Full wwPDB X-ray Structure Validation Report (i)

Nov 16, 2023 – 10:26 AM JST

PDB ID	:	6LFZ
Title	:	Crystal structure of SbCGTb in complex with UDPG
Authors	:	Gao, H.M.; Yun, C.H.
Deposited on	:	2019-12-04
Resolution	:	2.87 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 (i) 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 (i)) were used in the production of this report:

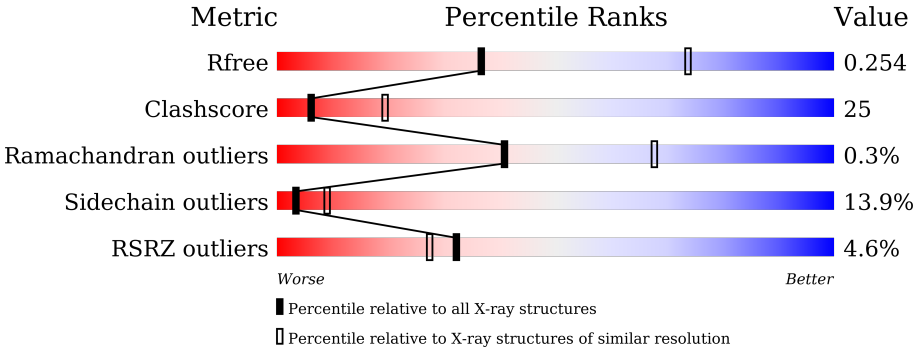
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

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}	130704	3168 (2.90-2.82)
Clashscore	141614	3438 (2.90-2.82)
Ramachandran outliers	138981	3348 (2.90-2.82)
Sidechain outliers	138945	3351 (2.90-2.82)
RSRZ outliers	127900	3103 (2.90-2.82)

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 <=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	459	3% 60% 31% • 5%
1	B	459	6% 55% 34% 7% •

2 Entry composition (i)

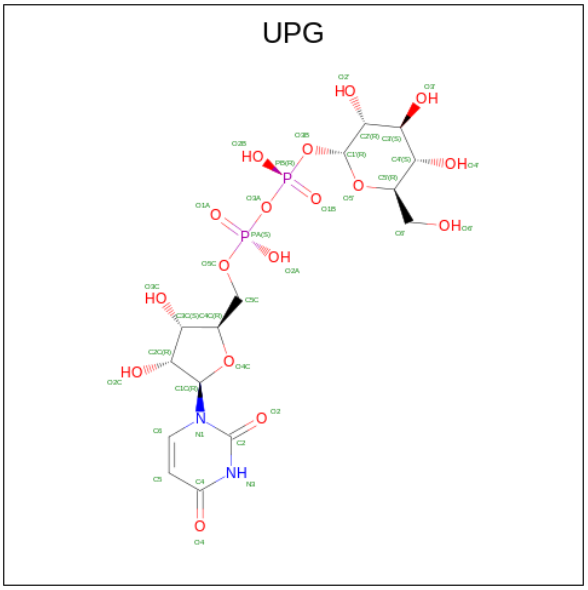
There are 3 unique types of molecules in this entry. The entry contains 6852 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 SbCGTb.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	435	3333	2143	579	603	8	0	0	0
1	B	439	3394	2180	596	609	9	0	0	0

- Molecule 2 is URIDINE-5'-DIPHOSPHATE-GLUCOSE (three-letter code: UPG) (formula: $C_{15}H_{24}N_2O_{17}P_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	36	15	2	17	2	0	0
2	B	1	36	15	2	17	2	0	0

- Molecule 3 is water.

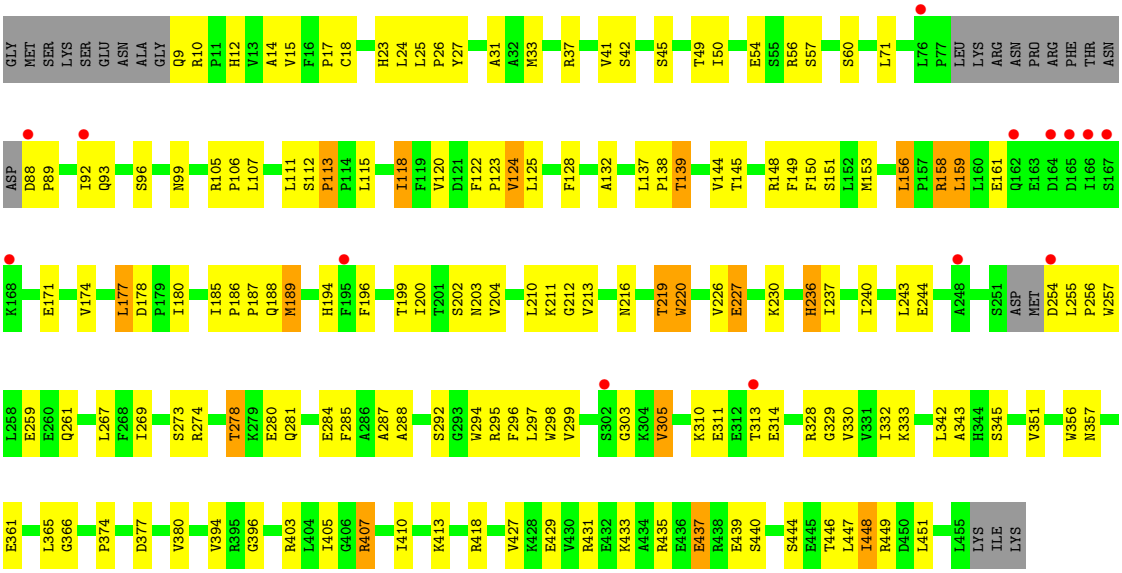
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	34	34	34	0	0
			Total	O		
3	B	19	19	19	0	0

3 Residue-property plots (i)

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.

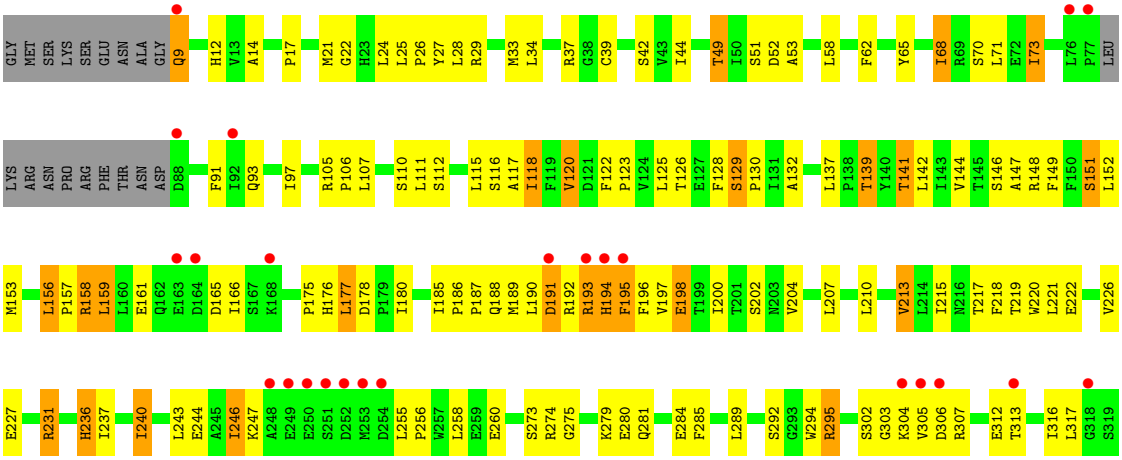
- Molecule 1: SbCGTb

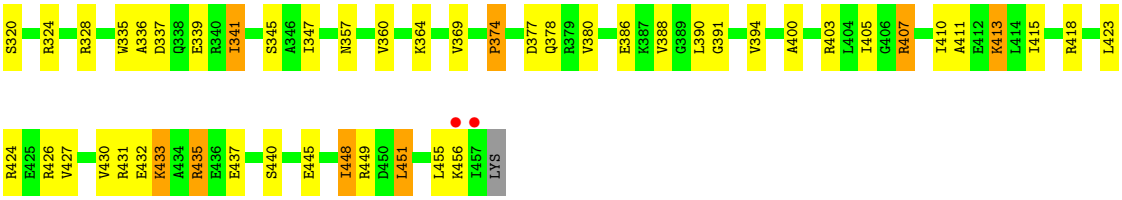
Chain A: 3% 60% 31% • 5%



- Molecule 1: SbCGTb

Chain B: 6% 55% 34% 7% •




S320
R324
R328
W335
A336
D337
Q338
E339
R340
I341
S345
A346
I347
N357
V360
K364
V369
P374
D377
Q378
R379
V380
E386
K387
V388
G389
L390
G391
V394
A400
R403
L404
I405
G406
R407
I410
A411
E412
K413
L414
I415
R418
L423
R424
E425
R426
V427
V430
R431
E432
K433
A434
R435
E436
E437
S440
E445
I448
R449
D450
L451
L455
K456
I457
LYS

4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α, β, γ	107.29Å 216.95Å 104.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.16 – 2.87 36.16 – 2.87	Depositor EDS
% Data completeness (in resolution range)	99.3 (36.16-2.87) 99.3 (36.16-2.87)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.68 (at 2.85Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
R, R_{free}	0.218 , 0.254 0.218 , 0.254	Depositor DCC
R_{free} test set	1337 reflections (4.73%)	wwPDB-VP
Wilson B-factor (Å^2)	44.2	Xtriage
Anisotropy	0.142	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 45.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.90	EDS
Total number of atoms	6852	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 5.06% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 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: UPG

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	0.81	0/3412	0.79	0/4636
1	B	0.70	0/3475	0.69	0/4720
All	All	0.76	0/6887	0.74	0/9356

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	3333	0	3267	136	0
1	B	3394	0	3358	197	0
2	A	36	0	22	1	0
2	B	36	0	22	4	0
3	A	34	0	0	6	0
3	B	19	0	0	5	0
All	All	6852	0	6669	331	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 25.

All (331) 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:112:SER:OG	1:A:113:PRO:HD3	1.17	1.27
1:B:193:ARG:HG3	1:B:193:ARG:HH11	1.04	1.15
1:B:295:ARG:HH21	1:B:295:ARG:CG	1.68	1.06
1:B:166:ILE:CB	1:B:190:LEU:HD22	1.88	1.04
1:B:117:ALA:HB2	1:B:455:LEU:HD21	1.42	1.01
1:B:295:ARG:HH21	1:B:295:ARG:HG2	0.87	1.00
1:A:112:SER:OG	1:A:113:PRO:CD	2.09	1.00
1:B:374:PRO:HD2	1:B:378:GLN:OE1	1.59	1.00
1:B:295:ARG:HG2	1:B:295:ARG:NH2	1.70	0.99
1:B:193:ARG:HG3	1:B:193:ARG:NH1	1.74	0.97
1:A:37:ARG:HG3	1:A:37:ARG:HH11	1.31	0.93
1:B:186:PRO:HD2	1:B:189:MET:HE3	1.51	0.92
1:A:185:ILE:HG23	1:A:189:MET:HE3	1.53	0.91
1:A:437:GLU:HG2	1:A:444:SER:OG	1.74	0.87
1:A:274:ARG:HD2	1:A:305:VAL:HA	1.55	0.87
1:B:149:PHE:CD1	1:B:377:ASP:HB2	2.10	0.87
1:B:213:VAL:HG22	1:B:237:ILE:HD12	1.55	0.86
1:A:49:THR:HG21	1:A:54:GLU:HB2	1.55	0.86
1:A:105:ARG:HB2	1:A:106:PRO:HD3	1.57	0.86
1:B:191:ASP:OD1	1:B:194:HIS:HB2	1.76	0.85
1:B:407:ARG:HG2	1:B:407:ARG:HH11	1.40	0.85
1:A:357:ASN:O	1:A:361:GLU:HG3	1.77	0.84
1:B:29:ARG:HG2	1:B:244:GLU:O	1.78	0.84
1:B:204:VAL:O	1:B:207:LEU:HB2	1.77	0.84
1:A:112:SER:CB	1:A:113:PRO:HD3	2.07	0.83
1:B:407:ARG:HG2	1:B:407:ARG:NH1	1.94	0.83
1:B:33:MET:HE1	1:B:243:LEU:HD23	1.60	0.82
1:B:73:ILE:HD11	1:B:107:LEU:HD22	1.62	0.82
1:A:112:SER:HG	1:A:113:PRO:HD3	1.04	0.81
1:B:394:VAL:CG2	1:B:413:LYS:HE3	2.10	0.81
1:B:220:TRP:CE2	1:B:433:LYS:HG2	2.16	0.81
1:A:356:TRP:HD1	3:A:622:HOH:O	1.64	0.79
1:B:200:ILE:O	1:B:204:VAL:HG23	1.82	0.79
1:B:49:THR:CG2	1:B:51:SER:O	2.30	0.79
1:A:174:VAL:HB	1:A:177:LEU:HD11	1.64	0.78
1:B:302:SER:CB	1:B:312:GLU:OE2	2.32	0.78
1:A:18:CYS:SG	1:A:124:VAL:HG22	2.23	0.77
1:B:407:ARG:HH11	1:B:407:ARG:CG	1.97	0.77
1:B:426:ARG:HD2	3:B:612:HOH:O	1.85	0.77
1:B:149:PHE:CG	1:B:377:ASP:HB2	2.20	0.76
1:B:426:ARG:O	1:B:430:VAL:HG23	1.84	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:278:THR:HG22	1:A:281:GLN:H	1.50	0.74
1:B:316:ILE:HG22	1:B:317:LEU:HD23	1.70	0.74
1:B:374:PRO:CB	1:B:400:ALA:HB2	2.18	0.73
1:B:149:PHE:HZ	1:B:189:MET:HE2	1.52	0.73
1:B:426:ARG:CD	3:B:612:HOH:O	2.37	0.73
1:B:159:LEU:HD23	1:B:159:LEU:O	1.89	0.72
1:B:148:ARG:HG2	1:B:380:VAL:CG1	2.20	0.72
1:B:105:ARG:HB2	1:B:106:PRO:HD3	1.69	0.72
1:B:144:VAL:HG23	2:B:500:UPG:O6'	1.90	0.72
1:B:148:ARG:NH1	1:B:178:ASP:HB3	2.05	0.72
1:B:394:VAL:HG22	1:B:413:LYS:HE3	1.70	0.72
1:A:255:LEU:N	1:A:256:PRO:HD2	2.05	0.71
1:B:374:PRO:HB2	1:B:400:ALA:HB2	1.70	0.71
1:B:451:LEU:O	1:B:455:LEU:HB2	1.91	0.71
1:A:88:ASP:N	1:A:89:PRO:CD	2.53	0.71
1:A:33:MET:HE1	1:A:243:LEU:HD23	1.73	0.71
1:B:149:PHE:CZ	1:B:189:MET:HE2	2.26	0.71
1:A:18:CYS:SG	1:A:124:VAL:CG2	2.79	0.70
1:B:33:MET:HE1	1:B:243:LEU:CD2	2.21	0.70
1:A:227:GLU:CG	3:A:602:HOH:O	2.40	0.70
1:A:216:ASN:HB3	3:A:622:HOH:O	1.92	0.70
1:B:159:LEU:O	1:B:159:LEU:CD2	2.40	0.69
1:A:148:ARG:HD3	1:A:380:VAL:HG13	1.72	0.69
1:A:149:PHE:CG	1:A:377:ASP:HB2	2.27	0.69
1:A:213:VAL:HG23	1:A:237:ILE:HD12	1.74	0.69
1:B:33:MET:HE2	1:B:243:LEU:HD22	1.75	0.69
1:B:280:GLU:O	1:B:284:GLU:HG2	1.93	0.69
1:B:33:MET:CE	1:B:243:LEU:CD2	2.71	0.68
1:B:33:MET:CE	1:B:243:LEU:HD22	2.24	0.68
1:B:149:PHE:CZ	1:B:189:MET:CE	2.77	0.68
1:B:451:LEU:HD12	1:B:451:LEU:C	2.13	0.68
1:A:37:ARG:HG3	1:A:37:ARG:NH1	2.03	0.68
1:A:159:LEU:O	1:A:159:LEU:HD22	1.95	0.67
1:B:49:THR:HG22	1:B:51:SER:O	1.94	0.67
1:A:356:TRP:CD1	3:A:622:HOH:O	2.42	0.67
1:A:88:ASP:N	1:A:89:PRO:HD3	2.10	0.66
1:A:227:GLU:HG2	3:A:602:HOH:O	1.94	0.66
1:A:303:GLY:HA3	1:A:310:LYS:HB2	1.78	0.66
1:A:278:THR:CG2	1:A:281:GLN:H	2.08	0.66
1:B:9:GLN:N	1:B:9:GLN:CD	2.48	0.66
1:B:117:ALA:CB	1:B:455:LEU:HD21	2.23	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:12:HIS:ND1	1:B:115:LEU:HD23	2.10	0.66
1:B:25:LEU:N	1:B:26:PRO:HD2	2.11	0.65
1:B:116:SER:HB2	1:B:455:LEU:HD11	1.78	0.65
1:A:49:THR:CG2	1:A:54:GLU:HB2	2.26	0.65
1:B:193:ARG:HH11	1:B:193:ARG:CG	1.91	0.64
1:B:423:LEU:O	1:B:427:VAL:HG23	1.96	0.64
1:A:138:PRO:HB3	1:A:211:LYS:HD3	1.79	0.64
1:B:53:ALA:H	1:B:307:ARG:CD	2.11	0.64
1:B:191:ASP:HB2	1:B:193:ARG:HH21	1.62	0.64
1:B:220:TRP:CD2	1:B:433:LYS:HG2	2.33	0.64
1:B:195:PHE:H	1:B:195:PHE:HD1	1.44	0.63
1:B:188:GLN:OE1	1:B:188:GLN:N	2.29	0.63
1:B:255:LEU:N	1:B:256:PRO:HD2	2.14	0.63
1:A:33:MET:HE2	1:A:243:LEU:HD22	1.80	0.63
1:A:188:GLN:CD	1:A:188:GLN:H	2.02	0.62
1:B:116:SER:HB2	1:B:455:LEU:CD1	2.30	0.62
1:B:159:LEU:CD2	1:B:159:LEU:C	2.66	0.62
1:B:189:MET:HG3	1:B:196:PHE:HE2	1.64	0.62
1:A:161:GLU:CB	3:A:618:HOH:O	2.48	0.62
1:B:122:PHE:N	1:B:123:PRO:HD2	2.15	0.62
1:A:33:MET:HE1	1:A:243:LEU:CD2	2.29	0.62
1:A:257:TRP:O	1:A:261:GLN:HG2	2.00	0.62
1:B:193:ARG:NH1	1:B:193:ARG:H	1.98	0.62
1:A:200:ILE:O	1:A:204:VAL:HG23	2.00	0.61
1:B:451:LEU:O	1:B:451:LEU:HD12	2.01	0.61
1:B:191:ASP:CB	1:B:193:ARG:NH2	2.63	0.61
1:A:186:PRO:HD2	1:A:189:MET:HE2	1.82	0.60
1:B:213:VAL:CG2	1:B:237:ILE:HD12	2.30	0.60
1:B:118:ILE:HD13	1:B:120:VAL:HG12	1.84	0.60
1:B:411:ALA:O	1:B:415:ILE:HG13	2.02	0.60
1:A:199:THR:O	1:A:203:ASN:ND2	2.35	0.59
1:B:73:ILE:CD1	1:B:107:LEU:HD22	2.31	0.59
1:A:311:GLU:OE2	1:A:311:GLU:HA	2.03	0.59
1:B:153:MET:HG2	1:B:200:ILE:HD12	1.85	0.59
1:B:193:ARG:NH1	1:B:193:ARG:CG	2.56	0.59
1:B:156:LEU:HB3	1:B:157:PRO:HD3	1.85	0.58
1:B:149:PHE:CZ	1:B:377:ASP:HB3	2.38	0.58
1:A:50:ILE:HD11	1:A:93:GLN:HB2	1.83	0.58
1:A:33:MET:CE	1:A:243:LEU:CD2	2.81	0.58
1:B:195:PHE:CD1	1:B:195:PHE:N	2.58	0.58
1:B:360:VAL:O	1:B:364:LYS:HB3	2.03	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:279:LYS:HD3	1:B:316:ILE:HA	1.87	0.57
1:A:149:PHE:CD1	1:A:377:ASP:HB2	2.39	0.57
1:A:396:GLY:O	1:A:403:ARG:NH2	2.38	0.57
1:B:120:VAL:CG2	1:B:125:LEU:HD23	2.35	0.57
1:A:14:ALA:HB3	1:A:118:ILE:HG13	1.86	0.57
1:A:178:ASP:N	1:A:178:ASP:OD1	2.36	0.57
1:B:128:PHE:CD1	1:B:128:PHE:O	2.56	0.57
1:A:285:PHE:O	1:A:288:ALA:HB3	2.05	0.57
1:B:148:ARG:HH11	1:B:178:ASP:HB3	1.68	0.57
1:A:137:LEU:O	1:A:139:THR:HG22	2.04	0.56
1:B:295:ARG:CG	1:B:295:ARG:NH2	2.38	0.56
1:A:120:VAL:HG22	1:A:125:LEU:HD23	1.85	0.56
1:B:195:PHE:O	1:B:198:GLU:N	2.38	0.56
1:A:88:ASP:O	1:A:92:ILE:HG13	2.06	0.56
1:A:25:LEU:N	1:A:26:PRO:HD2	2.20	0.56
1:A:437:GLU:HG3	1:A:444:SER:H	1.70	0.56
1:B:141:THR:OG1	1:B:210:LEU:HD13	2.06	0.56
1:A:297:LEU:HD11	1:A:332:ILE:HG13	1.87	0.55
1:A:437:GLU:CG	1:A:444:SER:OG	2.53	0.55
1:A:254:ASP:OD2	1:B:435:ARG:HD3	2.06	0.55
1:A:255:LEU:N	1:A:256:PRO:CD	2.69	0.55
1:A:150:PHE:HA	1:A:153:MET:HE2	1.89	0.55
1:A:194:HIS:HD2	1:A:196:PHE:H	1.54	0.55
1:A:149:PHE:CE2	1:A:377:ASP:HB3	2.41	0.54
1:A:122:PHE:N	1:A:123:PRO:HD2	2.23	0.54
1:A:213:VAL:HG23	1:A:213:VAL:O	2.07	0.54
1:B:149:PHE:CE2	1:B:377:ASP:HB3	2.42	0.54
1:B:149:PHE:CE1	1:B:377:ASP:CB	2.90	0.54
1:B:258:LEU:HD11	1:B:341:ILE:HG23	1.90	0.54
1:B:149:PHE:CD1	1:B:377:ASP:CB	2.87	0.54
1:B:149:PHE:CZ	1:B:189:MET:HE1	2.44	0.53
1:A:37:ARG:NH1	1:A:37:ARG:CG	2.71	0.53
1:B:28:LEU:HB3	1:B:246:ILE:HD11	1.90	0.53
1:B:122:PHE:HB3	1:B:123:PRO:HD3	1.91	0.53
1:B:126:THR:O	1:B:129:SER:HB2	2.08	0.53
1:A:17:PRO:HG3	1:A:27:TYR:CB	2.38	0.53
1:B:378:GLN:HG3	2:B:500:UPG:O3'	2.09	0.53
1:B:120:VAL:O	1:B:141:THR:HG23	2.08	0.53
1:A:278:THR:HG23	1:A:280:GLU:H	1.73	0.53
1:A:33:MET:CE	1:A:243:LEU:HD22	2.39	0.52
1:A:71:LEU:HB3	1:A:107:LEU:HD21	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:191:ASP:O	1:B:197:VAL:HG21	2.09	0.52
1:B:29:ARG:O	1:B:33:MET:HG3	2.09	0.52
1:A:429:GLU:O	1:A:433:LYS:HG2	2.09	0.52
1:B:148:ARG:HG2	1:B:380:VAL:HG12	1.90	0.52
1:B:65:TYR:OH	1:B:247:LYS:HB2	2.09	0.52
1:B:207:LEU:HB2	3:B:604:HOH:O	2.09	0.52
1:A:278:THR:HG23	1:A:280:GLU:N	2.25	0.52
1:A:314:GLU:N	1:A:314:GLU:OE1	2.43	0.52
1:A:405:ILE:HG22	1:A:410:ILE:HG13	1.91	0.52
1:A:56:ARG:HG2	1:B:449:ARG:NH2	2.24	0.52
1:B:189:MET:HG3	1:B:196:PHE:CE2	2.42	0.52
1:B:191:ASP:HB2	1:B:193:ARG:NH2	2.25	0.52
1:A:118:ILE:HD13	1:A:120:VAL:HG12	1.92	0.52
1:A:303:GLY:HA3	1:A:310:LYS:CB	2.40	0.52
1:A:292:SER:O	1:A:418:ARG:NH1	2.43	0.52
1:B:116:SER:CB	1:B:455:LEU:CD1	2.88	0.51
1:A:37:ARG:NH2	1:A:244:GLU:OE2	2.42	0.51
1:B:129:SER:HB3	1:B:130:PRO:HD3	1.92	0.51
1:B:191:ASP:CB	1:B:193:ARG:HH21	2.24	0.51
1:A:394:VAL:CG2	1:A:413:LYS:HE2	2.41	0.51
1:B:337:ASP:O	1:B:341:ILE:HG13	2.10	0.51
1:A:227:GLU:HA	1:A:230:LYS:HE2	1.92	0.51
1:B:129:SER:N	1:B:130:PRO:CD	2.73	0.51
1:A:144:VAL:HG23	2:A:500:UPG:O6'	2.11	0.50
1:B:65:TYR:HB3	1:B:68:ILE:HG13	1.94	0.50
1:A:366:GLY:O	1:A:427:VAL:HG11	2.11	0.50
1:B:191:ASP:OD1	1:B:191:ASP:O	2.30	0.50
1:B:144:VAL:HG23	2:B:500:UPG:C6'	2.42	0.50
1:A:33:MET:CE	1:A:243:LEU:HD23	2.42	0.49
1:B:159:LEU:O	1:B:159:LEU:HD22	2.12	0.49
1:B:122:PHE:N	1:B:123:PRO:CD	2.76	0.49
1:A:159:LEU:HD22	1:A:159:LEU:C	2.32	0.49
1:B:191:ASP:OD2	1:B:193:ARG:NE	2.46	0.49
1:A:145:THR:HG21	1:A:153:MET:HE1	1.95	0.49
1:B:24:LEU:HD11	1:B:58:LEU:HD21	1.93	0.49
1:B:65:TYR:CZ	1:B:247:LYS:HB2	2.47	0.49
1:A:254:ASP:OD1	1:A:254:ASP:O	2.30	0.49
1:A:23:HIS:O	1:A:26:PRO:HG2	2.13	0.48
1:A:303:GLY:HA3	1:A:310:LYS:CG	2.43	0.48
1:A:212:GLY:HA2	1:A:236:HIS:HB2	1.95	0.48
1:B:292:SER:O	1:B:418:ARG:NH1	2.46	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:281:GLN:O	1:A:284:GLU:HB2	2.14	0.48
1:B:313:THR:O	1:B:313:THR:OG1	2.21	0.48
1:B:215:ILE:HG23	1:B:217:THR:HG22	1.94	0.48
1:A:269:ILE:HG23	1:A:351:VAL:HB	1.95	0.48
1:B:149:PHE:CE1	1:B:189:MET:HE1	2.48	0.48
1:A:50:ILE:HD11	1:A:89:PRO:O	2.14	0.48
1:B:116:SER:HB3	1:B:455:LEU:HD13	1.95	0.48
1:B:156:LEU:N	1:B:157:PRO:CD	2.77	0.48
1:B:159:LEU:C	1:B:159:LEU:HD22	2.34	0.48
1:B:305:VAL:C	1:B:306:ASP:OD1	2.53	0.48
1:A:112:SER:HG	1:A:113:PRO:CD	1.98	0.47
1:A:255:LEU:O	1:A:259:GLU:N	2.40	0.47
1:A:177:LEU:O	1:A:177:LEU:HD12	2.15	0.47
1:A:313:THR:HG23	1:A:313:THR:O	2.14	0.47
1:B:175:PRO:O	1:B:176:HIS:HB2	2.15	0.47
1:A:105:ARG:HB2	1:A:106:PRO:CD	2.38	0.47
1:B:220:TRP:CD2	1:B:433:LYS:CG	2.98	0.47
1:A:138:PRO:HB3	1:A:211:LYS:CD	2.44	0.47
1:A:149:PHE:CZ	1:A:377:ASP:HB3	2.50	0.47
1:B:147:ALA:O	1:B:151:SER:OG	2.32	0.47
1:A:12:HIS:O	1:A:115:LEU:HA	2.15	0.47
1:A:18:CYS:SG	1:A:124:VAL:HG21	2.54	0.46
1:B:448:ILE:CG2	1:B:449:ARG:N	2.78	0.46
1:B:25:LEU:N	1:B:26:PRO:CD	2.79	0.46
1:B:213:VAL:HG22	1:B:237:ILE:CD1	2.36	0.46
1:A:366:GLY:O	1:A:427:VAL:CG1	2.63	0.46
1:B:390:LEU:HD12	1:B:427:VAL:HG13	1.97	0.46
1:A:177:LEU:CD1	1:A:177:LEU:C	2.84	0.46
1:A:287:ALA:HB3	1:A:407:ARG:HG2	1.98	0.46
1:A:237:ILE:HG23	1:A:237:ILE:O	2.15	0.46
1:B:14:ALA:HB3	1:B:118:ILE:HG13	1.98	0.46
1:A:111:LEU:CD1	1:A:115:LEU:HD21	2.46	0.45
1:A:177:LEU:HD12	1:A:177:LEU:C	2.37	0.45
1:B:22:GLY:O	1:B:26:PRO:HG3	2.17	0.45
1:B:128:PHE:CE1	1:B:132:ALA:HB2	2.51	0.45
1:B:62:PHE:CE2	1:B:70:SER:HB2	2.51	0.45
1:A:267:LEU:HB2	1:A:294:TRP:CE3	2.52	0.45
1:B:186:PRO:HA	1:B:187:PRO:HD3	1.85	0.45
1:A:297:LEU:HD13	1:A:330:VAL:HB	1.98	0.45
1:B:25:LEU:HB2	1:B:26:PRO:CD	2.47	0.45
1:B:445:GLU:O	1:B:448:ILE:HG22	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:24:LEU:HD22	1:A:54:GLU:HB3	1.99	0.44
1:A:296:PHE:CZ	1:A:329:GLY:HA3	2.52	0.44
1:A:343:ALA:HA	1:A:365:LEU:HD22	1.99	0.44
1:B:149:PHE:O	1:B:152:LEU:HB3	2.17	0.44
1:B:22:GLY:O	1:B:357:ASN:ND2	2.48	0.44
1:B:218:PHE:CB	1:B:437:GLU:HG3	2.47	0.44
1:B:191:ASP:OD1	1:B:191:ASP:C	2.55	0.44
1:A:219:THR:CG2	1:A:226:VAL:HG11	2.47	0.44
1:B:307:ARG:H	1:B:307:ARG:HG2	1.50	0.44
1:B:27:TYR:CZ	1:B:142:LEU:HB3	2.52	0.44
1:B:156:LEU:N	1:B:157:PRO:HD2	2.32	0.44
1:A:148:ARG:NH2	1:A:180:ILE:HG12	2.33	0.44
1:A:274:ARG:NE	1:A:305:VAL:HG12	2.33	0.44
1:B:22:GLY:O	1:B:26:PRO:CG	2.65	0.44
1:B:191:ASP:CA	1:B:193:ARG:NH2	2.81	0.44
1:B:222:GLU:O	1:B:226:VAL:HG23	2.18	0.44
1:A:111:LEU:HD13	1:A:115:LEU:HD21	1.99	0.43
1:A:158:ARG:HA	1:A:158:ARG:HD3	1.34	0.43
1:A:448:ILE:CG2	1:A:449:ARG:N	2.81	0.43
1:B:156:LEU:CB	1:B:157:PRO:HD3	2.48	0.43
1:A:125:LEU:HD13	1:A:210:LEU:HD11	2.00	0.43
1:B:195:PHE:O	1:B:196:PHE:C	2.57	0.43
1:A:236:HIS:ND1	1:A:236:HIS:N	2.59	0.43
1:B:191:ASP:HA	1:B:193:ARG:HH22	1.82	0.43
1:B:105:ARG:N	1:B:106:PRO:CD	2.82	0.43
1:B:275:GLY:HA3	1:B:303:GLY:HA3	2.01	0.43
1:A:125:LEU:CD1	1:A:210:LEU:HD11	2.49	0.43
1:B:341:ILE:O	1:B:347:ILE:HG12	2.19	0.43
1:A:448:ILE:HG22	1:A:449:ARG:N	2.34	0.42
1:B:185:ILE:HA	1:B:186:PRO:HD3	1.89	0.42
1:B:347:ILE:HD13	1:B:347:ILE:N	2.34	0.42
1:B:320:SER:OG	1:B:324:ARG:NH2	2.52	0.42
1:A:128:PHE:O	1:A:128:PHE:CG	2.72	0.42
1:A:132:ALA:CB	1:A:139:THR:HG21	2.50	0.42
1:B:137:LEU:O	1:B:139:THR:HG22	2.19	0.42
1:A:31:ALA:HB1	1:A:41:VAL:HG11	2.01	0.42
1:A:186:PRO:HD2	1:A:189:MET:CE	2.49	0.42
1:B:17:PRO:HD2	1:B:44:ILE:O	2.20	0.42
1:A:105:ARG:CB	1:A:106:PRO:HD3	2.39	0.42
1:A:220:TRP:CE2	1:A:433:LYS:HG3	2.55	0.42
1:B:149:PHE:CE1	1:B:377:ASP:HA	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:204:VAL:O	1:B:207:LEU:CB	2.58	0.42
1:B:207:LEU:HD23	1:B:207:LEU:HA	1.85	0.42
1:B:221:LEU:HD12	1:B:388:VAL:HG21	2.02	0.42
1:B:12:HIS:CE1	1:B:115:LEU:HD23	2.54	0.42
1:B:227:GLU:O	1:B:231:ARG:HG3	2.20	0.41
1:B:122:PHE:HB3	1:B:123:PRO:CD	2.48	0.41
1:B:158:ARG:O	1:B:161:GLU:HB2	2.20	0.41
1:B:37:ARG:HG2	1:B:449:ARG:CZ	2.50	0.41
1:B:304:LYS:HD3	1:B:304:LYS:HA	1.54	0.41
1:B:306:ASP:OD1	1:B:306:ASP:N	2.50	0.41
1:B:407:ARG:H	1:B:407:ARG:HG3	1.49	0.41
1:B:28:LEU:HD23	1:B:28:LEU:HA	1.85	0.41
1:B:34:LEU:O	1:B:39:CYS:HB2	2.20	0.41
1:B:116:SER:CB	1:B:455:LEU:HD11	2.50	0.41
1:B:149:PHE:CG	1:B:377:ASP:CB	3.00	0.41
1:B:294:TRP:CE2	1:B:418:ARG:HG3	2.55	0.41
1:A:413:LYS:HD2	1:A:413:LYS:HA	1.61	0.41
1:A:447:LEU:HA	1:A:447:LEU:HD23	1.82	0.41
1:B:177:LEU:HD23	1:B:177:LEU:HA	1.72	0.41
1:B:284:GLU:OE2	1:B:407:ARG:NH1	2.54	0.41
1:A:213:VAL:CG2	1:A:237:ILE:HA	2.51	0.41
1:B:21:MET:HG3	1:B:335:TRP:CH2	2.56	0.41
1:B:42:SER:HB3	1:B:71:LEU:CD1	2.51	0.41
1:B:129:SER:CB	1:B:130:PRO:HD3	2.51	0.41
1:B:255:LEU:N	1:B:256:PRO:CD	2.83	0.41
1:B:285:PHE:HD1	1:B:410:ILE:HG21	1.86	0.41
1:B:369:VAL:O	1:B:391:GLY:HA2	2.21	0.41
1:B:93:GLN:O	1:B:97:ILE:HG13	2.21	0.41
1:B:213:VAL:O	1:B:237:ILE:HA	2.20	0.41
1:B:240:ILE:HG23	1:B:240:ILE:O	2.20	0.41
1:A:342:LEU:HB3	1:A:365:LEU:HD13	2.02	0.41
1:A:366:GLY:C	1:A:427:VAL:CG1	2.89	0.41
1:A:186:PRO:HA	1:A:187:PRO:HD3	1.97	0.41
1:B:374:PRO:HB2	1:B:400:ALA:CB	2.46	0.41
1:B:451:LEU:C	1:B:451:LEU:CD1	2.86	0.41
1:B:236:HIS:ND1	1:B:236:HIS:N	2.64	0.41
1:A:156:LEU:HA	1:A:156:LEU:HD23	1.77	0.40
1:A:298:TRP:HD1	1:A:299:VAL:N	2.19	0.40
1:A:213:VAL:O	1:A:213:VAL:CG2	2.70	0.40
1:B:111:LEU:HD23	1:B:111:LEU:HA	1.89	0.40
1:B:281:GLN:O	1:B:284:GLU:HB2	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:336:ALA:O	2:B:500:UPG:N3	2.36	0.40
1:B:339:GLU:N	3:B:602:HOH:O	2.35	0.40
1:B:426:ARG:HB3	3:B:608:HOH:O	2.21	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	429/459 (94%)	420 (98%)	7 (2%)	2 (0%)	29	57
1	B	435/459 (95%)	428 (98%)	6 (1%)	1 (0%)	47	75
All	All	864/918 (94%)	848 (98%)	13 (2%)	3 (0%)	41	68

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	374	PRO
1	B	374	PRO
1	A	113	PRO

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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	343/396 (87%)	302 (88%)	41 (12%)	5 13
1	B	355/396 (90%)	299 (84%)	56 (16%)	2 6
All	All	698/792 (88%)	601 (86%)	97 (14%)	3 9

All (97) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	GLN
1	A	10	ARG
1	A	15	VAL
1	A	42	SER
1	A	45	SER
1	A	57	SER
1	A	60	SER
1	A	96	SER
1	A	99	ASN
1	A	118	ILE
1	A	124	VAL
1	A	139	THR
1	A	151	SER
1	A	156	LEU
1	A	158	ARG
1	A	159	LEU
1	A	171	GLU
1	A	177	LEU
1	A	189	MET
1	A	202	SER
1	A	219	THR
1	A	220	TRP
1	A	227	GLU
1	A	236	HIS
1	A	240	ILE
1	A	273	SER
1	A	278	THR
1	A	295	ARG
1	A	305	VAL
1	A	328	ARG
1	A	333	LYS
1	A	345	SER

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Mol	Chain	Res	Type
1	A	407	ARG
1	A	431	ARG
1	A	435	ARG
1	A	437	GLU
1	A	439	GLU
1	A	440	SER
1	A	446	THR
1	A	448	ILE
1	A	451	LEU
1	B	9	GLN
1	B	49	THR
1	B	52	ASP
1	B	68	ILE
1	B	73	ILE
1	B	91	PHE
1	B	110	SER
1	B	112	SER
1	B	118	ILE
1	B	120	VAL
1	B	129	SER
1	B	139	THR
1	B	141	THR
1	B	146	SER
1	B	151	SER
1	B	156	LEU
1	B	158	ARG
1	B	159	LEU
1	B	165	ASP
1	B	177	LEU
1	B	180	ILE
1	B	191	ASP
1	B	192	ARG
1	B	193	ARG
1	B	194	HIS
1	B	195	PHE
1	B	198	GLU
1	B	202	SER
1	B	213	VAL
1	B	219	THR
1	B	231	ARG
1	B	236	HIS
1	B	240	ILE

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Mol	Chain	Res	Type
1	B	246	ILE
1	B	260	GLU
1	B	273	SER
1	B	274	ARG
1	B	289	LEU
1	B	295	ARG
1	B	328	ARG
1	B	341	ILE
1	B	345	SER
1	B	386	GLU
1	B	403	ARG
1	B	405	ILE
1	B	407	ARG
1	B	413	LYS
1	B	424	ARG
1	B	431	ARG
1	B	432	GLU
1	B	433	LYS
1	B	435	ARG
1	B	440	SER
1	B	448	ILE
1	B	451	LEU
1	B	456	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	74	GLN
1	A	194	HIS
1	B	35	HIS
1	B	194	HIS
1	B	261	GLN
1	B	338	GLN

5.3.3 RNA (i)

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

2 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	UPG	A	500	-	35,38,38	1.24	4 (11%)	53,58,58	1.69	11 (20%)
2	UPG	B	500	-	35,38,38	0.35	0	53,58,58	0.58	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
2	UPG	A	500	-	-	7/23/59/59	0/3/3/3
2	UPG	B	500	-	-	3/23/59/59	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	UPG	C4-N3	-3.63	1.32	1.38
2	A	500	UPG	C2-N3	-3.31	1.32	1.38
2	A	500	UPG	C5-C4	-2.62	1.37	1.43
2	A	500	UPG	C6-N1	-2.61	1.31	1.38

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	500	UPG	C4-N3-C2	-5.57	119.23	126.58
2	A	500	UPG	N3-C2-N1	5.13	121.70	114.89
2	A	500	UPG	C5-C4-N3	3.74	120.43	114.84
2	A	500	UPG	O4-C4-C5	-2.96	119.96	125.16
2	A	500	UPG	O2-C2-N1	-2.89	118.95	122.79
2	A	500	UPG	O6'-C6'-C5'	-2.57	102.48	111.29
2	A	500	UPG	C6'-C5'-C4'	-2.50	107.16	113.00
2	A	500	UPG	C5C-C4C-C3C	-2.19	106.97	115.18
2	A	500	UPG	C5-C6-N1	-2.15	118.20	121.81
2	A	500	UPG	PB-O3B-C1'	2.14	128.01	119.74
2	A	500	UPG	C2C-C1C-N1	-2.01	107.51	113.22

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	500	UPG	C3C-C4C-C5C-O5C
2	A	500	UPG	C5C-O5C-PA-O1A
2	B	500	UPG	PA-O3A-PB-O3B
2	B	500	UPG	O5'-C1'-O3B-PB
2	A	500	UPG	O4C-C4C-C5C-O5C
2	A	500	UPG	C1'-O3B-PB-O3A
2	A	500	UPG	C5C-O5C-PA-O3A
2	A	500	UPG	C5C-O5C-PA-O2A
2	A	500	UPG	C1'-O3B-PB-O2B
2	B	500	UPG	O4C-C4C-C5C-O5C

There are no ring outliers.

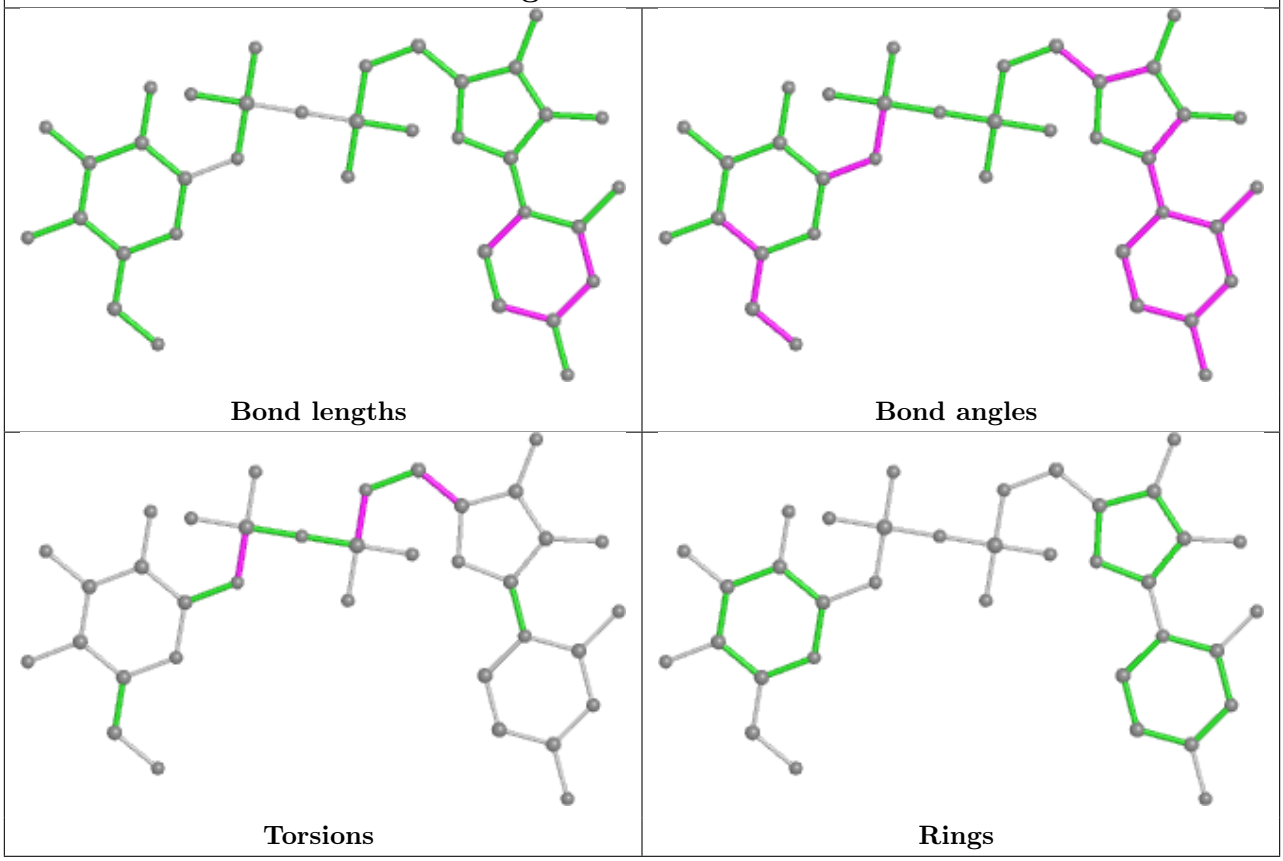
2 monomers are involved in 5 short contacts:

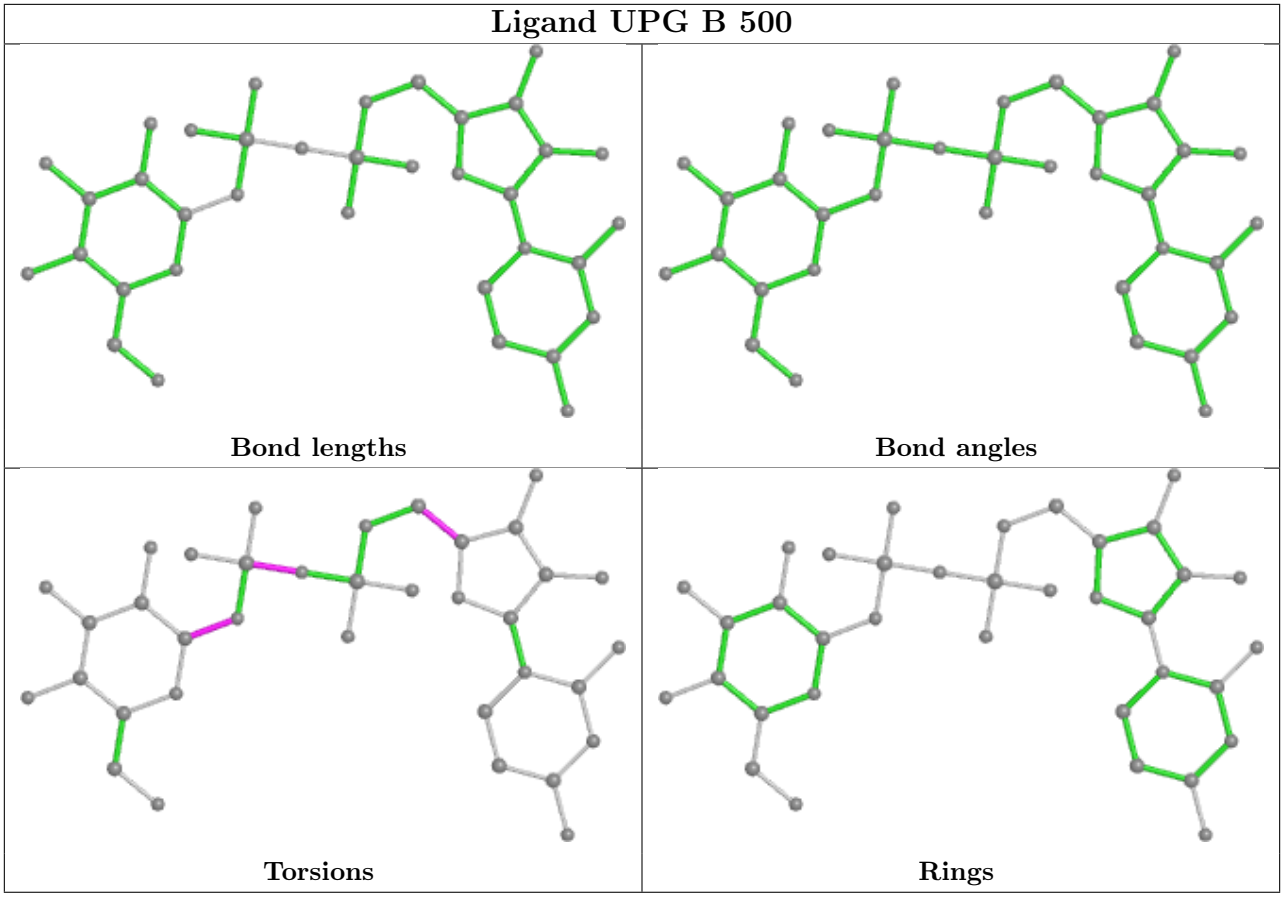
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	500	UPG	1	0
2	B	500	UPG	4	0

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 then 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.

Ligand UPG A 500





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	435/459 (94%)	-0.07	14 (3%)	47	42	22, 34, 61, 103	0
1	B	439/459 (95%)	0.16	26 (5%)	22	17	21, 39, 72, 106	0
All	All	874/918 (95%)	0.05	40 (4%)	32	27	21, 37, 69, 106	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	168	LYS	6.9
1	A	164	ASP	6.5
1	B	252	ASP	6.1
1	B	457	ILE	6.1
1	B	253	MET	5.2
1	B	76	LEU	5.1
1	B	250	GLU	5.1
1	A	167	SER	4.5
1	B	251	SER	4.4
1	B	164	ASP	4.3
1	A	248	ALA	4.1
1	B	163	GLU	4.1
1	A	165	ASP	4.1
1	B	168	LYS	4.0
1	A	88	ASP	3.8
1	A	254	ASP	3.8
1	B	194	HIS	3.7
1	B	77	PRO	3.6
1	B	254	ASP	3.6
1	A	195	PHE	3.4
1	B	318	GLY	3.4
1	B	88	ASP	3.3
1	B	195	PHE	3.2
1	B	92	ILE	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	193	ARG	3.0
1	B	306	ASP	2.8
1	A	166	ILE	2.8
1	A	302	SER	2.7
1	B	248	ALA	2.7
1	A	76	LEU	2.6
1	B	249	GLU	2.6
1	A	162	GLN	2.5
1	B	305	VAL	2.4
1	A	313	THR	2.3
1	B	191	ASP	2.3
1	B	456	LYS	2.2
1	B	304	LYS	2.2
1	A	92	ILE	2.2
1	B	9	GLN	2.1
1	B	313	THR	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 monosaccharides in this entry.

6.4 Ligands (i)

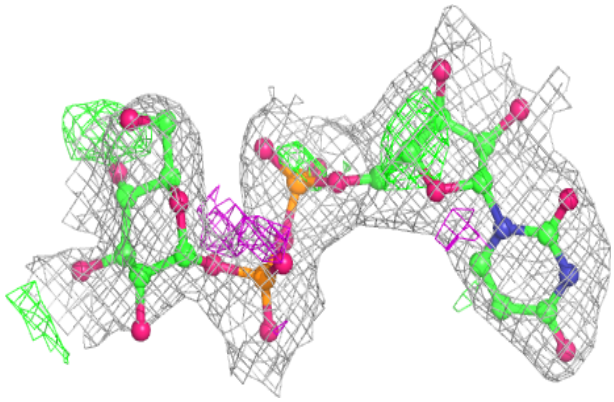
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

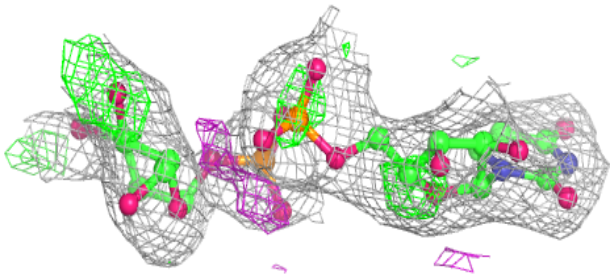
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	UPG	A	500	36/36	0.85	0.33	55,55,59,69	0
2	UPG	B	500	36/36	0.86	0.31	55,55,68,69	0

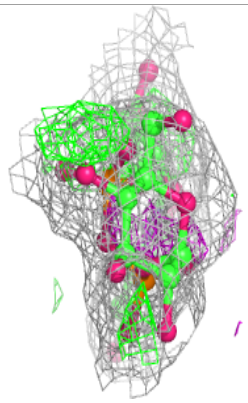
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around UPG A 500:

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

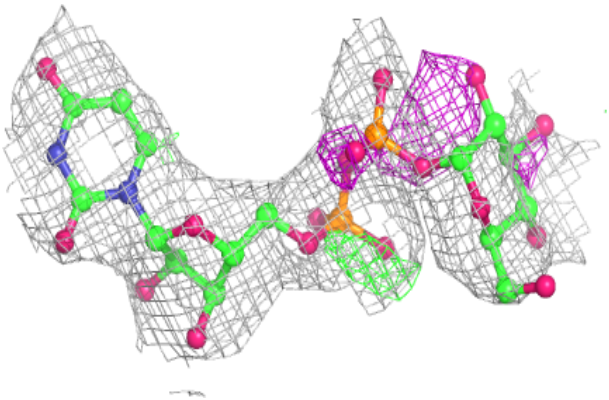


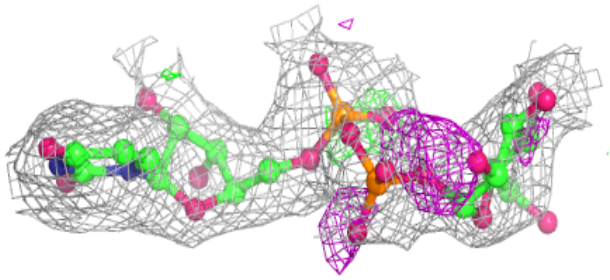


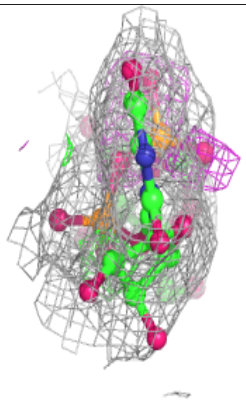


Electron density around UPG B 500:

$2mF_o$-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.