Full wwPDB X-ray Structure Validation Report (i)

Jan 29, 2024 – 07:16 PM EST

PDB ID	:	1EVL
Title	:	CRYSTAL STRUCTURE OF A TRUNCATED FORM OF THREONYL-TRNA SYNTHETASE WITH A THREONYL ADENYLATE ANALOG
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Deposited on	:	2000-04-20
Resolution	:	1.55 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:

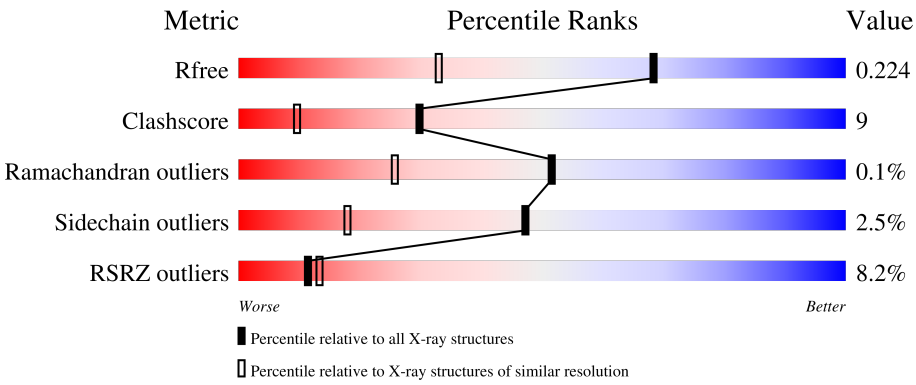
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (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 1.55 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	1483 (1.56-1.56)
Clashscore	141614	1529 (1.56-1.56)
Ramachandran outliers	138981	1498 (1.56-1.56)
Sidechain outliers	138945	1495 (1.56-1.56)
RSRZ outliers	127900	1465 (1.56-1.56)

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	401	7% 85% 15%
1	B	401	8% 83% 17%
1	C	401	8% 79% 19% •
1	D	401	9% 79% 20% •

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 14730 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.

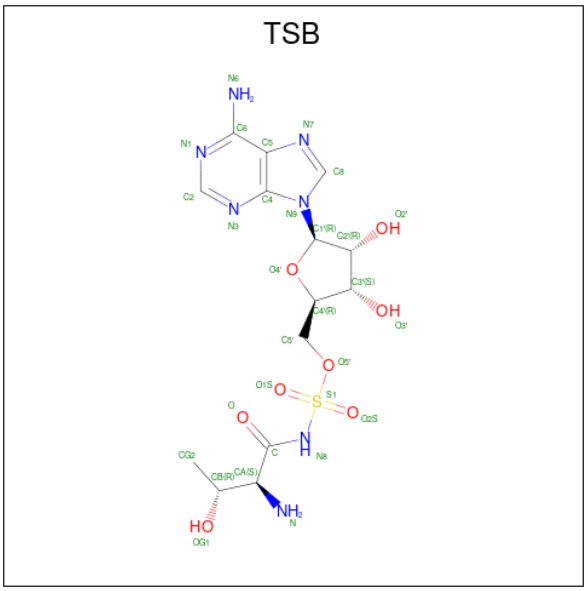
- Molecule 1 is a protein called THREONYL-TRNA SYNTHETASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	401	3265	2063	576	603	23	0	0	0
1	B	401	3268	2064	576	605	23	0	0	0
1	C	401	3268	2064	576	605	23	0	0	0
1	D	401	3261	2061	576	601	23	0	0	0

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 3 is 5'-O-(N-(L-THREONYL)-SULFAMOYL)ADENOSINE (three-letter code: TSB) (formula: $C_{14}H_{21}N_7O_8S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
3	A	1	30	14	7	8	1	0	0
			Total	C	N	O	S		
3	B	1	30	14	7	8	1	0	0
			Total	C	N	O	S		
3	C	1	30	14	7	8	1	0	0
			Total	C	N	O	S		
3	D	1	30	14	7	8	1	0	0

- Molecule 4 is water.

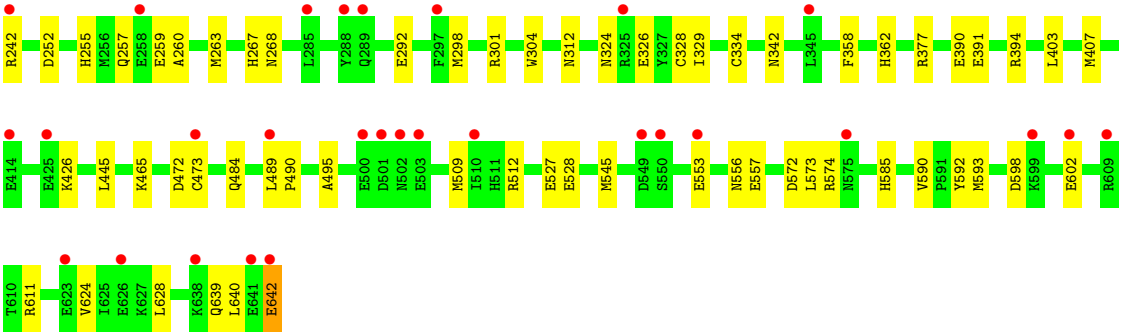
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	395	395	395	0	0
			Total	O		
4	B	390	390	390	0	0
			Total	O		
4	C	397	397	397	0	0
			Total	O		
4	D	362	362	362	0	0

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.

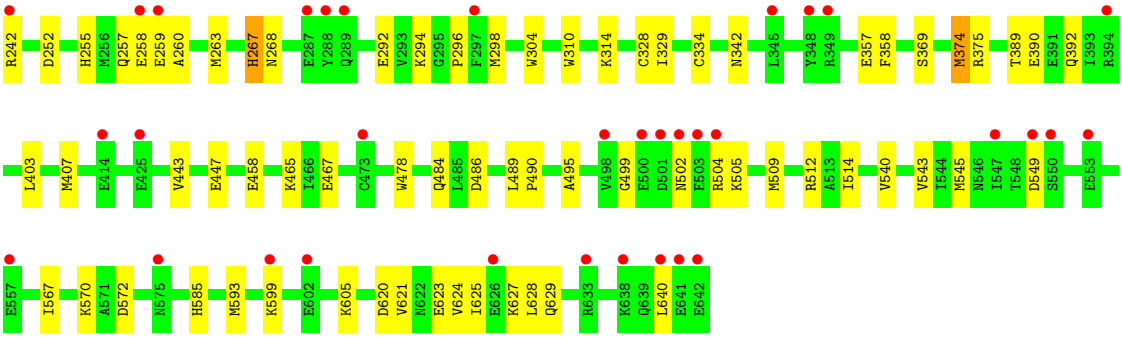
• Molecule 1: THREONYL-TRNA SYNTHETASE

Chain A: 7% 85% 15%



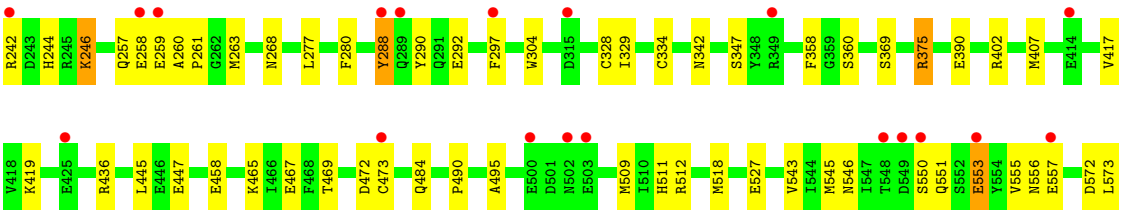
• Molecule 1: THREONYL-TRNA SYNTHETASE

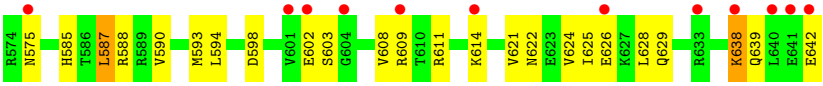
Chain B: 8% 83% 17%



• Molecule 1: THREONYL-TRNA SYNTHETASE

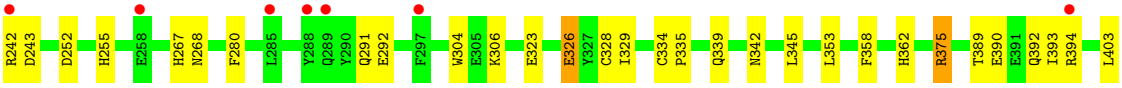
Chain C: 8% 79% 19%

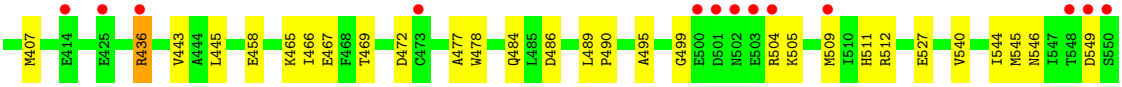


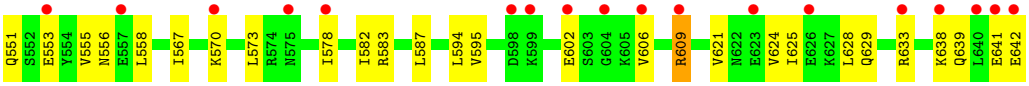


- Molecule 1: THREONYL-TRNA SYNTHETASE

Chain D:







4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	86.70Å 111.10Å 135.20Å 90.00° 93.50° 90.00°	Depositor
Resolution (Å)	30.00 – 1.55 29.89 – 1.55	Depositor EDS
% Data completeness (in resolution range)	(Not available) (30.00-1.55) 95.4 (29.89-1.55)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.80 (at 1.54Å)	Xtriage
Refinement program	CNS 0.4	Depositor
R, R_{free}	0.215 , 0.226 0.213 , 0.224	Depositor DCC
R_{free} test set	18259 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å^2)	16.5	Xtriage
Anisotropy	0.171	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.38 , 41.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	14730	wwPDB-VP
Average B, all atoms (Å^2)	21.0	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 50.85 % 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 6.1009e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[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: TSB, ZN

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	0/3336	0.63	1/4492 (0.0%)
1	B	0.35	0/3339	0.63	1/4496 (0.0%)
1	C	0.34	0/3339	0.62	1/4496 (0.0%)
1	D	0.35	0/3332	0.63	1/4487 (0.0%)
All	All	0.35	0/13346	0.63	4/17971 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	358	PHE	N-CA-C	-5.61	95.86	111.00
1	D	358	PHE	N-CA-C	-5.45	96.29	111.00
1	A	358	PHE	N-CA-C	-5.43	96.35	111.00
1	B	358	PHE	N-CA-C	-5.22	96.91	111.00

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	3265	0	3197	55	0
1	B	3268	0	3199	60	0
1	C	3268	0	3199	67	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	3261	0	3193	83	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	30	0	21	1	0
3	B	30	0	20	1	0
3	C	30	0	20	1	0
3	D	30	0	21	1	0
4	A	395	0	0	1	0
4	B	390	0	0	4	0
4	C	397	0	0	7	0
4	D	362	0	0	12	0
All	All	14730	0	12870	245	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 9.

All (245) 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:D:375:ARG:HH11	1:D:375:ARG:HB2	1.03	1.12
1:D:594:LEU:HB3	1:D:606:VAL:HG21	1.36	1.04
1:A:263:MET:HE1	1:B:298:MET:H	1.30	0.93
1:D:582:ILE:HD12	1:D:609:ARG:HD3	1.54	0.90
1:D:375:ARG:HH11	1:D:375:ARG:CB	1.85	0.90
1:D:466:ILE:HG22	4:D:5173:HOH:O	1.72	0.89
1:B:499:GLY:HA3	1:B:505:LYS:HE2	1.55	0.88
1:B:627:LYS:HD3	1:B:640:LEU:HD21	1.53	0.88
1:A:263:MET:CE	1:B:298:MET:H	1.87	0.87
1:D:375:ARG:HB2	1:D:375:ARG:NH1	1.89	0.86
1:A:390:GLU:HG2	1:A:391:GLU:OE2	1.80	0.82
1:D:594:LEU:HB3	1:D:606:VAL:CG2	2.10	0.81
1:B:465:LYS:HB3	1:B:484:GLN:HG2	1.63	0.81
1:A:574:ARG:HH11	1:A:574:ARG:HG3	1.43	0.80
1:B:540:VAL:HG11	1:B:570:LYS:HD3	1.64	0.80
1:B:543:VAL:HG22	1:B:570:LYS:HG3	1.63	0.79
1:A:263:MET:HE2	1:B:298:MET:O	1.84	0.77
1:B:260:ALA:HB1	1:B:263:MET:HB2	1.66	0.75
1:A:263:MET:HE1	1:B:298:MET:N	2.02	0.75
1:A:257:GLN:HE21	1:A:259:GLU:CG	1.99	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:465:LYS:HG2	1:A:484:GLN:HG2	1.69	0.74
1:A:292:GLU:H	1:B:268:ASN:HD22	1.36	0.74
1:C:614:LYS:HD3	4:C:4209:HOH:O	1.88	0.73
1:C:268:ASN:HD22	1:D:292:GLU:H	1.35	0.72
1:D:393:ILE:HD12	1:D:436:ARG:HH22	1.55	0.72
1:C:246:LYS:HE3	1:C:246:LYS:HA	1.73	0.71
1:C:292:GLU:H	1:D:268:ASN:HD22	1.39	0.71
1:C:484:GLN:HE22	3:C:4002:TSB:HN8	1.38	0.71
1:A:553:GLU:O	1:A:557:GLU:HG3	1.91	0.71
1:A:255:HIS:ND1	1:A:267:HIS:HE1	1.90	0.70
1:D:255:HIS:ND1	1:D:267:HIS:HE1	1.90	0.69
1:B:621:VAL:O	1:B:624:VAL:HG22	1.93	0.69
1:A:329:ILE:HD11	1:B:329:ILE:HD11	1.75	0.68
1:A:268:ASN:HD22	1:B:292:GLU:H	1.41	0.68
1:D:465:LYS:HG2	1:D:484:GLN:HG2	1.76	0.68
1:A:484:GLN:HE22	3:A:2002:TSB:HN8	1.43	0.67
1:D:582:ILE:CD1	1:D:609:ARG:HD3	2.25	0.67
1:B:255:HIS:ND1	1:B:267:HIS:HE1	1.93	0.66
1:C:469:THR:HG22	4:C:4199:HOH:O	1.96	0.66
1:D:390:GLU:CD	1:D:390:GLU:H	1.99	0.65
1:C:553:GLU:O	1:C:557:GLU:HG3	1.97	0.65
1:B:484:GLN:HE22	3:B:3002:TSB:HN8	1.44	0.65
1:B:403:LEU:HG	1:B:407:MET:CE	2.26	0.64
1:B:623:GLU:HG3	1:B:627:LYS:HE2	1.79	0.64
1:B:567:ILE:HD13	1:B:628:LEU:HD12	1.78	0.64
1:C:260:ALA:HB1	1:C:263:MET:HB2	1.79	0.64
1:C:621:VAL:O	1:C:624:VAL:HG22	1.98	0.64
1:B:252:ASP:OD1	1:B:267:HIS:HD2	1.81	0.63
1:C:419:LYS:HE3	1:C:469:THR:HG21	1.81	0.63
1:D:252:ASP:OD1	1:D:267:HIS:HD2	1.82	0.63
1:D:638:LYS:HB2	1:D:642:GLU:HG3	1.81	0.62
1:C:465:LYS:HE3	1:C:467:GLU:OE2	2.00	0.62
1:D:490:PRO:HG3	1:D:509:MET:SD	2.40	0.62
1:A:572:ASP:OD2	1:A:585:HIS:HE1	1.81	0.61
1:A:242:ARG:HD2	1:A:473:CYS:SG	2.41	0.61
1:A:252:ASP:OD1	1:A:267:HIS:HD2	1.83	0.60
1:A:403:LEU:HG	1:A:407:MET:HE3	1.83	0.60
1:D:362:HIS:HD2	4:D:5011:HOH:O	1.85	0.60
1:A:257:GLN:HE21	1:A:259:GLU:HG2	1.67	0.59
1:D:484:GLN:HE22	3:D:5002:TSB:HN8	1.50	0.59
1:A:585:HIS:HD2	4:A:2169:HOH:O	1.86	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:329:ILE:HD11	1:D:329:ILE:HD11	1.83	0.59
1:C:390:GLU:H	1:C:390:GLU:CD	2.04	0.59
1:B:624:VAL:O	1:B:628:LEU:HG	2.03	0.59
1:C:257:GLN:HE21	1:C:259:GLU:CG	2.15	0.58
1:A:390:GLU:CD	1:A:390:GLU:H	2.05	0.58
1:B:257:GLN:HE21	1:B:259:GLU:HG2	1.69	0.58
1:C:242:ARG:HG3	1:C:473:CYS:HB3	1.86	0.58
1:C:594:LEU:HD22	1:C:608:VAL:HG22	1.85	0.58
1:C:638:LYS:H	1:C:638:LYS:HD3	1.69	0.58
1:B:242:ARG:HD2	4:B:3344:HOH:O	2.03	0.57
1:D:558:LEU:HD11	1:D:621:VAL:HG21	1.85	0.57
1:A:598:ASP:O	1:A:602:GLU:HG2	2.04	0.57
1:C:588:ARG:HG3	4:C:4080:HOH:O	2.04	0.57
1:B:572:ASP:OD2	1:B:585:HIS:HE1	1.88	0.57
1:B:627:LYS:CD	1:B:640:LEU:HD21	2.32	0.57
1:C:375:ARG:HG2	1:C:375:ARG:HH11	1.70	0.57
1:D:465:LYS:HD2	1:D:484:GLN:NE2	2.19	0.57
1:B:443:VAL:O	1:B:447:GLU:HG3	2.04	0.56
1:A:324:ASN:ND2	1:D:323:GLU:H	2.02	0.56
1:C:572:ASP:OD2	1:C:585:HIS:HE1	1.86	0.56
1:C:624:VAL:O	1:C:628:LEU:HG	2.05	0.56
1:A:312:ASN:OD1	1:A:426:LYS:HG3	2.06	0.56
1:A:574:ARG:HG3	1:A:574:ARG:NH1	2.19	0.56
1:D:291:GLN:OE1	1:D:353:LEU:HD21	2.06	0.56
1:B:294:LYS:HD2	1:B:357:GLU:OE2	2.05	0.56
1:C:417:VAL:HB	1:C:469:THR:OG1	2.05	0.56
1:C:465:LYS:CB	1:C:484:GLN:HG2	2.36	0.56
1:C:490:PRO:HG3	1:C:509:MET:CE	2.36	0.56
1:C:609:ARG:HD3	1:C:614:LYS:O	2.05	0.56
1:C:280:PHE:CE2	1:C:407:MET:HG2	2.41	0.55
1:A:624:VAL:O	1:A:628:LEU:HG	2.06	0.55
1:D:345:LEU:HD12	1:D:504:ARG:HH22	1.71	0.55
1:D:583:ARG:O	1:D:587:LEU:HD13	2.06	0.55
1:A:592:TYR:OH	1:A:640:LEU:HD13	2.07	0.55
1:B:465:LYS:CB	1:B:484:GLN:HG2	2.35	0.55
1:C:402:ARG:HG3	4:C:4359:HOH:O	2.06	0.55
1:D:465:LYS:CG	1:D:484:GLN:HG2	2.35	0.55
1:D:490:PRO:HG3	1:D:509:MET:HG3	1.88	0.55
1:D:578:ILE:HD11	1:D:595:VAL:HG21	1.88	0.54
1:C:465:LYS:HB3	1:C:484:GLN:HG2	1.88	0.54
1:B:390:GLU:CD	1:B:390:GLU:H	2.09	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:490:PRO:HG3	1:A:509:MET:CE	2.37	0.54
1:B:504:ARG:HG2	1:B:504:ARG:HH11	1.72	0.53
1:B:543:VAL:HG22	1:B:570:LYS:CG	2.35	0.53
1:B:504:ARG:HG2	1:B:504:ARG:NH1	2.23	0.53
1:B:260:ALA:HB1	1:B:263:MET:CB	2.35	0.53
1:D:306:LYS:HE2	4:D:5243:HOH:O	2.08	0.53
1:A:301:ARG:HG3	1:A:328:CYS:SG	2.48	0.53
1:C:268:ASN:ND2	1:D:292:GLU:H	2.05	0.52
1:D:394:ARG:HH21	1:D:443:VAL:HG11	1.75	0.52
1:D:389:THR:OG1	1:D:392:GLN:HG3	2.10	0.52
1:C:277:LEU:HB2	1:C:518:MET:CE	2.40	0.52
1:A:472:ASP:HB2	1:A:527:GLU:OE1	2.08	0.51
1:B:490:PRO:HD3	1:B:509:MET:HE3	1.93	0.51
1:B:540:VAL:HG11	1:B:570:LYS:CD	2.38	0.51
1:D:545:MET:CE	1:D:595:VAL:HG22	2.40	0.51
1:B:304:TRP:CD1	1:B:328:CYS:HB2	2.46	0.51
1:C:556:ASN:HD21	1:C:573:LEU:HD11	1.74	0.51
1:B:374:MET:HG2	1:B:478:TRP:CE2	2.46	0.51
1:D:540:VAL:HG11	1:D:570:LYS:HZ3	1.75	0.50
1:D:578:ILE:HD12	1:D:595:VAL:HG11	1.92	0.50
1:A:298:MET:HG2	1:B:263:MET:CE	2.42	0.50
1:D:469:THR:HG21	1:D:477:ALA:HB1	1.93	0.50
1:A:304:TRP:CD1	1:A:328:CYS:HB2	2.47	0.50
1:A:391:GLU:CD	1:A:391:GLU:H	2.14	0.50
1:B:257:GLN:HE21	1:B:259:GLU:CG	2.25	0.50
1:D:472:ASP:HB2	1:D:527:GLU:OE1	2.11	0.50
1:C:280:PHE:CZ	1:C:407:MET:HG2	2.47	0.50
1:B:342:ASN:HD21	1:B:495:ALA:HA	1.76	0.49
1:D:403:LEU:HG	1:D:407:MET:HE3	1.93	0.49
1:D:486:ASP:OD2	1:D:489:LEU:HG	2.13	0.49
1:D:545:MET:CE	1:D:582:ILE:HD11	2.43	0.49
1:A:292:GLU:H	1:B:268:ASN:ND2	2.07	0.49
1:C:369:SER:O	1:C:375:ARG:HG2	2.13	0.48
1:C:447:GLU:HG2	4:C:4377:HOH:O	2.13	0.48
1:D:602:GLU:HG3	4:D:5308:HOH:O	2.13	0.48
1:C:417:VAL:O	1:C:469:THR:HG23	2.13	0.48
1:C:550:SER:HB3	1:C:551:GLN:NE2	2.28	0.48
1:A:260:ALA:HB1	1:A:263:MET:HB2	1.95	0.48
1:D:553:GLU:HB3	4:D:5077:HOH:O	2.13	0.48
1:D:304:TRP:CD1	1:D:328:CYS:HB2	2.49	0.48
1:A:298:MET:HG2	1:B:263:MET:HE1	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:258:GLU:O	1:C:261:PRO:HD3	2.14	0.47
1:C:304:TRP:CD1	1:C:328:CYS:HB2	2.49	0.47
1:D:499:GLY:HA3	1:D:505:LYS:HE2	1.97	0.47
1:D:242:ARG:HA	4:D:5092:HOH:O	2.14	0.47
1:D:545:MET:HE3	1:D:545:MET:HB2	1.51	0.47
1:D:545:MET:HE1	1:D:582:ILE:HD11	1.96	0.47
1:B:623:GLU:CG	1:B:627:LYS:HE2	2.43	0.47
1:D:567:ILE:HD13	1:D:628:LEU:HD13	1.97	0.47
1:C:590:VAL:O	1:C:611:ARG:HB3	2.14	0.47
1:D:255:HIS:ND1	1:D:267:HIS:CE1	2.78	0.47
1:D:353:LEU:HD23	4:D:5150:HOH:O	2.14	0.47
1:B:369:SER:O	1:B:375:ARG:HD2	2.15	0.47
1:C:543:VAL:HG11	1:C:585:HIS:CE1	2.50	0.47
1:B:458:GLU:O	1:B:465:LYS:HE2	2.15	0.46
1:D:342:ASN:HD21	1:D:495:ALA:HA	1.80	0.46
1:D:490:PRO:HG3	1:D:509:MET:CG	2.45	0.46
1:A:394:ARG:NH1	1:A:394:ARG:HB2	2.31	0.46
1:A:394:ARG:HB2	1:A:394:ARG:HH11	1.80	0.46
1:B:599:LYS:HB2	1:B:599:LYS:NZ	2.30	0.46
1:C:260:ALA:HB1	1:C:263:MET:CB	2.44	0.46
1:A:545:MET:CE	1:A:593:MET:HB3	2.46	0.46
1:A:263:MET:HE3	1:B:296:PRO:HB2	1.98	0.46
1:B:403:LEU:HG	1:B:407:MET:HE3	1.95	0.46
1:C:545:MET:CE	1:C:593:MET:HB3	2.46	0.46
1:B:407:MET:HE1	1:B:514:ILE:HG13	1.97	0.46
1:C:342:ASN:HD21	1:C:495:ALA:HA	1.79	0.46
1:C:472:ASP:HB2	1:C:527:GLU:OE1	2.16	0.46
1:A:490:PRO:HG3	1:A:509:MET:HE2	1.96	0.46
1:B:625:ILE:O	1:B:629:GLN:HG3	2.16	0.46
1:C:258:GLU:HG3	1:C:259:GLU:OE2	2.16	0.46
1:C:257:GLN:HE21	1:C:259:GLU:HG2	1.78	0.45
1:A:342:ASN:HD21	1:A:495:ALA:HA	1.80	0.45
1:A:574:ARG:HH11	1:A:574:ARG:CG	2.18	0.45
1:D:375:ARG:NH1	1:D:478:TRP:HZ3	2.15	0.45
1:D:544:ILE:HG22	1:D:555:VAL:HG13	1.98	0.45
1:D:639:GLN:O	1:D:642:GLU:HG2	2.17	0.45
1:B:294:LYS:HD2	1:B:294:LYS:HA	1.77	0.45
1:C:490:PRO:HG3	1:C:509:MET:HE2	1.98	0.45
1:A:574:ARG:NH1	1:A:574:ARG:CG	2.78	0.45
1:D:345:LEU:CD1	1:D:504:ARG:NH2	2.80	0.45
1:D:625:ILE:O	1:D:629:GLN:HG3	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:297:PHE:HE2	1:D:329:ILE:HD13	1.82	0.44
1:C:257:GLN:NE2	1:C:259:GLU:CG	2.80	0.44
1:C:625:ILE:O	1:C:629:GLN:HG3	2.17	0.44
1:D:280:PHE:CE2	1:D:407:MET:HG2	2.52	0.44
1:D:345:LEU:HD12	1:D:504:ARG:NH2	2.33	0.44
1:C:277:LEU:HB2	1:C:518:MET:HE1	1.98	0.44
1:D:393:ILE:HD12	1:D:436:ARG:NH2	2.27	0.44
1:A:268:ASN:ND2	1:B:292:GLU:H	2.13	0.44
1:B:490:PRO:HG2	4:B:3126:HOH:O	2.17	0.44
1:D:621:VAL:O	1:D:624:VAL:CG1	2.66	0.44
1:C:292:GLU:H	1:D:268:ASN:ND2	2.11	0.43
1:C:484:GLN:HB2	1:C:511:HIS:HB2	2.00	0.43
1:B:465:LYS:HE3	1:B:467:GLU:OE2	2.18	0.43
1:C:546:ASN:HB3	1:C:555:VAL:HG21	2.00	0.43
1:B:486:ASP:OD2	1:B:489:LEU:HG	2.18	0.43
1:C:598:ASP:O	1:C:602:GLU:HG2	2.18	0.43
1:C:622:ASN:O	1:C:626:GLU:HG3	2.18	0.43
1:D:375:ARG:NH1	1:D:478:TRP:CZ3	2.86	0.43
1:A:260:ALA:HB1	1:A:263:MET:HG3	2.01	0.43
1:B:389:THR:OG1	1:B:392:GLN:HG3	2.19	0.43
1:A:556:ASN:HD21	1:A:573:LEU:HD11	1.84	0.42
1:D:465:LYS:NZ	4:D:5044:HOH:O	2.47	0.42
1:D:540:VAL:HG11	1:D:570:LYS:NZ	2.34	0.42
1:A:255:HIS:ND1	1:A:267:HIS:CE1	2.79	0.42
1:D:467:GLU:C	4:D:5173:HOH:O	2.58	0.42
1:A:362:HIS:HA	1:A:377:ARG:O	2.19	0.42
1:D:436:ARG:HD2	4:D:5191:HOH:O	2.19	0.42
1:D:545:MET:HE2	1:D:595:VAL:HG22	2.00	0.42
1:C:258:GLU:HG2	4:C:4257:HOH:O	2.20	0.42
1:D:633:ARG:NH1	1:D:633:ARG:HG2	2.35	0.42
1:C:242:ARG:HD3	1:C:527:GLU:O	2.19	0.42
1:D:466:ILE:C	4:D:5173:HOH:O	2.58	0.42
1:A:639:GLN:O	1:A:642:GLU:HG3	2.20	0.42
1:C:244:HIS:HB3	1:C:527:GLU:HG3	2.01	0.42
1:D:621:VAL:O	1:D:624:VAL:HG13	2.20	0.42
1:A:324:ASN:HD21	1:D:323:GLU:H	1.65	0.41
1:A:590:VAL:O	1:A:611:ARG:HB3	2.20	0.41
1:C:587:LEU:HD12	1:C:587:LEU:HA	1.83	0.41
1:D:469:THR:CG2	1:D:477:ALA:HB1	2.50	0.41
1:D:556:ASN:HD21	1:D:573:LEU:HD11	1.85	0.41
4:B:3315:HOH:O	1:D:326:GLU:HB2	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:360:SER:N	4:C:4337:HOH:O	2.48	0.41
1:B:504:ARG:NH2	4:B:3153:HOH:O	2.53	0.41
1:B:605:LYS:NZ	1:B:620:ASP:OD1	2.51	0.41
1:C:639:GLN:O	1:C:642:GLU:HB2	2.21	0.41
1:A:391:GLU:CD	1:A:391:GLU:N	2.74	0.41
1:D:484:GLN:HB2	1:D:511:HIS:HB2	2.02	0.41
1:D:335:PRO:O	1:D:339:GLN:HG2	2.20	0.41
1:A:489:LEU:HB3	1:A:490:PRO:HD3	2.03	0.41
1:C:277:LEU:HB2	1:C:518:MET:HE3	2.02	0.41
1:C:458:GLU:O	1:C:465:LYS:HE2	2.20	0.41
1:D:490:PRO:CG	1:D:509:MET:HG3	2.50	0.41
1:B:310:TRP:O	1:B:314:LYS:HB2	2.20	0.41
1:C:288:TYR:HB2	1:C:290:TYR:CD1	2.55	0.41
1:D:578:ILE:O	1:D:582:ILE:HG12	2.21	0.41
1:B:545:MET:CE	1:B:593:MET:HB3	2.51	0.41
1:A:465:LYS:HE2	1:A:484:GLN:NE2	2.35	0.40
1:C:259:GLU:HB2	4:D:5247:HOH:O	2.21	0.40
1:D:546:ASN:HB2	1:D:551:GLN:HB2	2.01	0.40
1:C:621:VAL:O	1:C:624:VAL:CG2	2.69	0.40
1:D:546:ASN:HB3	1:D:555:VAL:HG21	2.03	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	399/401 (100%)	392 (98%)	7 (2%)	0	100	100
1	B	399/401 (100%)	393 (98%)	6 (2%)	0	100	100
1	C	399/401 (100%)	389 (98%)	9 (2%)	1 (0%)	41	19
1	D	399/401 (100%)	391 (98%)	7 (2%)	1 (0%)	41	19
All	All	1596/1604 (100%)	1565 (98%)	29 (2%)	2 (0%)	51	26

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	603	SER
1	D	641	GLU

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	353/356 (99%)	347 (98%)	6 (2%)	60	32
1	B	354/356 (99%)	347 (98%)	7 (2%)	55	26
1	C	354/356 (99%)	342 (97%)	12 (3%)	37	9
1	D	352/356 (99%)	342 (97%)	10 (3%)	43	14
All	All	1413/1424 (99%)	1378 (98%)	35 (2%)	47	18

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

Mol	Chain	Res	Type
1	A	326	GLU
1	A	334	CYS
1	A	445	LEU
1	A	512	ARG
1	A	528	GLU
1	A	642	GLU
1	B	258	GLU
1	B	267	HIS
1	B	334	CYS
1	B	374	MET
1	B	502	ASN
1	B	512	ARG
1	B	549	ASP
1	C	246	LYS
1	C	288	TYR
1	C	334	CYS
1	C	347	SER
1	C	375	ARG

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Mol	Chain	Res	Type
1	C	436	ARG
1	C	445	LEU
1	C	512	ARG
1	C	553	GLU
1	C	575	ASN
1	C	587	LEU
1	C	638	LYS
1	D	243	ASP
1	D	326	GLU
1	D	334	CYS
1	D	375	ARG
1	D	436	ARG
1	D	445	LEU
1	D	458	GLU
1	D	512	ARG
1	D	549	ASP
1	D	609	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (40) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	250	GLN
1	A	257	GLN
1	A	267	HIS
1	A	268	ASN
1	A	289	GLN
1	A	291	GLN
1	A	324	ASN
1	A	342	ASN
1	A	343	GLN
1	A	381	GLN
1	A	455	GLN
1	A	484	GLN
1	A	556	ASN
1	A	585	HIS
1	B	267	HIS
1	B	268	ASN
1	B	291	GLN
1	B	312	ASN
1	B	339	GLN
1	B	342	ASN
1	B	381	GLN

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Mol	Chain	Res	Type
1	B	455	GLN
1	B	502	ASN
1	B	556	ASN
1	B	585	HIS
1	C	268	ASN
1	C	291	GLN
1	C	324	ASN
1	C	342	ASN
1	C	449	ASN
1	C	455	GLN
1	C	556	ASN
1	C	585	HIS
1	D	267	HIS
1	D	268	ASN
1	D	289	GLN
1	D	342	ASN
1	D	362	HIS
1	D	381	GLN
1	D	556	ASN

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)

Of 8 ligands modelled in this entry, 4 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	TSB	A	2002	2	29,32,32	2.11	4 (13%)	33,48,48	1.01	3 (9%)
3	TSB	D	5002	2	29,32,32	2.18	4 (13%)	33,48,48	0.98	2 (6%)
3	TSB	C	4002	2	29,32,32	2.26	4 (13%)	33,48,48	1.01	3 (9%)
3	TSB	B	3002	2	29,32,32	2.24	4 (13%)	33,48,48	1.05	3 (9%)

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	TSB	A	2002	2	-	0/18/39/39	0/3/3/3
3	TSB	D	5002	2	-	0/18/39/39	0/3/3/3
3	TSB	C	4002	2	-	0/18/39/39	0/3/3/3
3	TSB	B	3002	2	-	0/18/39/39	0/3/3/3

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	5002	TSB	O2S-S1	8.19	1.49	1.42
3	C	4002	TSB	O2S-S1	8.07	1.49	1.42
3	B	3002	TSB	O2S-S1	8.01	1.49	1.42
3	A	2002	TSB	O2S-S1	7.85	1.49	1.42
3	C	4002	TSB	O1S-S1	7.06	1.48	1.42
3	B	3002	TSB	O1S-S1	7.04	1.48	1.42
3	D	5002	TSB	O1S-S1	6.55	1.48	1.42
3	A	2002	TSB	O1S-S1	6.40	1.47	1.42
3	C	4002	TSB	CA-C	-3.55	1.49	1.53
3	B	3002	TSB	C-N8	-3.41	1.31	1.37
3	B	3002	TSB	CA-C	-3.41	1.49	1.53
3	C	4002	TSB	C-N8	-3.25	1.31	1.37
3	D	5002	TSB	C-N8	-3.16	1.31	1.37
3	A	2002	TSB	C-N8	-2.95	1.31	1.37
3	A	2002	TSB	CA-C	-2.88	1.50	1.53
3	D	5002	TSB	CA-C	-2.85	1.50	1.53

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	B	3002	TSB	C-CA-N	-2.72	106.13	110.28
3	B	3002	TSB	C-N8-S1	-2.58	120.42	124.61
3	A	2002	TSB	C-CA-N	-2.45	106.54	110.28
3	A	2002	TSB	C-N8-S1	-2.44	120.66	124.61
3	C	4002	TSB	C-CA-N	-2.35	106.69	110.28
3	D	5002	TSB	C5-C6-N6	2.32	123.87	120.35
3	C	4002	TSB	C-N8-S1	-2.28	120.93	124.61
3	A	2002	TSB	C5-C6-N6	2.25	123.77	120.35
3	D	5002	TSB	C-CA-N	-2.18	106.95	110.28
3	B	3002	TSB	C5-C6-N6	2.13	123.58	120.35
3	C	4002	TSB	C5-C6-N6	2.07	123.50	120.35

There are no chirality outliers.

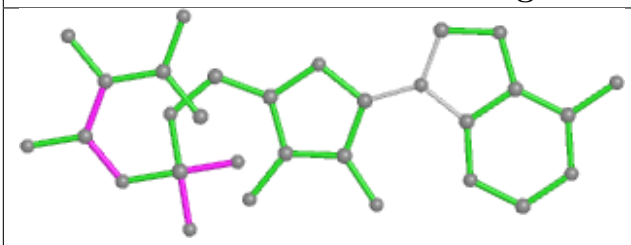
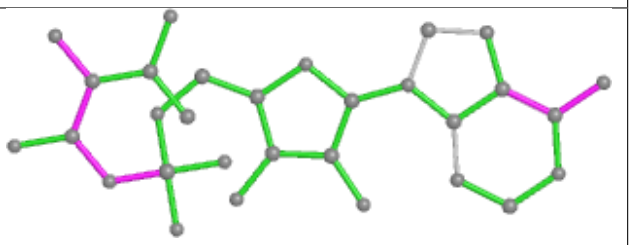
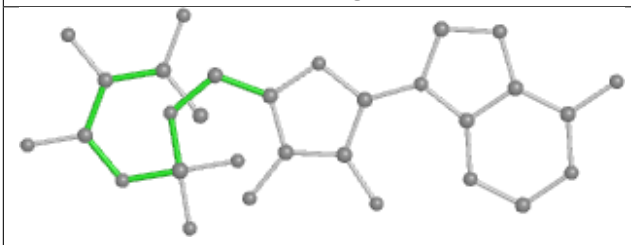
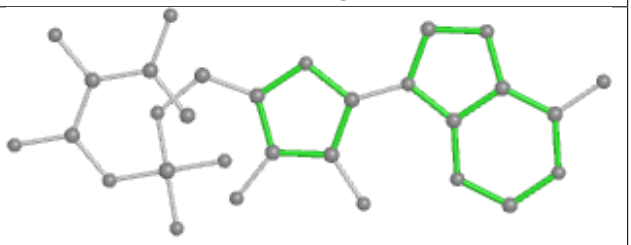
There are no torsion outliers.

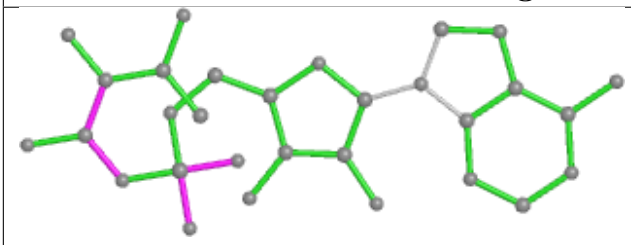
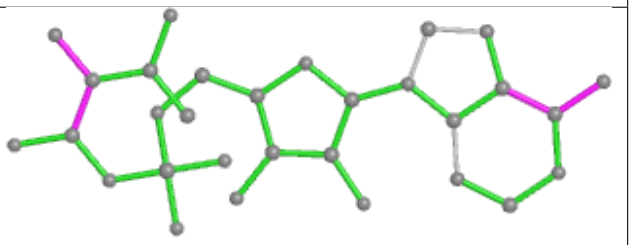
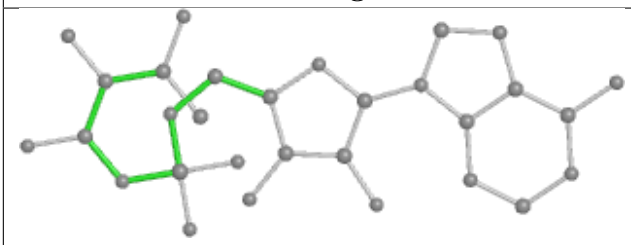
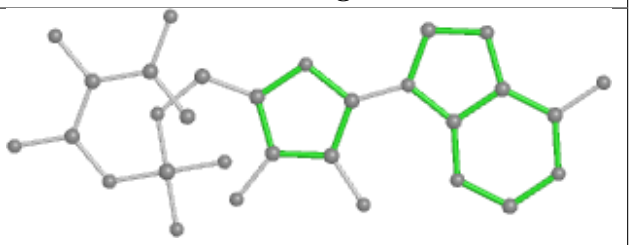
There are no ring outliers.

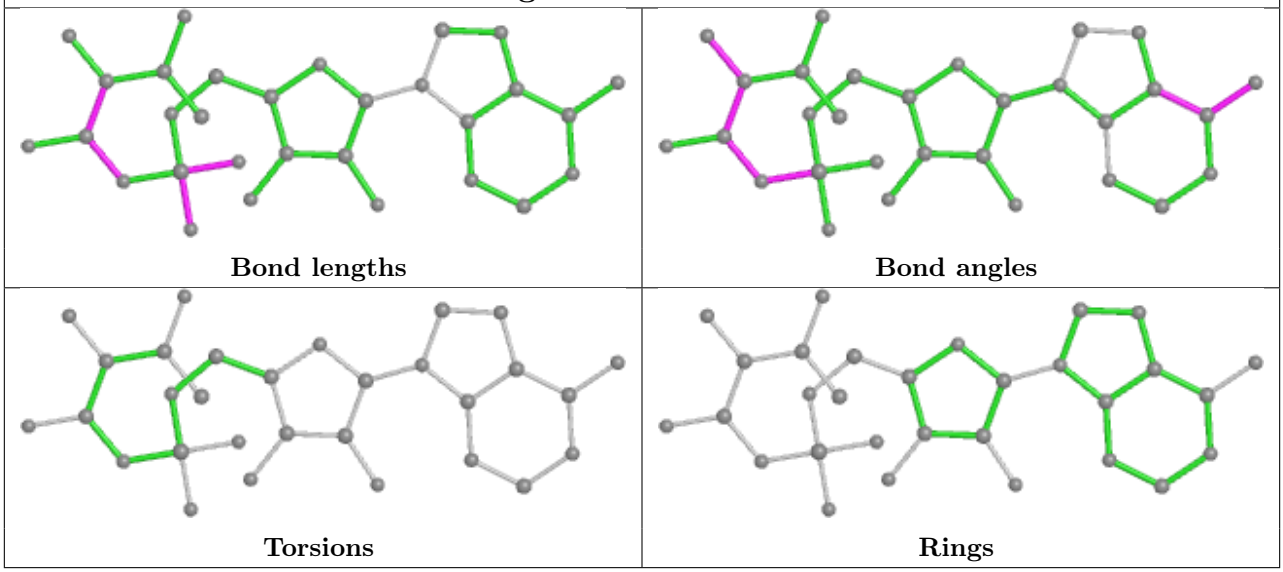
4 monomers are involved in 4 short contacts:

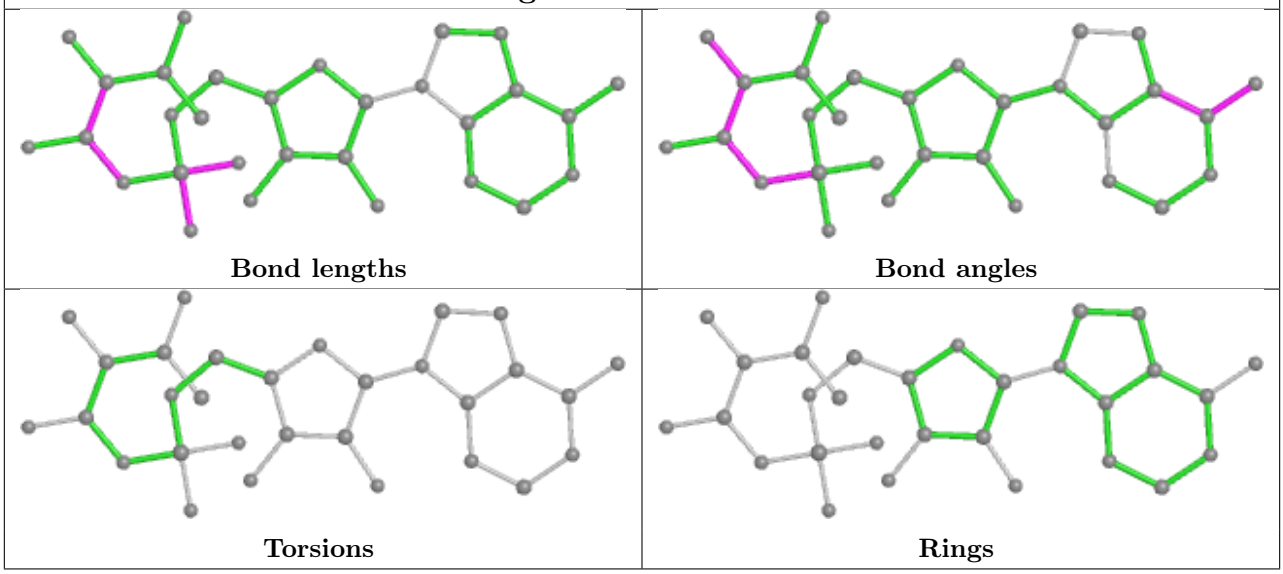
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	2002	TSB	1	0
3	D	5002	TSB	1	0
3	C	4002	TSB	1	0
3	B	3002	TSB	1	0

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 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 TSB A 2002	
 Bond lengths	 Bond angles
 Torsions	 Rings

Ligand TSB D 5002	
 Bond lengths	 Bond angles
 Torsions	 Rings

Ligand TSB C 4002	
	
Bond lengths	Bond angles
Torsions	Rings

Ligand TSB B 3002	
	
Bond lengths	Bond angles
Torsions	Rings

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	401/401 (100%)	0.40	29 (7%)	15	18	11, 19, 30, 36	0
1	B	401/401 (100%)	0.47	34 (8%)	10	12	11, 19, 32, 37	0
1	C	401/401 (100%)	0.45	31 (7%)	13	16	12, 20, 32, 38	0
1	D	401/401 (100%)	0.54	38 (9%)	8	9	12, 20, 33, 37	0
All	All	1604/1604 (100%)	0.47	132 (8%)	11	13	11, 20, 32, 38	0

All (132) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	602	GLU	5.8
1	A	502	ASN	5.6
1	D	640	LEU	5.2
1	D	642	GLU	4.8
1	A	575	ASN	4.6
1	A	642	GLU	4.6
1	D	549	ASP	4.5
1	C	642	GLU	4.5
1	B	502	ASN	4.5
1	C	549	ASP	4.4
1	B	602	GLU	4.3
1	C	633	ARG	4.2
1	D	550	SER	4.1
1	D	297	PHE	4.0
1	C	242	ARG	4.0
1	C	288	TYR	4.0
1	D	599	LYS	3.9
1	B	640	LEU	3.9
1	D	575	ASN	3.8
1	C	602	GLU	3.8
1	A	297	PHE	3.8

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Mol	Chain	Res	Type	RSRZ
1	D	504	ARG	3.8
1	B	242	ARG	3.7
1	D	503	GLU	3.7
1	B	349	ARG	3.7
1	B	503	GLU	3.6
1	A	602	GLU	3.6
1	B	641	GLU	3.5
1	D	242	ARG	3.5
1	C	473	CYS	3.5
1	D	501	ASP	3.5
1	C	502	ASN	3.5
1	A	414	GLU	3.4
1	D	578	ILE	3.4
1	B	288	TYR	3.4
1	C	258	GLU	3.4
1	D	609	ARG	3.4
1	A	503	GLU	3.4
1	A	473	CYS	3.3
1	C	575	ASN	3.3
1	A	641	GLU	3.3
1	B	642	GLU	3.3
1	C	349	ARG	3.3
1	D	557	GLU	3.2
1	C	503	GLU	3.2
1	D	633	ARG	3.2
1	D	641	GLU	3.2
1	B	549	ASP	3.2
1	C	425	GLU	3.1
1	C	500	GLU	3.1
1	A	549	ASP	3.1
1	A	425	GLU	3.1
1	B	500	GLU	3.1
1	A	501	ASP	3.0
1	B	501	ASP	3.0
1	D	502	ASN	3.0
1	A	258	GLU	2.9
1	D	436	ARG	2.9
1	B	258	GLU	2.9
1	B	498	VAL	2.9
1	B	414	GLU	2.9
1	D	500	GLU	2.9
1	C	604	GLY	2.9

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Mol	Chain	Res	Type	RSRZ
1	D	258	GLU	2.8
1	D	425	GLU	2.8
1	B	547	ILE	2.8
1	B	550	SER	2.8
1	A	242	ARG	2.8
1	A	345	LEU	2.8
1	C	640	LEU	2.8
1	D	285	LEU	2.8
1	D	289	GLN	2.7
1	C	550	SER	2.7
1	D	553	GLU	2.7
1	C	553	GLU	2.6
1	A	289	GLN	2.6
1	C	259	GLU	2.6
1	B	626	GLU	2.6
1	C	609	ARG	2.6
1	D	626	GLU	2.6
1	C	297	PHE	2.6
1	A	500	GLU	2.6
1	D	598	ASP	2.6
1	B	638	LYS	2.6
1	B	504	ARG	2.5
1	C	289	GLN	2.5
1	D	288	TYR	2.4
1	D	509	MET	2.4
1	D	473	CYS	2.4
1	B	553	GLU	2.4
1	B	425	GLU	2.4
1	A	599	LYS	2.4
1	A	285	LEU	2.4
1	A	288	TYR	2.4
1	B	473	CYS	2.4
1	D	414	GLU	2.4
1	B	348	TYR	2.4
1	B	297	PHE	2.3
1	A	325	ARG	2.3
1	B	345	LEU	2.3
1	C	641	GLU	2.3
1	D	548	THR	2.3
1	C	626	GLU	2.3
1	B	575	ASN	2.2
1	B	287	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	614	LYS	2.2
1	C	548	THR	2.2
1	D	394	ARG	2.2
1	B	557	GLU	2.2
1	C	638	LYS	2.2
1	B	633	ARG	2.2
1	C	414	GLU	2.2
1	B	599	LYS	2.2
1	D	604	GLY	2.2
1	A	489	LEU	2.2
1	C	601	VAL	2.2
1	D	623	GLU	2.2
1	D	606	VAL	2.1
1	A	510	ILE	2.1
1	D	570	LYS	2.1
1	B	289	GLN	2.1
1	A	609	ARG	2.1
1	B	394	ARG	2.1
1	A	553	GLU	2.1
1	A	623	GLU	2.1
1	B	259	GLU	2.1
1	A	550	SER	2.0
1	A	626	GLU	2.0
1	A	638	LYS	2.0
1	C	315	ASP	2.0
1	C	557	GLU	2.0
1	D	638	LYS	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 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,

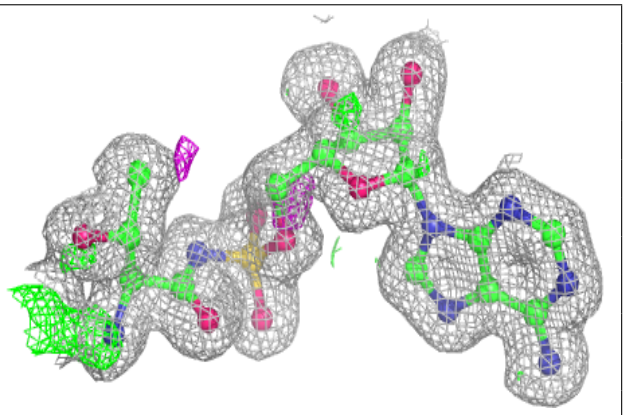
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

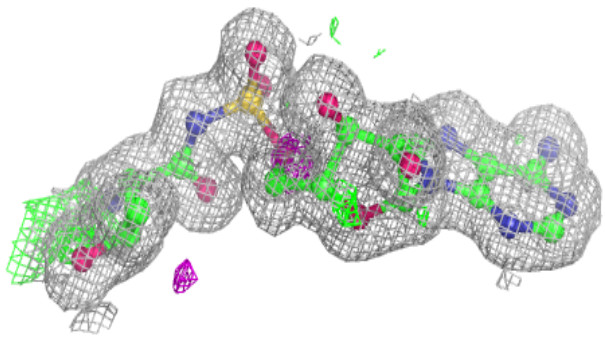
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	TSB	C	4002	30/30	0.96	0.10	11,13,14,14	0
3	TSB	B	3002	30/30	0.97	0.08	10,11,13,14	0
3	TSB	A	2002	30/30	0.97	0.10	11,13,14,14	0
3	TSB	D	5002	30/30	0.97	0.11	12,14,15,15	0
2	ZN	A	1	1/1	0.99	0.07	12,12,12,12	0
2	ZN	B	1	1/1	0.99	0.05	10,10,10,10	0
2	ZN	C	1	1/1	1.00	0.10	13,13,13,13	0
2	ZN	D	1	1/1	1.00	0.10	13,13,13,13	0

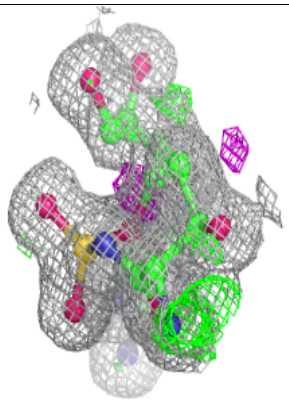
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around TSB C 4002:

2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative)
and green (positive)

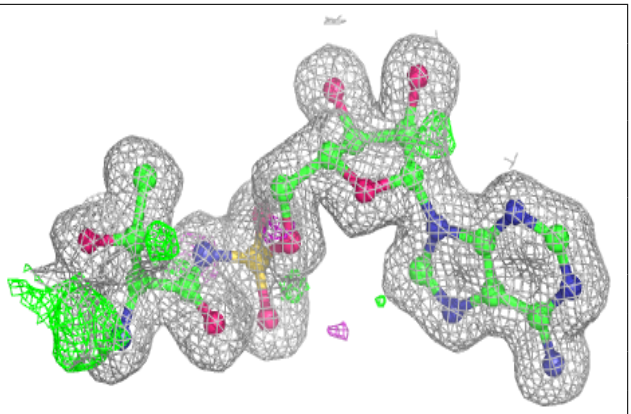


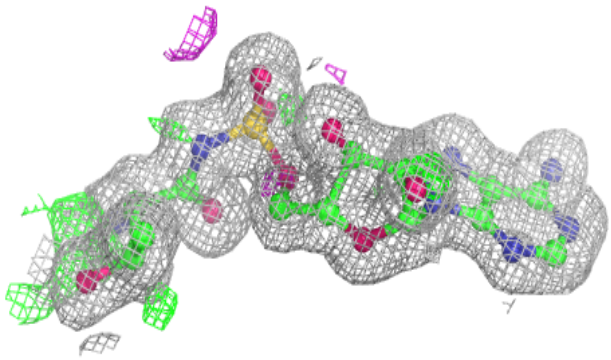


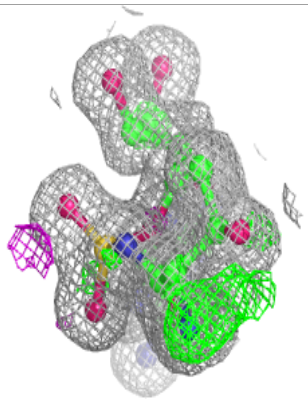


Electron density around TSB B 3002:

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

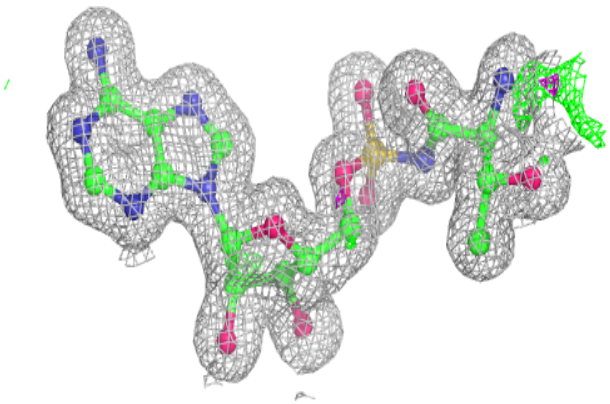


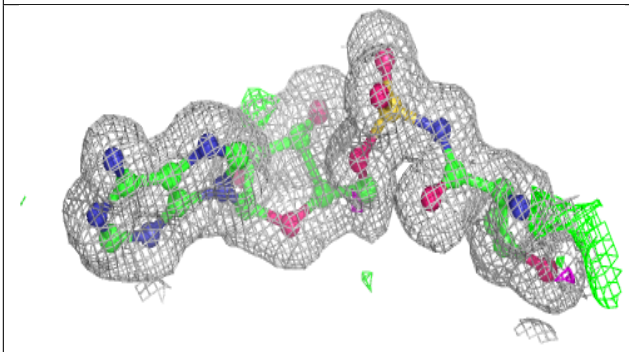


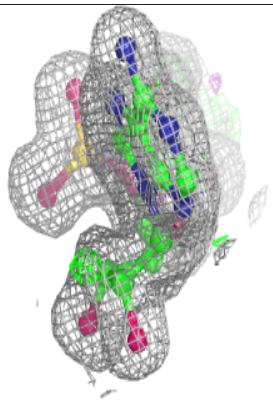


Electron density around TSB A 2002:

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

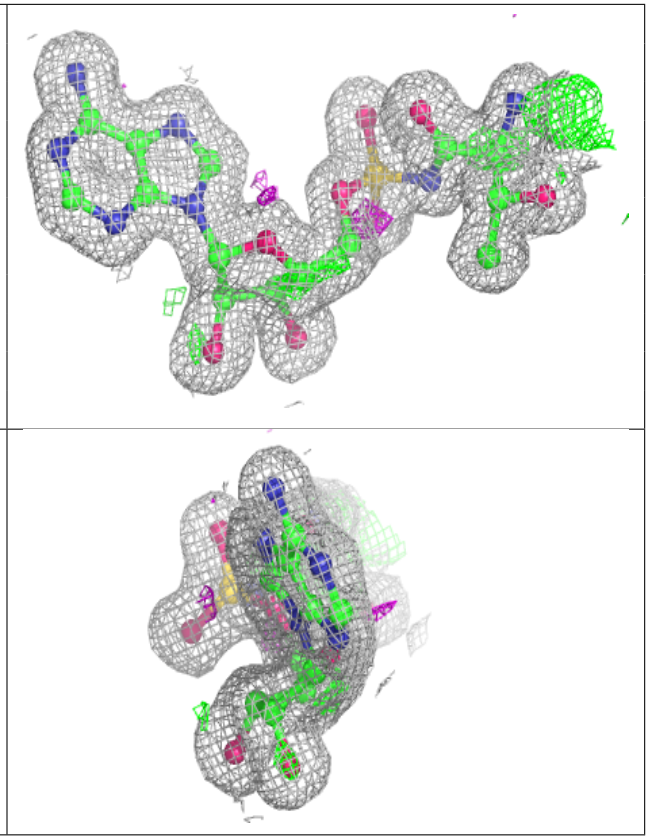






Electron density around TSB D 5002:

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.