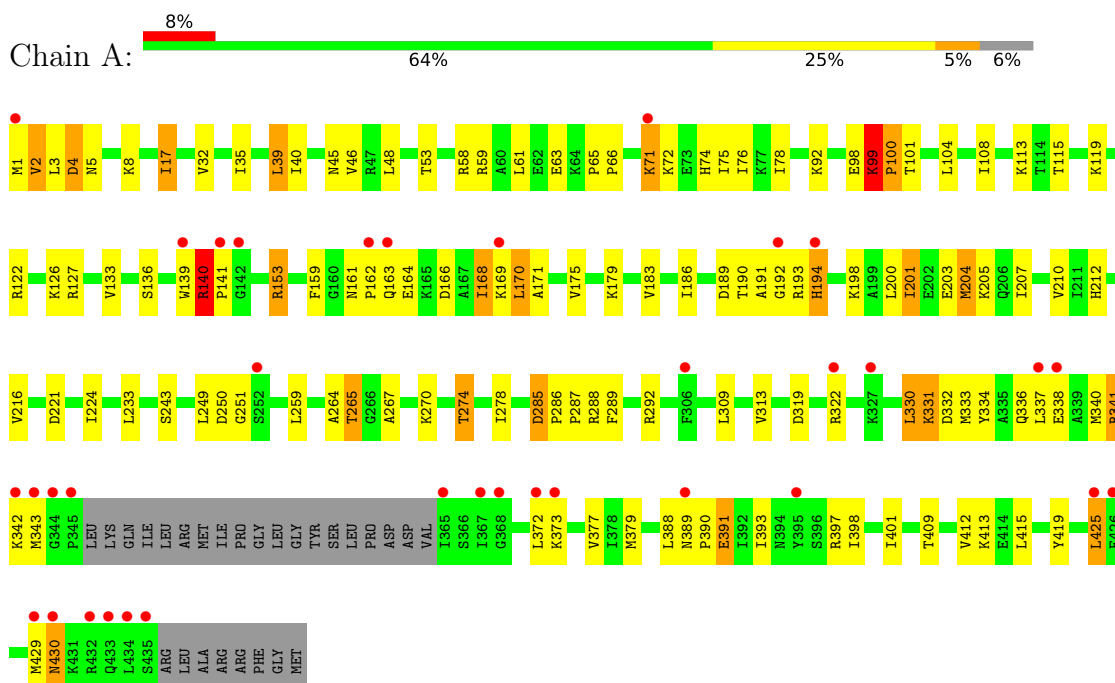
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	72	Total O 72 72	0	0
5	B	70	Total O 70 70	0	0

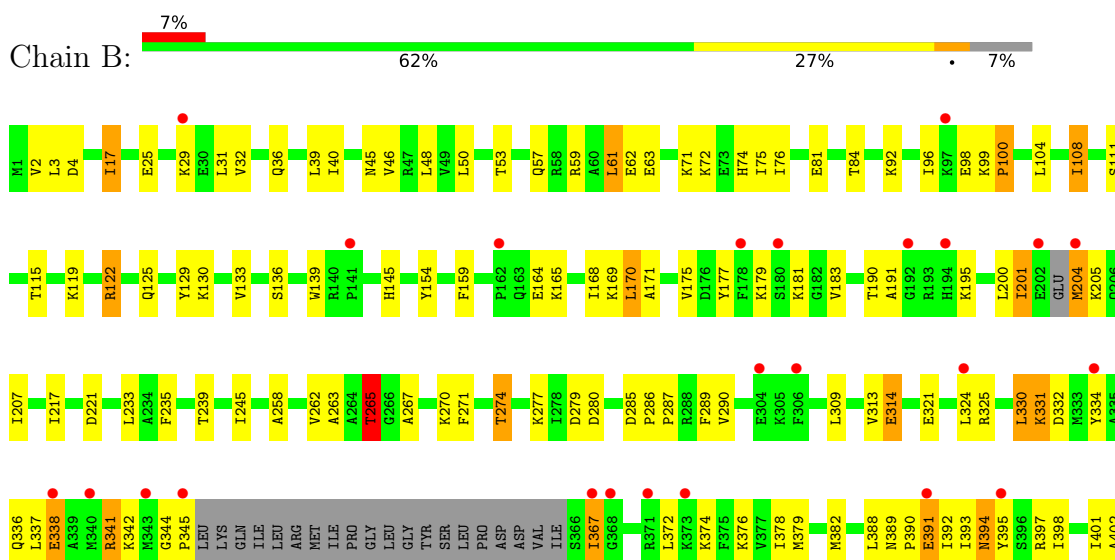
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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: Signal recognition 54 kDa protein



• Molecule 1: Signal recognition 54 kDa protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 42 21 2	Depositor
Cell constants a, b, c, α , β , γ	127.01Å 127.01Å 186.84Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	64.74 – 2.51 64.74 – 2.51	Depositor EDS
% Data completeness (in resolution range)	99.5 (64.74-2.51) 99.5 (64.74-2.51)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.76 (at 2.51Å)	Xtriage
Refinement program	PHENIX (phenix.refine), ELVES	Depositor
R, R_{free}	0.222 , 0.259 0.216 , 0.254	Depositor DCC
R_{free} test set	2000 reflections (3.80%)	wwPDB-VP
Wilson B-factor (Å ²)	55.2	Xtriage
Anisotropy	0.163	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 56.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6649	wwPDB-VP
Average B, all atoms (Å ²)	65.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.33% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ACT, GDP, 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	0.52	0/3234	0.91	8/4357 (0.2%)
1	B	0.52	0/3235	0.86	1/4348 (0.0%)
All	All	0.52	0/6469	0.88	9/8705 (0.1%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	265	THR	N-CA-C	-8.46	103.39	114.31
1	A	99	LYS	CA-C-N	-7.60	110.34	119.84
1	A	99	LYS	C-N-CA	-7.60	110.34	119.84
1	A	140	ARG	CA-C-N	5.97	127.30	119.84
1	A	140	ARG	C-N-CA	5.97	127.30	119.84
1	A	285	ASP	CA-C-N	5.68	126.23	120.38
1	A	285	ASP	C-N-CA	5.68	126.23	120.38
1	A	264	ALA	N-CA-C	-5.49	106.42	113.01
1	A	313	VAL	N-CA-C	5.32	116.05	110.62

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3192	0	3248	121	0
1	B	3196	0	3289	105	0
2	A	30	0	0	1	0
2	B	25	0	0	2	0
3	A	4	0	3	0	0
3	B	4	0	3	0	0
4	A	28	0	12	0	0
4	B	28	0	12	0	0
5	A	72	0	0	1	0
5	B	70	0	0	2	0
All	All	6649	0	6567	226	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 17.

All (226) 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:341:ARG:HG3	1:A:341:ARG:HH11	1.27	0.99
1:A:99:LYS:O	1:A:100:PRO:C	2.09	0.95
1:B:341:ARG:HG3	1:B:341:ARG:HH11	1.36	0.91
1:B:330:LEU:HD12	1:B:390:PRO:HG3	1.54	0.86
1:A:265:THR:HG22	1:A:267:ALA:H	1.40	0.85
1:B:324:LEU:HD22	1:B:429:MET:HB2	1.60	0.83
1:A:127:ARG:HG3	1:A:127:ARG:HH11	1.43	0.82
1:A:99:LYS:HG3	1:A:100:PRO:HD3	1.62	0.81
1:B:99:LYS:HG3	1:B:100:PRO:HD3	1.61	0.81
1:B:168:ILE:HD13	1:B:207:ILE:HG12	1.63	0.80
1:A:168:ILE:HD13	1:A:207:ILE:HG12	1.66	0.77
1:B:341:ARG:HG3	1:B:341:ARG:NH1	1.97	0.77
1:B:425:LEU:HD22	1:B:429:MET:HE2	1.67	0.77
1:A:113:LYS:HE2	1:A:189:ASP:OD1	1.85	0.76
1:A:341:ARG:HG3	1:A:341:ARG:NH1	1.97	0.74
1:A:330:LEU:HD12	1:A:390:PRO:HG3	1.70	0.72
1:A:45:ASN:HD22	1:A:48:LEU:H	1.38	0.72
1:B:99:LYS:O	1:B:100:PRO:C	2.33	0.72
1:A:193:ARG:HG3	1:A:200:LEU:HD12	1.71	0.71
1:B:398:ILE:HG23	1:B:412:VAL:HG11	1.74	0.70
1:A:92:LYS:O	1:A:270:LYS:HD3	1.93	0.68
1:A:265:THR:CG2	1:A:267:ALA:H	2.07	0.67
1:A:153:ARG:HG2	5:A:505:HOH:O	1.95	0.66
1:A:2:VAL:H	1:A:250:ASP:HA	1.61	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:390:PRO:HG2	1:A:419:TYR:CZ	2.29	0.66
1:B:71:LYS:O	1:B:75:ILE:HG12	1.96	0.66
1:B:104:LEU:HD11	1:B:190:THR:CG2	2.26	0.66
1:B:111:SER:HA	1:B:221:ASP:HB2	1.77	0.66
1:B:334:TYR:OH	1:B:376:LYS:HE2	1.96	0.65
1:B:382:MET:HG2	1:B:401:ILE:HD13	1.79	0.65
1:A:136:SER:HA	1:A:170:LEU:HD13	1.79	0.65
1:B:330:LEU:HD21	1:B:415:LEU:HD21	1.80	0.64
1:A:164:GLU:OE1	1:A:164:GLU:HA	1.97	0.63
1:A:341:ARG:HD3	1:A:372:LEU:HD22	1.80	0.63
1:A:1:MET:O	1:A:2:VAL:HG12	1.99	0.62
1:B:136:SER:HA	1:B:170:LEU:HD13	1.81	0.62
1:A:59:ARG:O	1:A:63:GLU:HB2	1.98	0.62
1:A:32:VAL:HG13	1:A:53:THR:HB	1.81	0.62
1:A:99:LYS:O	1:A:101:THR:N	2.32	0.62
1:B:245:ILE:HG12	1:B:271:PHE:HB2	1.81	0.62
1:A:203:GLU:O	1:A:207:ILE:HG13	1.99	0.62
1:B:100:PRO:HG2	1:B:179:LYS:HG3	1.81	0.61
1:B:324:LEU:CD2	1:B:429:MET:HB2	2.30	0.61
1:A:40:ILE:HD12	1:A:46:VAL:HG22	1.83	0.60
1:A:166:ASP:HB3	1:A:169:LYS:HB2	1.83	0.60
1:B:277:LYS:HE2	5:B:535:HOH:O	2.01	0.60
1:B:341:ARG:HH11	1:B:341:ARG:CG	2.14	0.60
1:B:398:ILE:HG23	1:B:412:VAL:CG1	2.32	0.60
1:A:127:ARG:HH11	1:A:127:ARG:CG	2.15	0.60
1:B:63:GLU:OE1	1:B:74:HIS:HE1	1.85	0.60
1:B:45:ASN:HD22	1:B:48:LEU:H	1.51	0.59
1:B:330:LEU:HD12	1:B:390:PRO:CG	2.31	0.59
1:A:98:GLU:C	1:A:99:LYS:O	2.43	0.59
1:A:285:ASP:OD1	1:A:287:PRO:HD2	2.03	0.59
1:B:337:LEU:HB3	1:B:372:LEU:HD12	1.85	0.58
1:A:71:LYS:HD2	1:A:71:LYS:H	1.68	0.58
1:A:286:PRO:HB2	1:A:287:PRO:HD3	1.85	0.58
1:A:139:TRP:HB2	1:A:191:ALA:HB1	1.85	0.57
1:A:58:ARG:HH12	1:A:59:ARG:HG2	1.69	0.57
1:A:99:LYS:CG	1:A:100:PRO:HD3	2.35	0.56
1:B:25:GLU:O	1:B:29:LYS:HG2	2.04	0.56
1:B:274:THR:HG22	1:B:280:ASP:HB3	1.87	0.56
1:A:127:ARG:HG3	1:A:127:ARG:NH1	2.19	0.56
1:B:29:LYS:HD2	1:B:57:GLN:HE21	1.71	0.55
1:B:59:ARG:O	1:B:63:GLU:HB2	2.05	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:3:LEU:HD22	1:A:289:PHE:HZ	1.72	0.54
1:B:139:TRP:HB2	1:B:191:ALA:HB1	1.89	0.53
1:A:2:VAL:H	1:A:251:GLY:H	1.57	0.53
1:A:249:LEU:HD11	1:A:259:LEU:HD21	1.91	0.53
1:B:342:LYS:C	1:B:344:GLY:H	2.16	0.53
1:B:100:PRO:CG	1:B:179:LYS:HG3	2.38	0.53
1:A:100:PRO:HG2	1:A:179:LYS:HG3	1.90	0.52
1:B:96:ILE:HG21	1:B:129:TYR:CE1	2.45	0.52
1:A:99:LYS:HB3	1:A:100:PRO:CD	2.39	0.52
1:A:265:THR:HG22	1:A:267:ALA:N	2.18	0.52
1:B:263:ALA:C	1:B:265:THR:H	2.17	0.52
1:B:430:ASN:C	1:B:432:ARG:H	2.16	0.52
1:B:334:TYR:HD1	1:B:379:MET:CE	2.23	0.52
1:B:130:LYS:HG2	1:B:183:VAL:HG12	1.90	0.52
1:A:104:LEU:HD11	1:A:190:THR:HG21	1.91	0.52
1:A:35:ILE:HG22	1:A:39:LEU:HD22	1.92	0.52
1:A:40:ILE:CD1	1:A:46:VAL:HG22	2.40	0.51
1:A:192:GLY:O	1:A:193:ARG:HG2	2.10	0.51
1:B:394:ASN:O	1:B:398:ILE:HG13	2.10	0.51
1:B:262:VAL:O	1:B:265:THR:HG22	2.11	0.51
1:A:340:MET:SD	1:A:343:MET:HE1	2.50	0.51
1:B:92:LYS:O	1:B:270:LYS:HD3	2.10	0.51
1:A:341:ARG:NH1	1:A:341:ARG:CG	2.72	0.51
1:A:288:ARG:O	1:A:292:ARG:HG2	2.10	0.51
1:B:164:GLU:HG2	1:B:170:LEU:HG	1.93	0.51
1:A:161:ASN:C	1:A:163:GLN:H	2.19	0.50
1:A:2:VAL:O	1:A:4:ASP:N	2.42	0.50
1:B:334:TYR:HE1	1:B:376:LYS:HG3	1.77	0.50
1:B:338:GLU:O	1:B:341:ARG:HG2	2.11	0.50
1:B:84:THR:HA	1:B:290:VAL:HG21	1.94	0.50
1:B:59:ARG:HH12	1:B:81:GLU:CD	2.20	0.50
1:A:198:LYS:HE2	2:A:448:SO4:O4	2.11	0.50
1:B:36:GLN:OE1	1:B:53:THR:HG21	2.11	0.49
1:B:334:TYR:HD1	1:B:379:MET:HE3	1.77	0.49
1:A:74:HIS:O	1:A:78:ILE:HG13	2.12	0.49
1:B:99:LYS:HG3	1:B:100:PRO:CD	2.37	0.49
1:B:235:PHE:O	1:B:239:THR:HG22	2.11	0.49
1:B:397:ARG:O	1:B:401:ILE:HG13	2.12	0.49
1:A:122:ARG:CZ	1:A:126:LYS:HG3	2.41	0.49
1:B:285:ASP:OD1	1:B:287:PRO:HD2	2.13	0.49
1:A:17:ILE:CD1	1:A:75:ILE:HG21	2.43	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:258:ALA:O	1:B:262:VAL:HG23	2.13	0.48
1:A:104:LEU:HD23	1:A:216:VAL:HG22	1.95	0.48
1:A:115:THR:HG22	1:A:119:LYS:HE3	1.95	0.48
1:A:161:ASN:C	1:A:163:GLN:N	2.72	0.48
1:B:36:GLN:CG	1:B:53:THR:HG21	2.44	0.48
1:A:164:GLU:CG	1:A:170:LEU:HG	2.43	0.48
1:B:425:LEU:CD2	1:B:429:MET:HE2	2.41	0.48
1:A:127:ARG:CG	1:A:127:ARG:NH1	2.73	0.48
1:B:331:LYS:HE3	1:B:388:LEU:HD21	1.96	0.47
1:B:390:PRO:HG2	1:B:419:TYR:CZ	2.50	0.47
1:A:99:LYS:HD2	1:A:99:LYS:HA	1.49	0.47
1:B:332:ASP:O	1:B:336:GLN:HG2	2.15	0.47
1:A:104:LEU:HD11	1:A:190:THR:CG2	2.44	0.47
1:A:397:ARG:O	1:A:401:ILE:HG13	2.15	0.47
1:A:71:LYS:H	1:A:71:LYS:CD	2.23	0.47
1:A:331:LYS:HE3	1:A:388:LEU:HD21	1.96	0.47
1:B:390:PRO:O	1:B:393:ILE:HG13	2.15	0.47
1:A:161:ASN:O	1:A:163:GLN:N	2.48	0.47
1:A:204:MET:HG3	1:A:205:LYS:N	2.30	0.47
1:A:243:SER:HB3	1:A:270:LYS:HE3	1.97	0.47
1:B:99:LYS:HD2	1:B:99:LYS:HA	1.59	0.47
1:B:25:GLU:HG3	1:B:61:LEU:HD11	1.97	0.46
1:B:29:LYS:HD2	1:B:57:GLN:NE2	2.30	0.46
1:B:104:LEU:HD11	1:B:190:THR:HG21	1.97	0.46
1:B:177:TYR:C	1:B:177:TYR:CD2	2.92	0.46
1:A:1:MET:HG3	1:A:250:ASP:HB2	1.97	0.46
1:A:390:PRO:O	1:A:393:ILE:HG13	2.15	0.46
1:B:108:ILE:HG22	1:B:200:LEU:CD1	2.45	0.46
1:B:159:PHE:C	1:B:159:PHE:CD2	2.93	0.46
1:A:175:VAL:O	1:A:179:LYS:HB2	2.14	0.46
1:B:201:ILE:HA	1:B:204:MET:HG2	1.98	0.46
1:A:340:MET:HA	1:A:343:MET:HE3	1.97	0.46
1:B:342:LYS:C	1:B:344:GLY:N	2.72	0.46
1:B:367:ILE:HD11	1:B:372:LEU:HB2	1.97	0.46
1:A:333:MET:HE1	1:A:415:LEU:HD21	1.98	0.46
1:B:122:ARG:HD2	1:B:122:ARG:O	2.15	0.46
1:A:319:ASP:O	1:A:322:ARG:HB2	2.15	0.45
1:B:430:ASN:O	1:B:432:ARG:N	2.48	0.45
1:A:373:LYS:O	1:A:377:VAL:HG23	2.17	0.45
1:B:3:LEU:HD22	1:B:289:PHE:HZ	1.81	0.45
1:B:425:LEU:O	1:B:429:MET:HG2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:3:LEU:HD22	1:A:289:PHE:CZ	2.52	0.45
1:A:99:LYS:CB	1:A:100:PRO:CD	2.94	0.45
1:A:164:GLU:OE2	1:A:169:LYS:HE2	2.16	0.45
1:B:115:THR:HG22	1:B:119:LYS:HE3	1.99	0.45
1:B:374:LYS:O	1:B:378:ILE:HG13	2.17	0.45
1:B:145:HIS:HE1	2:B:445:SO4:O4	1.99	0.44
1:B:389:ASN:HB3	1:B:391:GLU:OE2	2.17	0.44
1:A:425:LEU:HD22	1:A:429:MET:CE	2.48	0.44
1:A:415:LEU:HD23	1:A:415:LEU:C	2.43	0.44
1:B:171:ALA:O	1:B:175:VAL:HG23	2.18	0.44
1:A:65:PRO:HG3	1:A:71:LYS:HA	1.99	0.44
1:A:99:LYS:HG3	1:A:100:PRO:CD	2.42	0.44
1:A:1:MET:HA	1:A:251:GLY:H	1.83	0.44
1:B:62:GLU:O	1:B:62:GLU:HG3	2.19	0.43
1:B:72:LYS:O	1:B:76:ILE:HG12	2.19	0.43
1:A:4:ASP:CG	1:A:5:ASN:H	2.24	0.43
1:A:1:MET:HA	1:A:251:GLY:N	2.33	0.43
1:A:65:PRO:HA	1:A:66:PRO:HD2	1.86	0.43
1:A:72:LYS:O	1:A:76:ILE:HG12	2.18	0.43
1:B:313:VAL:HG13	1:B:314:GLU:H	1.82	0.43
1:A:2:VAL:N	1:A:250:ASP:HA	2.29	0.43
1:B:17:ILE:HD13	1:B:75:ILE:HG21	2.00	0.43
1:A:243:SER:CB	1:A:270:LYS:HE3	2.49	0.43
1:B:279:ASP:HB2	5:B:545:HOH:O	2.18	0.43
1:A:159:PHE:C	1:A:159:PHE:CD2	2.96	0.43
1:A:164:GLU:HG3	1:A:170:LEU:HG	2.00	0.43
1:B:98:GLU:C	1:B:99:LYS:O	2.58	0.43
1:B:125:GLN:NE2	1:B:154:TYR:O	2.51	0.43
1:B:367:ILE:HD11	1:B:372:LEU:CA	2.49	0.43
1:A:274:THR:O	1:A:274:THR:CG2	2.67	0.43
1:A:425:LEU:HD22	1:A:429:MET:HE2	2.00	0.43
1:A:343:MET:HG2	1:A:343:MET:O	2.19	0.42
1:A:221:ASP:O	1:A:224:ILE:HG12	2.18	0.42
1:A:415:LEU:HD23	1:A:415:LEU:O	2.19	0.42
1:A:164:GLU:OE1	1:A:164:GLU:CA	2.64	0.42
1:B:177:TYR:OH	1:B:181:LYS:HE2	2.19	0.42
1:B:195:LYS:HB3	1:B:195:LYS:HE2	1.76	0.42
1:B:344:GLY:HA2	1:B:345:PRO:C	2.42	0.42
1:B:389:ASN:O	1:B:392:ILE:HG12	2.20	0.42
1:B:408:SER:C	1:B:410:LYS:N	2.75	0.42
1:A:183:VAL:HG21	1:A:186:ILE:HG12	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:278:ILE:O	1:A:278:ILE:HG13	2.19	0.42
1:B:337:LEU:HB3	1:B:372:LEU:CD1	2.49	0.42
1:A:389:ASN:HB3	1:A:391:GLU:OE2	2.20	0.42
1:B:50:LEU:HD23	1:B:50:LEU:HA	1.84	0.42
1:B:265:THR:CG2	1:B:267:ALA:H	2.33	0.42
1:B:313:VAL:HG13	1:B:314:GLU:N	2.35	0.42
1:A:168:ILE:HG21	1:A:210:VAL:HG21	2.02	0.41
1:A:171:ALA:O	1:A:175:VAL:HG23	2.20	0.41
1:A:337:LEU:HB3	1:A:372:LEU:HD11	2.02	0.41
1:B:337:LEU:O	1:B:338:GLU:C	2.61	0.41
1:A:198:LYS:HA	1:A:201:ILE:HD11	2.01	0.41
1:A:338:GLU:O	1:A:342:LYS:HD3	2.20	0.41
1:B:36:GLN:O	1:B:40:ILE:HG12	2.20	0.41
1:B:402:ALA:HB1	1:B:407:THR:O	2.20	0.41
1:B:71:LYS:HG3	2:B:448:SO4:O1	2.20	0.41
1:A:372:LEU:HD23	1:A:372:LEU:C	2.46	0.41
1:A:221:ASP:OD1	1:A:221:ASP:C	2.63	0.41
1:A:398:ILE:HG23	1:A:412:VAL:HG11	2.02	0.41
1:A:139:TRP:HH2	1:A:194:HIS:ND1	2.18	0.41
1:A:409:THR:O	1:A:413:LYS:HG3	2.20	0.41
1:A:99:LYS:HB3	1:A:100:PRO:HD3	2.02	0.41
1:B:204:MET:HG3	1:B:205:LYS:N	2.35	0.41
1:A:2:VAL:C	1:A:4:ASP:N	2.78	0.41
1:A:332:ASP:O	1:A:336:GLN:HG2	2.21	0.41
1:A:337:LEU:HB3	1:A:372:LEU:CD1	2.50	0.41
1:B:63:GLU:OE1	1:B:74:HIS:CE1	2.69	0.41
1:B:286:PRO:HB2	1:B:287:PRO:HD3	2.03	0.41
1:B:274:THR:HG22	1:B:280:ASP:CB	2.51	0.41
1:A:8:LYS:HE3	1:A:8:LYS:HB3	1.85	0.40
1:A:99:LYS:CB	1:A:100:PRO:HD3	2.51	0.40
1:A:334:TYR:CD1	1:A:379:MET:HE3	2.56	0.40
1:A:2:VAL:O	1:A:2:VAL:HG13	2.20	0.40
1:A:98:GLU:O	1:A:99:LYS:O	2.39	0.40
1:A:140:ARG:HA	1:A:141:PRO:HD2	1.87	0.40
1:A:309:LEU:HD12	1:A:309:LEU:HA	1.85	0.40
1:B:40:ILE:HD11	1:B:46:VAL:HG23	2.02	0.40
1:A:379:MET:HG3	1:A:415:LEU:HD11	2.03	0.40
1:B:321:GLU:O	1:B:325:ARG:HG3	2.22	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	412/443 (93%)	383 (93%)	23 (6%)	6 (2%)	8	16
1	B	408/443 (92%)	380 (93%)	23 (6%)	5 (1%)	10	20
All	All	820/886 (93%)	763 (93%)	46 (6%)	11 (1%)	9	18

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	99	LYS
1	B	4	ASP
1	B	100	PRO
1	B	394	ASN
1	A	430	ASN
1	B	431	LYS
1	A	4	ASP
1	B	421	GLN
1	A	100	PRO
1	A	2	VAL
1	A	162	PRO

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	332/378 (88%)	308 (93%)	24 (7%)	13	28
1	B	333/378 (88%)	305 (92%)	28 (8%)	10	22

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	665/756 (88%)	613 (92%)	52 (8%)	11	24

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	ILE
1	A	39	LEU
1	A	61	LEU
1	A	71	LYS
1	A	99	LYS
1	A	108	ILE
1	A	133	VAL
1	A	140	ARG
1	A	153	ARG
1	A	168	ILE
1	A	170	LEU
1	A	194	HIS
1	A	201	ILE
1	A	204	MET
1	A	212	HIS
1	A	233	LEU
1	A	265	THR
1	A	274	THR
1	A	330	LEU
1	A	331	LYS
1	A	341	ARG
1	A	391	GLU
1	A	425	LEU
1	A	430	ASN
1	B	2	VAL
1	B	17	ILE
1	B	31	LEU
1	B	32	VAL
1	B	39	LEU
1	B	61	LEU
1	B	108	ILE
1	B	122	ARG
1	B	133	VAL
1	B	165	LYS
1	B	169	LYS
1	B	170	LEU
1	B	201	ILE

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Mol	Chain	Res	Type
1	B	204	MET
1	B	217	ILE
1	B	233	LEU
1	B	265	THR
1	B	274	THR
1	B	309	LEU
1	B	314	GLU
1	B	330	LEU
1	B	331	LYS
1	B	338	GLU
1	B	341	ARG
1	B	367	ILE
1	B	391	GLU
1	B	395	TYR
1	B	425	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	GLN
1	A	45	ASN
1	A	74	HIS
1	A	209	ASN
1	A	226	GLN
1	A	300	GLN
1	A	418	GLN
1	B	45	ASN
1	B	57	GLN
1	B	74	HIS
1	B	145	HIS
1	B	163	GLN
1	B	226	GLN
1	B	336	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

15 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	SO4	B	447	-	4,4,4	0.28	0	6,6,6	0.32	0
2	SO4	A	447	-	4,4,4	0.26	0	6,6,6	0.30	0
3	ACT	A	450	-	3,3,3	0.75	0	3,3,3	1.58	1 (33%)
2	SO4	A	449	-	4,4,4	0.23	0	6,6,6	0.15	0
2	SO4	B	448	-	4,4,4	0.22	0	6,6,6	0.25	0
2	SO4	A	445	-	4,4,4	0.25	0	6,6,6	0.14	0
2	SO4	B	445	-	4,4,4	0.30	0	6,6,6	0.30	0
4	GDP	A	501	-	29,30,30	1.19	3 (10%)	45,47,47	1.71	7 (15%)
4	GDP	B	501	-	29,30,30	1.10	3 (10%)	45,47,47	1.77	9 (20%)
2	SO4	A	448	-	4,4,4	0.30	0	6,6,6	0.14	0
3	ACT	B	449	-	3,3,3	0.78	0	3,3,3	1.38	0
2	SO4	B	446	-	4,4,4	0.25	0	6,6,6	0.05	0
2	SO4	A	446	-	4,4,4	0.26	0	6,6,6	0.20	0
2	SO4	B	444	-	4,4,4	0.30	0	6,6,6	0.15	0
2	SO4	A	444	-	4,4,4	0.25	0	6,6,6	0.10	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
4	GDP	A	501	-	-	0/16/32/32	0/3/3/3
4	GDP	B	501	-	-	0/16/32/32	0/3/3/3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	501	GDP	C5-C4	3.10	1.47	1.38
4	A	501	GDP	C5-C4	3.05	1.47	1.38
4	A	501	GDP	C6-N1	-2.42	1.34	1.38
4	A	501	GDP	C4-N9	-2.33	1.32	1.38
4	B	501	GDP	C5-N7	-2.29	1.34	1.39
4	B	501	GDP	C6-N1	-2.16	1.34	1.38

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	501	GDP	C5-C4-N3	-5.55	119.56	128.39
4	A	501	GDP	C5-C4-N3	-5.11	120.25	128.39
4	B	501	GDP	C2-N3-C4	4.61	120.25	112.30
4	A	501	GDP	C2-N3-C4	4.56	120.15	112.30
4	B	501	GDP	N9-C4-N3	4.43	134.81	125.95
4	A	501	GDP	C6-C5-N7	4.01	137.59	130.29
4	A	501	GDP	N9-C4-N3	3.54	133.02	125.95
4	B	501	GDP	C6-C5-N7	3.22	136.15	130.29
4	B	501	GDP	O6-C6-C5	-2.83	119.05	126.53
4	A	501	GDP	C4-C5-N7	-2.79	106.25	110.67
4	B	501	GDP	O2A-PA-O3A	2.50	114.03	107.27
4	A	501	GDP	O3B-PB-O3A	2.39	112.65	104.64
4	B	501	GDP	C4-C5-N7	-2.12	107.31	110.67
3	A	450	ACT	OXT-C-CH3	2.10	123.88	115.05
4	B	501	GDP	C5-C6-N1	2.03	118.43	113.25
4	A	501	GDP	O3'-C3'-C4'	-2.02	105.28	111.08
4	B	501	GDP	O3'-C3'-C2'	-2.01	105.39	111.82

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 3 short contacts:

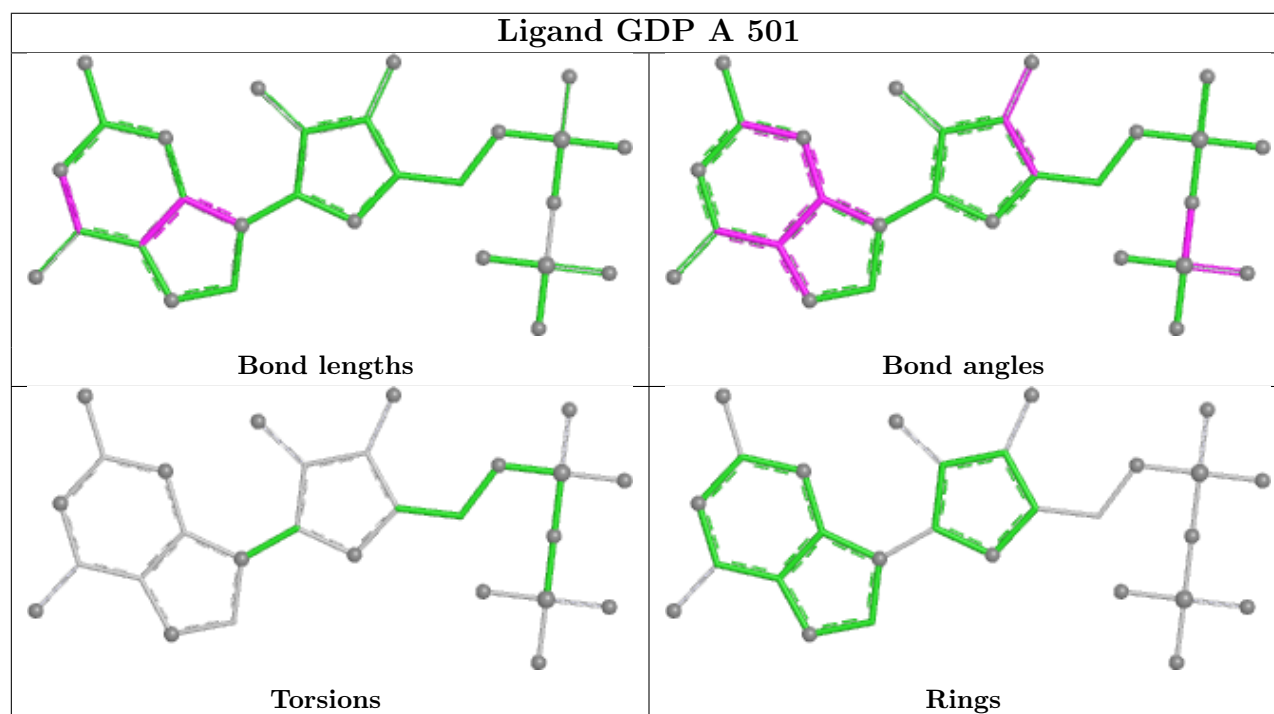
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	448	SO4	1	0

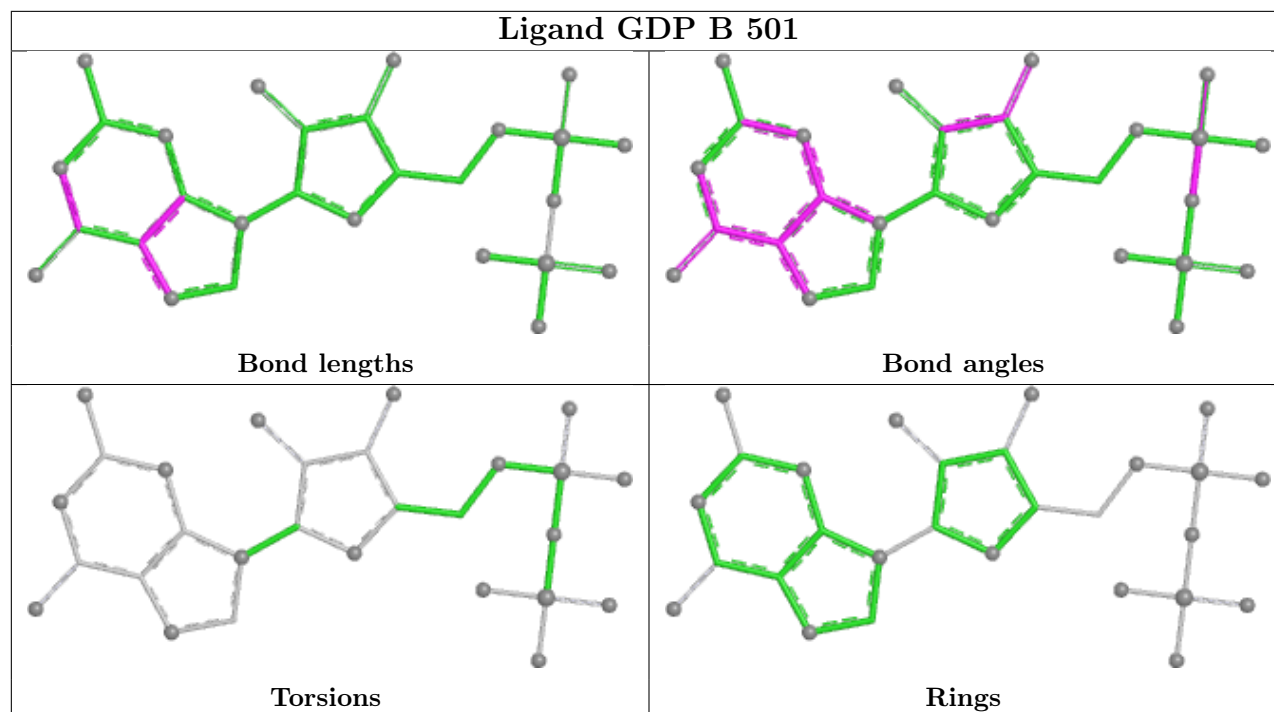
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	445	SO4	1	0
2	A	448	SO4	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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	416/443 (93%)	0.42	35 (8%) 17 15	35, 59, 112, 141	0
1	B	412/443 (93%)	0.52	33 (8%) 18 16	33, 64, 111, 132	0
All	All	828/886 (93%)	0.47	68 (8%) 17 15	33, 61, 112, 141	0

All (68) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	345	PRO	5.8
1	B	434	LEU	5.8
1	B	430	ASN	5.7
1	B	345	PRO	5.3
1	A	435	SER	4.9
1	B	141	PRO	4.2
1	A	365	ILE	4.2
1	B	432	ARG	3.8
1	B	395	TYR	3.7
1	B	202	GLU	3.7
1	B	306	PHE	3.6
1	A	306	PHE	3.4
1	B	334	TYR	3.3
1	B	340	MET	3.2
1	A	373	LYS	3.2
1	A	434	LEU	3.1
1	B	338	GLU	3.0
1	A	192	GLY	3.0
1	A	389	ASN	3.0
1	B	367	ILE	2.9
1	B	324	LEU	2.8
1	A	139	TRP	2.8
1	A	426	PHE	2.8
1	A	372	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	1	MET	2.8
1	A	162	PRO	2.8
1	A	425	LEU	2.8
1	A	337	LEU	2.7
1	A	433	GLN	2.7
1	A	142	GLY	2.7
1	B	368	GLY	2.7
1	A	367	ILE	2.7
1	A	71	LYS	2.6
1	B	343	MET	2.6
1	B	433	GLN	2.6
1	A	368	GLY	2.6
1	A	163	GLN	2.6
1	A	322	ARG	2.5
1	A	141	PRO	2.5
1	B	391	GLU	2.5
1	B	178	PHE	2.5
1	B	373	LYS	2.5
1	B	304	GLU	2.5
1	A	169	LYS	2.4
1	A	252	SER	2.4
1	A	194	HIS	2.4
1	A	429	MET	2.4
1	B	162	PRO	2.4
1	A	395	TYR	2.3
1	B	97	LYS	2.3
1	B	204	MET	2.3
1	A	432	ARG	2.3
1	B	405	SER	2.3
1	B	429	MET	2.3
1	A	327	LYS	2.2
1	B	424	LYS	2.2
1	A	338	GLU	2.1
1	B	414	GLU	2.1
1	A	342	LYS	2.1
1	A	430	ASN	2.1
1	B	192	GLY	2.1
1	B	180	SER	2.1
1	A	344	GLY	2.1
1	B	408	SER	2.1
1	A	343	MET	2.0
1	B	371	ARG	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	29	LYS	2.0
1	B	194	HIS	2.0

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 oligosaccharides in this entry.

6.4 Ligands [i](#)

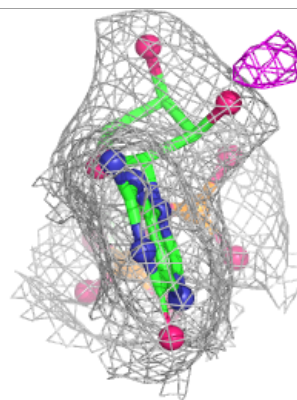
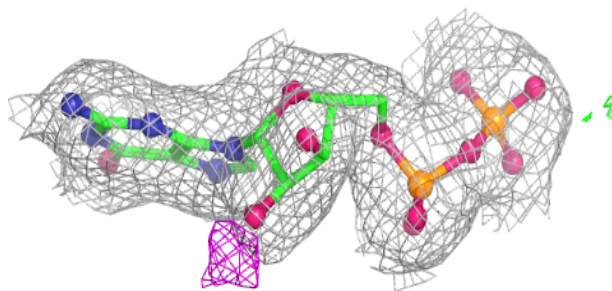
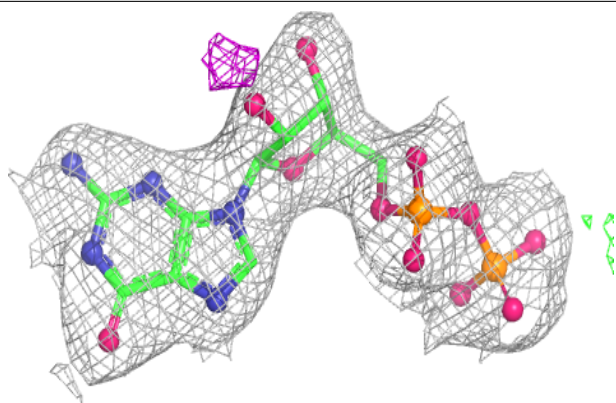
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	B	446	5/5	0.54	0.11	107,109,125,135	0
2	SO4	A	449	5/5	0.67	0.12	88,99,117,124	0
3	ACT	A	450	4/4	0.67	0.23	59,78,84,87	0
2	SO4	B	448	5/5	0.75	0.10	105,106,116,132	0
3	ACT	B	449	4/4	0.78	0.20	58,78,79,80	0
2	SO4	A	446	5/5	0.80	0.10	90,92,109,123	0
2	SO4	B	447	5/5	0.81	0.12	79,85,89,101	0
2	SO4	A	448	5/5	0.83	0.08	87,89,105,112	0
2	SO4	B	445	5/5	0.84	0.12	73,73,103,111	0
2	SO4	A	445	5/5	0.86	0.10	89,91,110,118	0
2	SO4	B	444	5/5	0.87	0.09	87,89,112,122	0
2	SO4	A	444	5/5	0.88	0.09	97,101,106,122	0
2	SO4	A	447	5/5	0.94	0.12	65,76,87,90	0
4	GDP	A	501	28/28	0.98	0.06	37,45,49,51	0
4	GDP	B	501	28/28	0.98	0.06	35,41,51,64	0

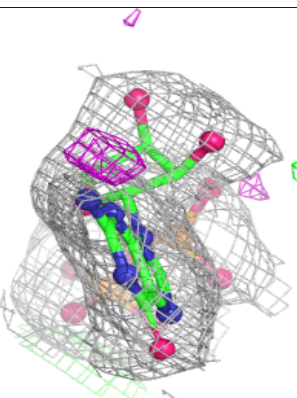
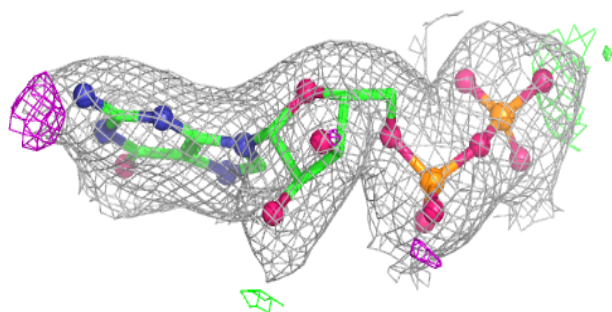
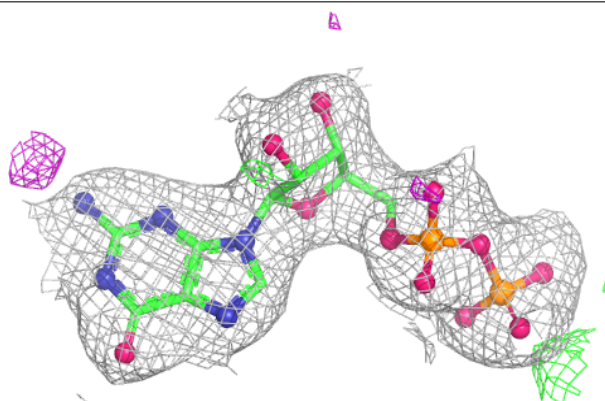
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around GDP A 501:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around GDP B 501:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.