Full wwPDB X-ray Structure Validation Report (i)

Jan 3, 2024 – 12:00 am GMT

PDB ID	:	5AO2
Title	:	Crystal structure of human SAMHD1 (amino acid residues 115-583) R164A variant bound to dGTP
Authors	:	Schwefel, D.; Taylor, I.A.
Deposited on	:	2015-09-09
Resolution	:	2.97 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:

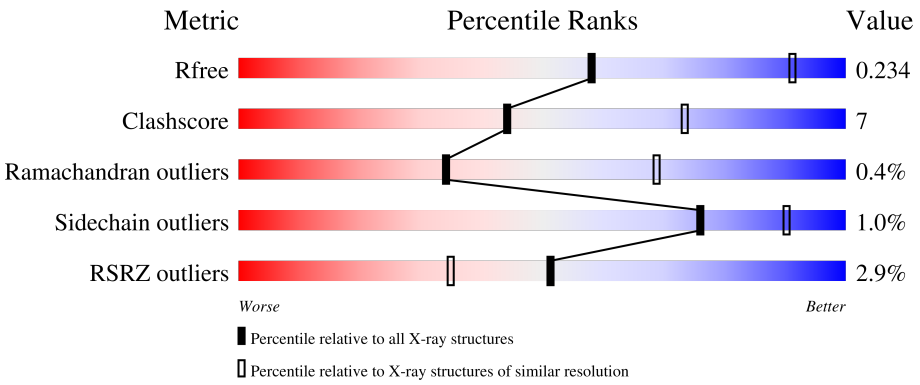
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
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 2.97 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	3104 (3.00-2.92)
Clashscore	141614	3462 (3.00-2.92)
Ramachandran outliers	138981	3340 (3.00-2.92)
Sidechain outliers	138945	3343 (3.00-2.92)
RSRZ outliers	127900	2986 (3.00-2.92)

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	491	2%, 75%, 13%, 11%
1	B	491	2%, 72%, 13%, •, 15%
1	C	491	3%, 74%, 14%, •, 11%
1	D	491	3%, 73%, 12%, 15%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 13606 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 DEOXYNUCLEOSIDE TRIPHOSPHATE TRIPHOSPHOHYDROLASE SAMHD1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	436	3433	2202	594	618	19	0	0	0
1	B	419	3317	2130	568	601	18	0	0	0
1	C	438	3452	2213	593	627	19	0	0	1
1	D	419	3272	2099	560	595	18	0	0	0

There are 92 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	93	MET	-	expression tag	UNP Q9Y3Z3
A	94	ALA	-	expression tag	UNP Q9Y3Z3
A	95	SER	-	expression tag	UNP Q9Y3Z3
A	96	TRP	-	expression tag	UNP Q9Y3Z3
A	97	SER	-	expression tag	UNP Q9Y3Z3
A	98	HIS	-	expression tag	UNP Q9Y3Z3
A	99	PRO	-	expression tag	UNP Q9Y3Z3
A	100	GLN	-	expression tag	UNP Q9Y3Z3
A	101	PHE	-	expression tag	UNP Q9Y3Z3
A	102	GLU	-	expression tag	UNP Q9Y3Z3
A	103	LYS	-	expression tag	UNP Q9Y3Z3
A	104	GLY	-	expression tag	UNP Q9Y3Z3
A	105	ALA	-	expression tag	UNP Q9Y3Z3
A	106	LEU	-	expression tag	UNP Q9Y3Z3
A	107	GLU	-	expression tag	UNP Q9Y3Z3
A	108	VAL	-	expression tag	UNP Q9Y3Z3
A	109	LEU	-	expression tag	UNP Q9Y3Z3
A	110	PHE	-	expression tag	UNP Q9Y3Z3
A	111	GLN	-	expression tag	UNP Q9Y3Z3
A	112	GLY	-	expression tag	UNP Q9Y3Z3

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Chain	Residue	Modelled	Actual	Comment	Reference
A	113	PRO	-	expression tag	UNP Q9Y3Z3
A	114	GLY	-	expression tag	UNP Q9Y3Z3
A	164	ALA	ARG	engineered mutation	UNP Q9Y3Z3
B	93	MET	-	expression tag	UNP Q9Y3Z3
B	94	ALA	-	expression tag	UNP Q9Y3Z3
B	95	SER	-	expression tag	UNP Q9Y3Z3
B	96	TRP	-	expression tag	UNP Q9Y3Z3
B	97	SER	-	expression tag	UNP Q9Y3Z3
B	98	HIS	-	expression tag	UNP Q9Y3Z3
B	99	PRO	-	expression tag	UNP Q9Y3Z3
B	100	GLN	-	expression tag	UNP Q9Y3Z3
B	101	PHE	-	expression tag	UNP Q9Y3Z3
B	102	GLU	-	expression tag	UNP Q9Y3Z3
B	103	LYS	-	expression tag	UNP Q9Y3Z3
B	104	GLY	-	expression tag	UNP Q9Y3Z3
B	105	ALA	-	expression tag	UNP Q9Y3Z3
B	106	LEU	-	expression tag	UNP Q9Y3Z3
B	107	GLU	-	expression tag	UNP Q9Y3Z3
B	108	VAL	-	expression tag	UNP Q9Y3Z3
B	109	LEU	-	expression tag	UNP Q9Y3Z3
B	110	PHE	-	expression tag	UNP Q9Y3Z3
B	111	GLN	-	expression tag	UNP Q9Y3Z3
B	112	GLY	-	expression tag	UNP Q9Y3Z3
B	113	PRO	-	expression tag	UNP Q9Y3Z3
B	114	GLY	-	expression tag	UNP Q9Y3Z3
B	164	ALA	ARG	engineered mutation	UNP Q9Y3Z3
C	93	MET	-	expression tag	UNP Q9Y3Z3
C	94	ALA	-	expression tag	UNP Q9Y3Z3
C	95	SER	-	expression tag	UNP Q9Y3Z3
C	96	TRP	-	expression tag	UNP Q9Y3Z3
C	97	SER	-	expression tag	UNP Q9Y3Z3
C	98	HIS	-	expression tag	UNP Q9Y3Z3
C	99	PRO	-	expression tag	UNP Q9Y3Z3
C	100	GLN	-	expression tag	UNP Q9Y3Z3
C	101	PHE	-	expression tag	UNP Q9Y3Z3
C	102	GLU	-	expression tag	UNP Q9Y3Z3
C	103	LYS	-	expression tag	UNP Q9Y3Z3
C	104	GLY	-	expression tag	UNP Q9Y3Z3
C	105	ALA	-	expression tag	UNP Q9Y3Z3
C	106	LEU	-	expression tag	UNP Q9Y3Z3
C	107	GLU	-	expression tag	UNP Q9Y3Z3
C	108	VAL	-	expression tag	UNP Q9Y3Z3

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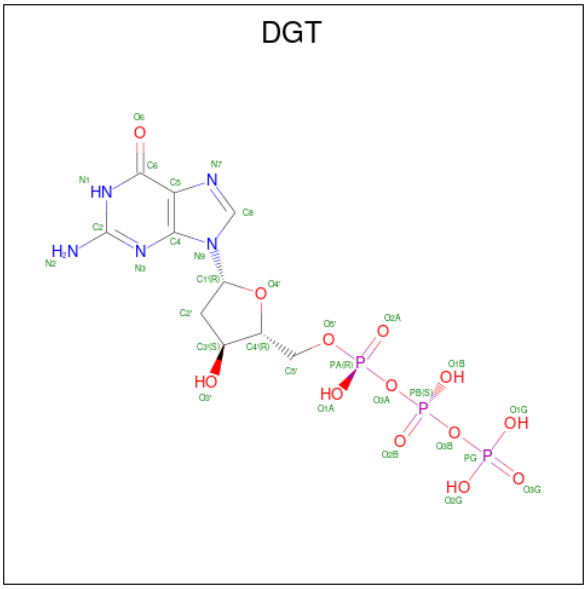
Chain	Residue	Modelled	Actual	Comment	Reference
C	109	LEU	-	expression tag	UNP Q9Y3Z3
C	110	PHE	-	expression tag	UNP Q9Y3Z3
C	111	GLN	-	expression tag	UNP Q9Y3Z3
C	112	GLY	-	expression tag	UNP Q9Y3Z3
C	113	PRO	-	expression tag	UNP Q9Y3Z3
C	114	GLY	-	expression tag	UNP Q9Y3Z3
C	164	ALA	ARG	engineered mutation	UNP Q9Y3Z3
D	93	MET	-	expression tag	UNP Q9Y3Z3
D	94	ALA	-	expression tag	UNP Q9Y3Z3
D	95	SER	-	expression tag	UNP Q9Y3Z3
D	96	TRP	-	expression tag	UNP Q9Y3Z3
D	97	SER	-	expression tag	UNP Q9Y3Z3
D	98	HIS	-	expression tag	UNP Q9Y3Z3
D	99	PRO	-	expression tag	UNP Q9Y3Z3
D	100	GLN	-	expression tag	UNP Q9Y3Z3
D	101	PHE	-	expression tag	UNP Q9Y3Z3
D	102	GLU	-	expression tag	UNP Q9Y3Z3
D	103	LYS	-	expression tag	UNP Q9Y3Z3
D	104	GLY	-	expression tag	UNP Q9Y3Z3
D	105	ALA	-	expression tag	UNP Q9Y3Z3
D	106	LEU	-	expression tag	UNP Q9Y3Z3
D	107	GLU	-	expression tag	UNP Q9Y3Z3
D	108	VAL	-	expression tag	UNP Q9Y3Z3
D	109	LEU	-	expression tag	UNP Q9Y3Z3
D	110	PHE	-	expression tag	UNP Q9Y3Z3
D	111	GLN	-	expression tag	UNP Q9Y3Z3
D	112	GLY	-	expression tag	UNP Q9Y3Z3
D	113	PRO	-	expression tag	UNP Q9Y3Z3
D	114	GLY	-	expression tag	UNP Q9Y3Z3
D	164	ALA	ARG	engineered mutation	UNP Q9Y3Z3

- Molecule 2 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Fe		
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 2'-DEOXYGUANOSINE-5'-TRIPHOSPHATE (three-letter code: DGT) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



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

- Molecule 4 is water.

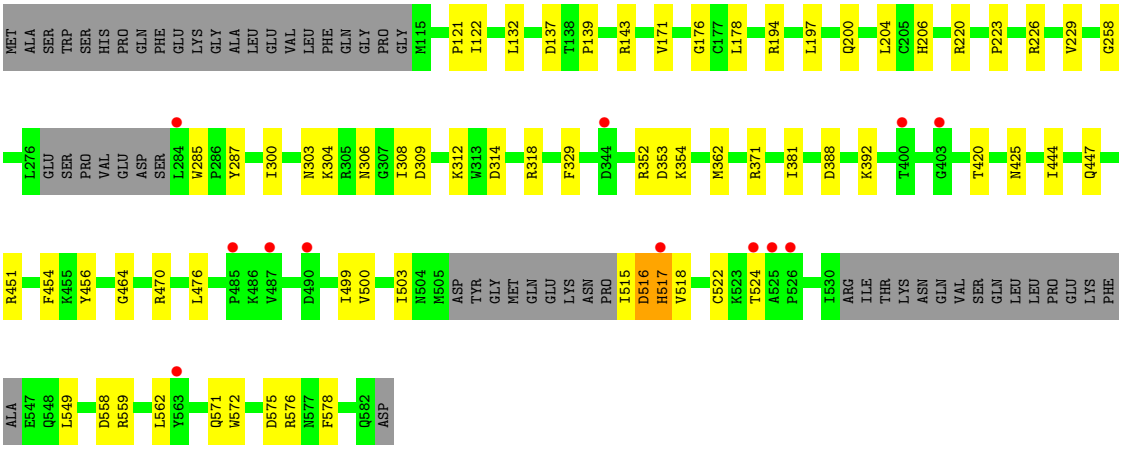
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	O	0	0
			1	1		
4	B	2	Total	O	0	0
			2	2		
4	C	1	Total	O	0	0
			1	1		

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.

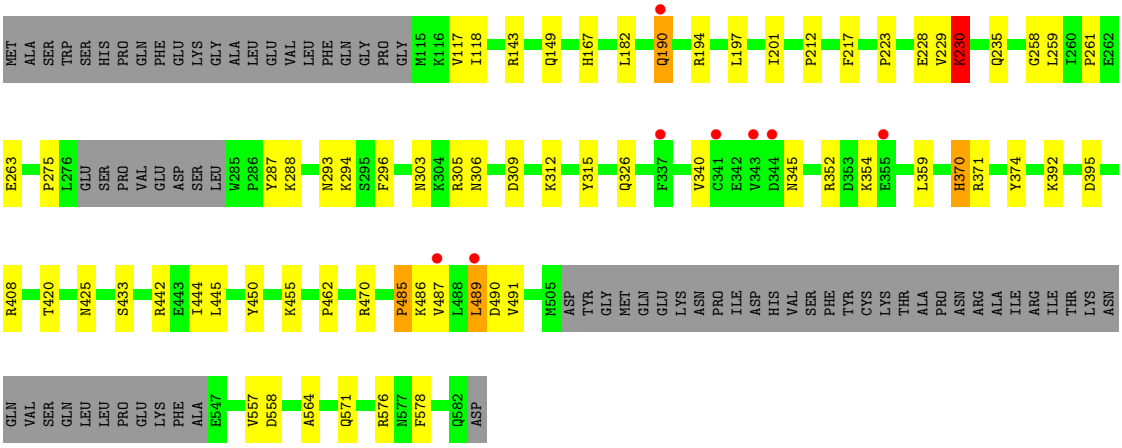
• Molecule 1: DEOXYNUCLEOSIDE TRIPHOSPHATE TRIPHOSPHOHYDROLASE SAMHD1

Chain A:



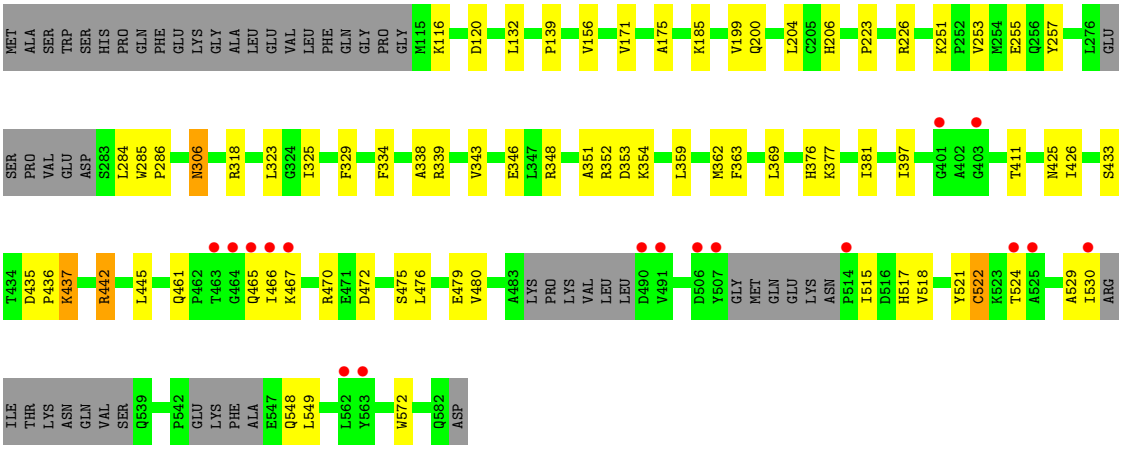
• Molecule 1: DEOXYNUCLEOSIDE TRIPHOSPHATE TRIPHOSPHOHYDROLASE SAMHD1

Chain B:



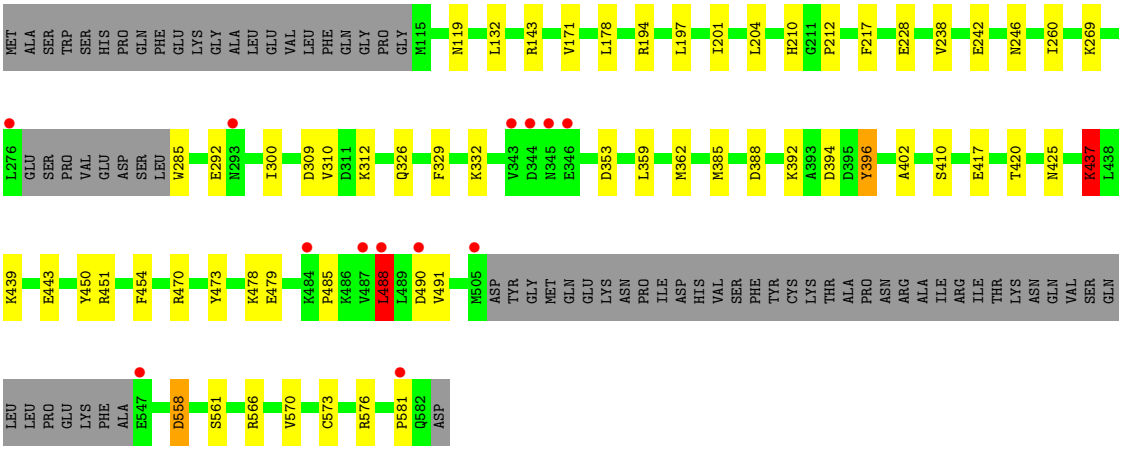
• Molecule 1: DEOXYNUCLEOSIDE TRIPHOSPHATE TRIPHOSPHOHYDROLASE SAMHD1

Chain C:



• Molecule 1: DEOXYNUCLEOSIDE TRIPHOSPHATE TRIPHOSPHOHYDROLASE SAMHD1

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	69.71Å 199.32Å 81.80Å 90.00° 100.79° 90.00°	Depositor
Resolution (Å)	29.56 – 2.97 29.56 – 2.97	Depositor EDS
% Data completeness (in resolution range)	96.0 (29.56-2.97) 96.0 (29.56-2.97)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.15 (at 2.95Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.156 , 0.234 0.159 , 0.234	Depositor DCC
R_{free} test set	2197 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å^2)	64.0	Xtriage
Anisotropy	0.303	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 81.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	13606	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 6.55% 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: FE, DGT

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	0.50	0/3515	0.69	2/4758 (0.0%)
1	B	0.52	0/3397	0.73	3/4596 (0.1%)
1	C	0.48	0/3535	0.70	1/4783 (0.0%)
1	D	0.54	2/3352 (0.1%)	0.73	3/4539 (0.1%)
All	All	0.51	2/13799 (0.0%)	0.71	9/18676 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	2
1	C	0	1
All	All	0	3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	573	CYS	CB-SG	-8.11	1.68	1.82
1	D	396	TYR	CE1-CZ	-8.11	1.28	1.38

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	437	LYS	CD-CE-NZ	-8.66	91.78	111.70
1	B	489	LEU	CA-CB-CG	8.60	135.09	115.30
1	D	488	LEU	CA-CB-CG	7.40	132.33	115.30
1	A	516	ASP	N-CA-C	-5.83	95.26	111.00
1	C	442	ARG	NE-CZ-NH2	5.77	123.19	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	516	ASP	CB-CG-OD1	5.48	123.23	118.30
1	B	489	LEU	CB-CG-CD2	-5.29	102.02	111.00
1	B	486	LYS	C-N-CA	5.27	134.88	121.70
1	D	353	ASP	C-N-CA	-5.12	108.89	121.70

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	230	LYS	Peptide
1	B	485	PRO	Peptide
1	C	284	LEU	Peptide

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	3433	0	3290	49	0
1	B	3317	0	3192	48	0
1	C	3452	0	3291	49	0
1	D	3272	0	3112	36	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	31	0	12	3	0
3	B	31	0	12	1	0
3	C	62	0	24	2	0
4	A	1	0	0	0	0
4	B	2	0	0	0	0
4	C	1	0	0	0	0
All	All	13606	0	12933	176	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 7.

All (176) 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:306:ASN:HB2	1:A:517:HIS:ND1	1.80	0.95
1:C:185:LYS:NZ	1:C:338:ALA:O	2.07	0.86
1:C:433:SER:OG	1:C:442:ARG:NH1	2.09	0.86
1:A:558:ASP:CA	1:A:559:ARG:N	2.39	0.85
1:B:370:HIS:HD1	1:B:374:TYR:HD2	1.32	0.76
1:B:190:GLN:HG2	1:B:294:LYS:HE3	1.68	0.76
1:A:517:HIS:CE1	1:A:518:VAL:HG23	2.23	0.74
1:B:485:PRO:HA	1:B:571:GLN:HG2	1.71	0.71
1:B:489:LEU:HD11	1:B:564:ALA:HB2	1.73	0.70
1:C:352:ARG:NE	1:C:354:LYS:HE2	2.07	0.70
1:C:515:ILE:HD12	1:C:515:ILE:H	1.57	0.69
1:D:479:GLU:OE1	1:D:576:ARG:NH1	2.26	0.68
1:B:228:GLU:HG2	1:B:229:VAL:HG23	1.76	0.67
1:B:326:GLN:OE1	1:D:326:GLN:NE2	2.26	0.66
1:B:194:ARG:NH2	1:B:263:GLU:OE1	2.29	0.65
3:C:1583:DGT:H5'	1:D:451:ARG:HH21	1.60	0.64
1:C:306:ASN:ND2	1:C:517:HIS:O	2.31	0.64
1:A:303:ASN:ND2	1:A:306:ASN:OD1	2.30	0.64
1:B:303:ASN:OD1	1:B:305:ARG:HG2	1.99	0.62
1:D:285:TRP:CG	1:D:292:GLU:HG3	2.35	0.61
1:B:370:HIS:ND1	1:B:374:TYR:HD2	1.98	0.61
1:B:305:ARG:HG3	1:B:306:ASN:N	2.16	0.60
1:C:522:CYS:HG	1:C:524:THR:HG1	1.49	0.60
1:A:143:ARG:HD2	1:A:420:THR:HA	1.83	0.60
1:D:285:TRP:CD2	1:D:292:GLU:HG3	2.36	0.60
1:C:522:CYS:HB2	1:C:530:ILE:HD11	1.84	0.59
1:B:190:GLN:HG2	1:B:190:GLN:O	2.02	0.59
1:A:515:ILE:N	1:A:517:HIS:CD2	2.71	0.59
1:B:230:LYS:HD2	1:B:235:GLN:HE22	1.68	0.58
1:C:352:ARG:HE	1:C:354:LYS:HE2	1.69	0.57
1:C:343:VAL:HG22	1:C:529:ALA:HB2	1.86	0.57
1:C:339:ARG:HD3	1:C:521:TYR:CZ	2.40	0.57
1:C:175:ALA:HB1	1:C:199:VAL:HG12	1.86	0.57
1:C:353:ASP:OD1	1:C:353:ASP:N	2.38	0.56
1:D:132:LEU:HD22	1:D:204:LEU:HD22	1.86	0.56
1:A:352:ARG:NH2	1:A:354:LYS:HE2	2.19	0.56
1:B:305:ARG:HG3	1:B:306:ASN:H	1.71	0.55
1:A:515:ILE:N	1:A:517:HIS:HD2	2.03	0.55
1:A:425:ASN:ND2	1:B:425:ASN:OD1	2.40	0.54
1:B:223:PRO:HG3	1:B:470:ARG:HG2	1.90	0.54
1:A:476:LEU:HB3	1:A:500:VAL:HG11	1.89	0.54
1:D:143:ARG:HD2	1:D:420:THR:HA	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:352:ARG:CZ	1:A:354:LYS:HE2	2.37	0.53
1:B:489:LEU:HD12	1:B:490:ASP:H	1.74	0.53
1:D:242:GLU:OE1	1:D:246:ASN:ND2	2.40	0.53
1:B:190:GLN:CG	1:B:294:LYS:HE3	2.39	0.52
1:B:352:ARG:HG3	1:B:354:LYS:H	1.75	0.52
1:A:571:GLN:NE2	1:A:575:ASP:OD1	2.43	0.51
1:B:197:LEU:O	1:B:201:ILE:HG13	2.09	0.51
1:C:354:LYS:N	1:C:354:LYS:HD3	2.24	0.51
1:A:503:ILE:HB	1:A:549:LEU:HB2	1.92	0.51
1:D:385:MET:HG2	1:D:454:PHE:CE2	2.46	0.50
1:A:308:ILE:CD1	1:A:517:HIS:HE1	2.24	0.50
1:A:353:ASP:N	1:A:353:ASP:OD1	2.45	0.50
1:A:381:ILE:HD11	1:A:456:TYR:HB2	1.94	0.50
1:B:558:ASP:N	1:B:558:ASP:OD1	2.44	0.50
1:A:314:ASP:OD2	1:A:318:ARG:NH1	2.45	0.50
1:A:522:CYS:HG	1:A:524:THR:HG1	1.60	0.50
1:A:572:TRP:O	1:A:576:ARG:HB2	2.11	0.50
1:B:190:GLN:HE22	1:B:293:ASN:HD21	1.59	0.49
1:A:306:ASN:CB	1:A:517:HIS:ND1	2.66	0.49
1:C:397:ILE:HG21	1:C:426:ILE:HD11	1.94	0.49
1:D:402:ALA:HA	1:D:417:GLU:OE2	2.12	0.49
1:A:516:ASP:O	1:A:517:HIS:HB3	2.11	0.49
1:A:220:ARG:NH2	1:A:500:VAL:O	2.47	0.48
1:C:433:SER:O	1:C:442:ARG:NH1	2.45	0.48
1:C:363:PHE:CE2	1:C:515:ILE:HG13	2.49	0.48
1:D:470:ARG:HG3	1:D:473:TYR:CE2	2.48	0.48
1:A:194:ARG:HD2	1:A:258:GLY:O	2.13	0.48
1:A:451:ARG:HH21	3:A:1583:DGT:C5'	2.27	0.48
1:B:229:VAL:HG13	1:B:230:LYS:HG2	1.95	0.48
1:D:194:ARG:NE	1:D:260:ILE:HD12	2.28	0.48
1:C:461:GLN:HG2	1:C:549:LEU:HD23	1.96	0.48
1:D:396:TYR:CG	1:D:437:LYS:HG2	2.49	0.48
1:B:190:GLN:HG2	1:B:294:LYS:CE	2.41	0.48
1:B:489:LEU:HD21	1:B:564:ALA:HB1	1.95	0.48
1:C:156:VAL:HG11	1:C:376:HIS:CE1	2.49	0.48
1:C:132:LEU:HD22	1:C:204:LEU:HD22	1.96	0.47
1:C:323:LEU:O	1:D:119:ASN:ND2	2.47	0.47
1:C:425:ASN:OD1	1:D:425:ASN:ND2	2.47	0.47
1:A:329:PHE:CD1	1:A:362:MET:HB2	2.49	0.47
1:A:451:ARG:HH21	3:A:1583:DGT:H5'A	1.79	0.47
1:C:132:LEU:HD11	1:C:200:GLN:NE2	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:454:PHE:CE2	1:A:499:ILE:HG13	2.50	0.47
1:A:226:ARG:NE	1:A:229:VAL:HG21	2.30	0.46
1:A:388:ASP:O	1:A:392:LYS:HG3	2.16	0.46
1:D:329:PHE:HB2	1:D:362:MET:HA	1.97	0.46
1:C:435:ASP:OD1	1:C:437:LYS:HB2	2.16	0.46
1:A:197:LEU:HA	1:A:197:LEU:HD23	1.73	0.46
1:C:223:PRO:CG	1:C:470:ARG:HG2	2.46	0.46
1:C:226:ARG:NH1	1:C:411:THR:HA	2.31	0.46
1:B:489:LEU:HG	1:B:491:VAL:H	1.81	0.46
1:D:388:ASP:O	1:D:392:LYS:HG3	2.15	0.46
1:A:223:PRO:HG3	1:A:470:ARG:HG2	1.98	0.46
1:D:171:VAL:HG13	1:D:310:VAL:HG23	1.97	0.45
1:C:472:ASP:O	1:C:475:SER:OG	2.33	0.45
1:D:558:ASP:HB2	1:D:561:SER:H	1.81	0.45
1:B:194:ARG:HD2	1:B:258:GLY:O	2.17	0.45
1:D:228:GLU:H	1:D:228:GLU:CD	2.18	0.45
1:A:132:LEU:HD11	1:A:200:GLN:OE1	2.17	0.45
1:C:329:PHE:CD1	1:C:362:MET:HB2	2.51	0.45
1:D:490:ASP:OD1	1:D:491:VAL:N	2.50	0.45
1:A:516:ASP:O	1:A:517:HIS:CD2	2.70	0.45
1:D:309:ASP:OD1	1:D:312:LYS:HG2	2.17	0.45
1:B:395:ASP:OD1	1:B:408:ARG:NH2	2.49	0.44
1:D:178:LEU:HD23	1:D:300:ILE:HG23	1.99	0.44
1:A:476:LEU:CB	1:A:500:VAL:HG11	2.47	0.44
1:D:396:TYR:CD1	1:D:437:LYS:HG2	2.52	0.44
1:B:167:HIS:CE1	1:B:315:TYR:HB3	2.52	0.44
1:C:329:PHE:HB2	1:C:362:MET:HA	2.00	0.44
1:B:149:GLN:HE22	1:B:167:HIS:CG	2.35	0.44
1:C:346:GLU:OE1	1:C:348:ARG:NH2	2.51	0.44
1:C:171:VAL:HG21	1:C:206:HIS:CE1	2.53	0.44
1:C:479:GLU:HB3	1:C:572:TRP:HE1	1.82	0.44
1:D:394:ASP:OD2	1:D:410:SER:OG	2.27	0.44
1:C:522:CYS:CB	1:C:530:ILE:HD11	2.48	0.44
1:B:359:LEU:HD23	1:B:359:LEU:HA	1.91	0.43
1:C:120:ASP:OD1	1:C:318:ARG:NH2	2.50	0.43
1:C:251:LYS:O	1:C:255:GLU:HG3	2.18	0.43
1:C:359:LEU:HD11	1:C:518:VAL:HG11	2.00	0.43
1:D:197:LEU:HD23	1:D:197:LEU:HA	1.81	0.43
1:D:212:PRO:HD2	1:D:217:PHE:CD1	2.53	0.43
1:D:566:ARG:O	1:D:570:VAL:HG23	2.19	0.43
1:B:455:LYS:HG2	1:B:557:VAL:HG12	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:212:PRO:HD2	1:B:217:PHE:CD1	2.53	0.43
1:C:116:LYS:HE2	3:C:1583:DGT:O3B	2.18	0.43
1:B:143:ARG:HD2	1:B:420:THR:HA	2.00	0.43
1:B:433:SER:OG	1:B:442:ARG:NH1	2.50	0.43
1:D:439:LYS:O	1:D:443:GLU:HG2	2.18	0.43
1:D:485:PRO:HD2	1:D:488:LEU:HD23	2.01	0.43
1:C:435:ASP:OD1	1:C:436:PRO:HD2	2.18	0.43
1:C:521:TYR:HD1	1:C:522:CYS:O	2.01	0.43
1:A:176:GLY:HA3	1:A:200:GLN:HE21	1.84	0.42
1:B:395:ASP:HA	1:B:408:ARG:HE	1.83	0.42
1:A:204:LEU:HD12	1:A:204:LEU:HA	1.90	0.42
1:C:476:LEU:O	1:C:480:VAL:HG23	2.19	0.42
1:A:444:ILE:O	1:A:447:GLN:HB2	2.19	0.42
1:D:478:LYS:H	1:D:478:LYS:HG2	1.65	0.42
1:C:467:LYS:HB3	1:C:548:GLN:OE1	2.19	0.42
1:A:139:PRO:HD3	1:B:450:TYR:CE1	2.54	0.42
1:B:182:LEU:HD13	1:B:340:VAL:CG2	2.50	0.42
1:A:178:LEU:HD23	1:A:300:ILE:HG12	2.00	0.42
1:A:392:LYS:NZ	1:A:444:ILE:HD11	2.34	0.42
1:C:445:LEU:HD23	1:C:445:LEU:HA	1.85	0.42
1:B:576:ARG:HH11	1:B:576:ARG:HG2	1.85	0.42
1:D:197:LEU:O	1:D:201:ILE:HG13	2.19	0.42
1:B:392:LYS:HD2	1:B:444:ILE:HD11	2.02	0.42
1:B:259:LEU:O	1:B:261:PRO:HD3	2.20	0.42
1:C:359:LEU:HD23	1:C:359:LEU:HA	1.67	0.41
1:C:466:ILE:HD12	1:C:466:ILE:N	2.34	0.41
1:A:121:PRO:HG2	1:A:122:ILE:HD12	2.01	0.41
1:D:329:PHE:CD1	1:D:362:MET:HB2	2.54	0.41
1:A:171:VAL:HG21	1:A:206:HIS:CE1	2.55	0.41
1:A:285:TRP:CZ2	1:A:287:TYR:HD2	2.38	0.41
1:A:576:ARG:HD2	1:A:578:PHE:CZ	2.55	0.41
1:A:137:ASP:CG	3:B:1583:DGT:HN2	2.24	0.41
1:B:182:LEU:HD23	1:B:182:LEU:HA	1.78	0.41
1:C:325:ILE:HD12	1:C:369:LEU:HD23	2.02	0.41
1:B:462:PRO:HA	1:B:578:PHE:CD1	2.56	0.41
1:C:377:LYS:O	1:C:381:ILE:HG13	2.21	0.41
1:B:371:ARG:O	1:B:371:ARG:HD3	2.21	0.41
1:B:294:LYS:HA	1:B:296:PHE:CE1	2.56	0.41
1:B:309:ASP:OD1	1:B:312:LYS:HG2	2.19	0.41
1:C:139:PRO:HD3	1:D:450:TYR:CE1	2.55	0.41
1:C:334:PHE:CZ	1:C:351:ALA:HB2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:359:LEU:HA	1:D:359:LEU:HD23	1.63	0.41
1:A:226:ARG:CZ	1:A:229:VAL:HG21	2.51	0.41
1:A:562:LEU:HD23	1:A:562:LEU:HA	1.83	0.40
3:A:1583:DGT:H2'	1:B:118:ILE:HG12	2.03	0.40
1:C:285:TRP:HA	1:C:286:PRO:HD3	1.93	0.40
1:D:238:VAL:HG13	1:D:269:LYS:HD3	2.03	0.40
1:A:516:ASP:O	1:A:517:HIS:CB	2.68	0.40
1:B:275:PRO:HD3	1:B:287:TYR:CE1	2.56	0.40
1:C:253:VAL:O	1:C:257:TYR:HD2	2.05	0.40
1:A:309:ASP:OD1	1:A:312:LYS:HG2	2.22	0.40
1:B:445:LEU:HD23	1:B:445:LEU:HA	1.83	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	426/491 (87%)	416 (98%)	7 (2%)	3 (1%)	22	56
1	B	413/491 (84%)	403 (98%)	7 (2%)	3 (1%)	22	56
1	C	426/491 (87%)	417 (98%)	9 (2%)	0	100	100
1	D	413/491 (84%)	405 (98%)	7 (2%)	1 (0%)	47	79
All	All	1678/1964 (85%)	1641 (98%)	30 (2%)	7 (0%)	34	69

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	487	VAL
1	A	464	GLY
1	A	517	HIS
1	B	230	LYS

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Mol	Chain	Res	Type
1	A	304	LYS
1	B	345	ASN
1	D	581	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.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	355/433 (82%)	354 (100%)	1 (0%)	92	97
1	B	344/433 (79%)	340 (99%)	4 (1%)	71	88
1	C	357/433 (82%)	353 (99%)	4 (1%)	73	89
1	D	336/433 (78%)	331 (98%)	5 (2%)	65	85
All	All	1392/1732 (80%)	1378 (99%)	14 (1%)	76	90

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	371	ARG
1	B	117	VAL
1	B	190	GLN
1	B	288	LYS
1	B	370	HIS
1	C	306	ASN
1	C	437	LYS
1	C	465	GLN
1	C	522	CYS
1	D	210	HIS
1	D	332	LYS
1	D	437	LYS
1	D	488	LEU
1	D	558	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	303	ASN
1	A	425	ASN
1	A	517	HIS
1	B	235	GLN
1	B	293	ASN
1	C	303	ASN
1	C	328	ASN
1	D	425	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	DGT	B	1583	-	26,33,33	1.84	7 (26%)	32,52,52	1.73	6 (18%)
3	DGT	A	1583	-	26,33,33	1.87	7 (26%)	32,52,52	1.91	6 (18%)
3	DGT	C	1584	-	26,33,33	1.81	8 (30%)	32,52,52	1.79	7 (21%)
3	DGT	C	1583	-	26,33,33	1.82	8 (30%)	32,52,52	1.75	6 (18%)

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	DGT	B	1583	-	-	2/18/34/34	0/3/3/3
3	DGT	A	1583	-	-	7/18/34/34	0/3/3/3
3	DGT	C	1584	-	-	3/18/34/34	0/3/3/3
3	DGT	C	1583	-	-	7/18/34/34	0/3/3/3

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1583	DGT	C2-N2	4.19	1.44	1.34
3	C	1584	DGT	C2-N2	3.99	1.43	1.34
3	A	1583	DGT	C8-N7	-3.88	1.28	1.35
3	B	1583	DGT	C2-N2	3.84	1.43	1.34
3	A	1583	DGT	C2-N2	3.84	1.43	1.34
3	B	1583	DGT	C8-N7	-3.81	1.28	1.35
3	C	1584	DGT	C8-N7	-3.66	1.28	1.35
3	C	1583	DGT	C8-N7	-3.57	1.28	1.35
3	A	1583	DGT	C5-C4	-3.36	1.34	1.43
3	B	1583	DGT	C5-C4	-3.17	1.35	1.43
3	C	1583	DGT	C3'-C4'	-3.03	1.44	1.53
3	C	1584	DGT	C3'-C4'	-2.96	1.44	1.53
3	C	1583	DGT	C5-C4	-2.94	1.35	1.43
3	A	1583	DGT	C3'-C4'	-2.90	1.45	1.53
3	C	1584	DGT	C5-C4	-2.85	1.35	1.43
3	B	1583	DGT	C1'-N9	-2.78	1.41	1.49
3	B	1583	DGT	C3'-C4'	-2.73	1.45	1.53
3	A	1583	DGT	C1'-N9	-2.67	1.41	1.49
3	A	1583	DGT	C4-N3	2.62	1.43	1.37
3	C	1584	DGT	C4-N3	2.61	1.43	1.37
3	B	1583	DGT	C4-N3	2.54	1.43	1.37
3	C	1583	DGT	C1'-N9	-2.54	1.41	1.49
3	C	1584	DGT	C1'-N9	-2.51	1.42	1.49
3	A	1583	DGT	O4'-C1'	2.51	1.48	1.42
3	C	1583	DGT	C4-N3	2.41	1.43	1.37
3	C	1584	DGT	O4'-C1'	2.38	1.47	1.42
3	B	1583	DGT	O4'-C1'	2.37	1.47	1.42
3	C	1583	DGT	O4'-C1'	2.30	1.47	1.42
3	C	1584	DGT	C2-N3	2.09	1.38	1.33
3	C	1583	DGT	C2-N3	2.07	1.38	1.33

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	1583	DGT	C8-N7-C5	6.65	115.65	102.99
3	B	1583	DGT	C8-N7-C5	6.47	115.31	102.99
3	C	1584	DGT	C8-N7-C5	6.28	114.95	102.99
3	C	1583	DGT	C8-N7-C5	6.06	114.54	102.99
3	A	1583	DGT	C5-C6-N1	3.99	121.00	113.95
3	B	1583	DGT	C5-C6-N1	3.48	120.10	113.95
3	C	1584	DGT	C5-C6-N1	3.43	120.01	113.95
3	C	1583	DGT	C5-C6-N1	3.27	119.73	113.95
3	B	1583	DGT	C2-N1-C6	-2.90	119.76	125.10
3	A	1583	DGT	C2-N1-C6	-2.85	119.85	125.10
3	A	1583	DGT	PA-O3A-PB	-2.79	123.26	132.83
3	C	1584	DGT	C2-N1-C6	-2.69	120.14	125.10
3	C	1583	DGT	C2-N1-C6	-2.68	120.16	125.10
3	C	1583	DGT	PB-O3B-PG	-2.52	124.17	132.83
3	C	1583	DGT	C2'-C3'-C4'	2.48	107.94	102.76
3	A	1583	DGT	PB-O3B-PG	-2.47	124.34	132.83
3	C	1584	DGT	O6-C6-N1	-2.42	117.79	120.65
3	B	1583	DGT	O6-C6-C5	-2.40	119.68	124.37
3	B	1583	DGT	C2'-C3'-C4'	2.30	107.55	102.76
3	C	1583	DGT	O6-C6-C5	-2.28	119.92	124.37
3	C	1584	DGT	C2'-C3'-C4'	2.17	107.28	102.76
3	A	1583	DGT	O6-C6-N1	-2.15	118.11	120.65
3	C	1584	DGT	PB-O3B-PG	-2.05	125.78	132.83
3	B	1583	DGT	PB-O3B-PG	-2.02	125.89	132.83
3	C	1584	DGT	C4'-O4'-C1'	-2.01	104.59	109.45

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1583	DGT	PB-O3B-PG-O1G
3	A	1583	DGT	PB-O3B-PG-O2G
3	A	1583	DGT	C5'-O5'-PA-O3A
3	A	1583	DGT	C5'-O5'-PA-O1A
3	C	1583	DGT	C5'-O5'-PA-O1A
3	C	1583	DGT	C5'-O5'-PA-O2A
3	C	1583	DGT	O4'-C4'-C5'-O5'
3	B	1583	DGT	O4'-C4'-C5'-O5'
3	B	1583	DGT	C3'-C4'-C5'-O5'
3	C	1584	DGT	O4'-C4'-C5'-O5'
3	C	1584	DGT	C3'-C4'-C5'-O5'

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Mol	Chain	Res	Type	Atoms
3	C	1583	DGT	C3'-C4'-C5'-O5'
3	A	1583	DGT	PB-O3A-PA-O5'
3	C	1583	DGT	PB-O3A-PA-O5'
3	A	1583	DGT	C5'-O5'-PA-O2A
3	C	1583	DGT	PB-O3B-PG-O3G
3	C	1583	DGT	C5'-O5'-PA-O3A
3	A	1583	DGT	O4'-C4'-C5'-O5'
3	C	1584	DGT	PB-O3A-PA-O2A

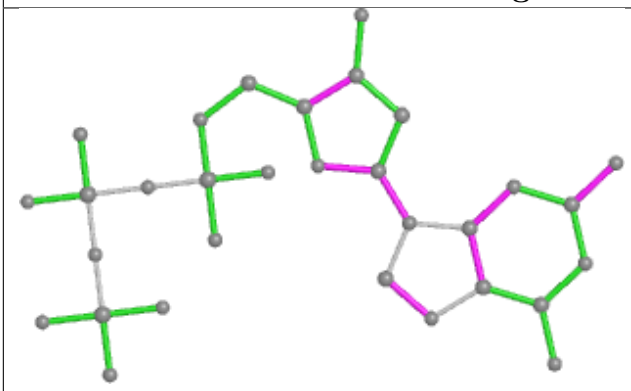
There are no ring outliers.

3 monomers are involved in 6 short contacts:

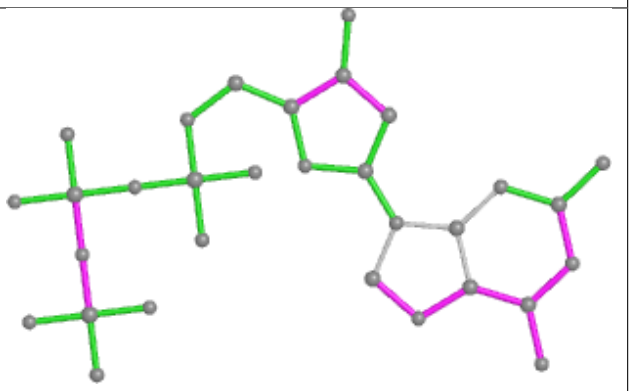
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1583	DGT	1	0
3	A	1583	DGT	3	0
3	C	1583	DGT	2	0

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.

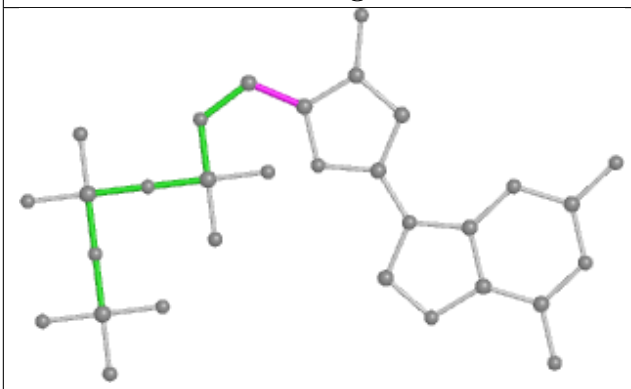
Ligand DGT B 1583



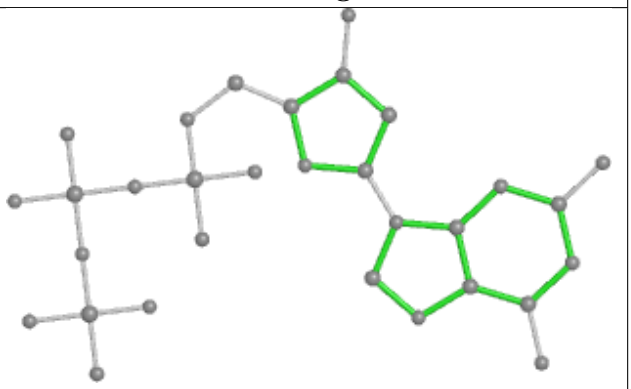
Bond lengths



Bond angles

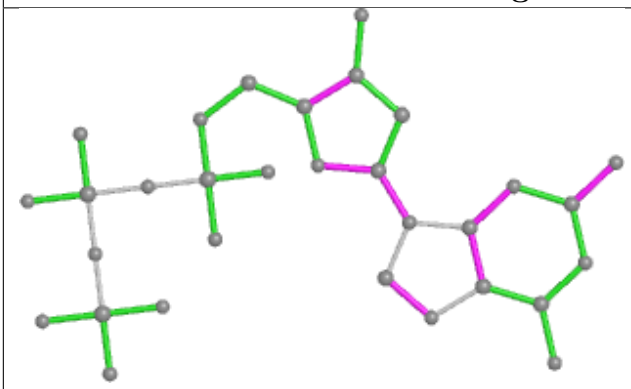


Torsions

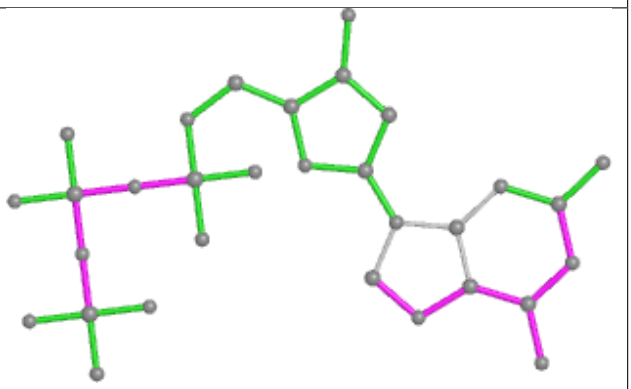


Rings

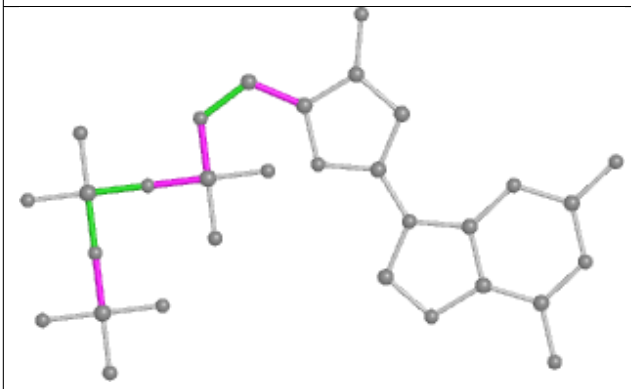
Ligand DGT A 1583



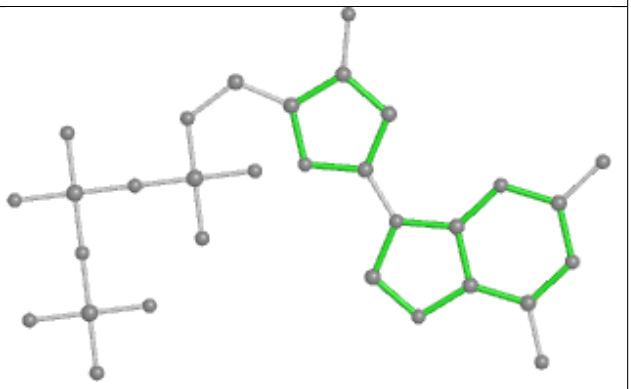
Bond lengths



Bond angles

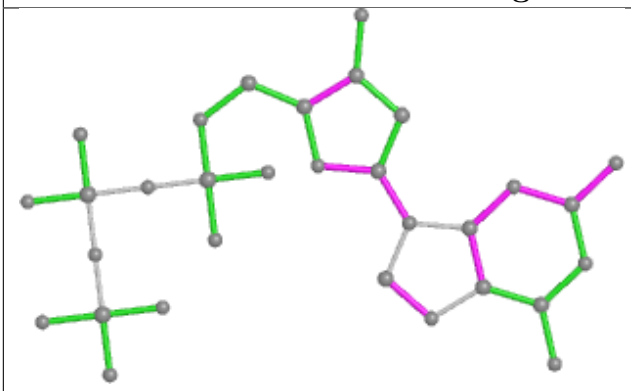


Torsions

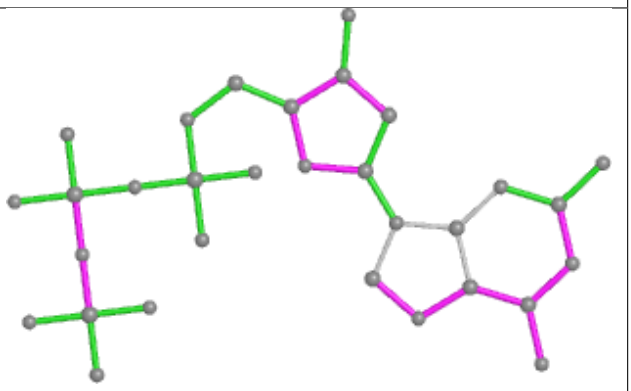


Rings

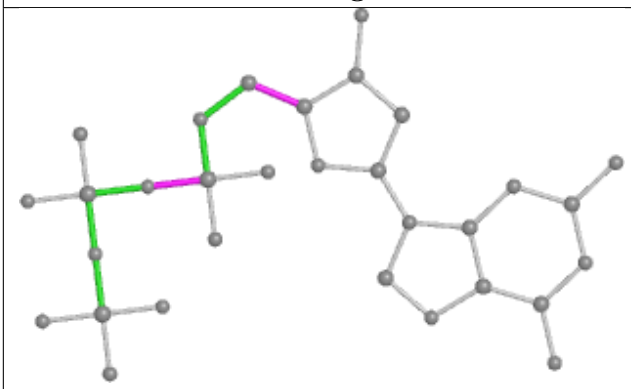
Ligand DGT C 1584



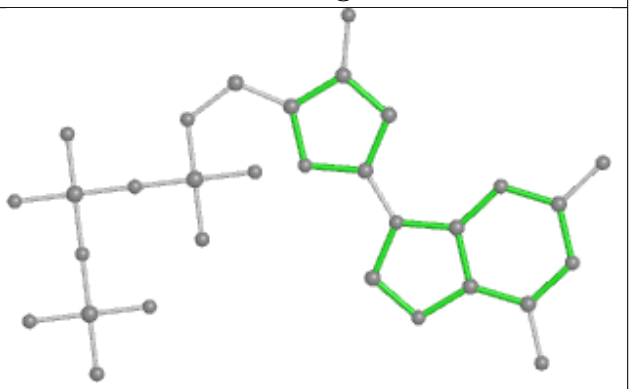
Bond lengths



Bond angles

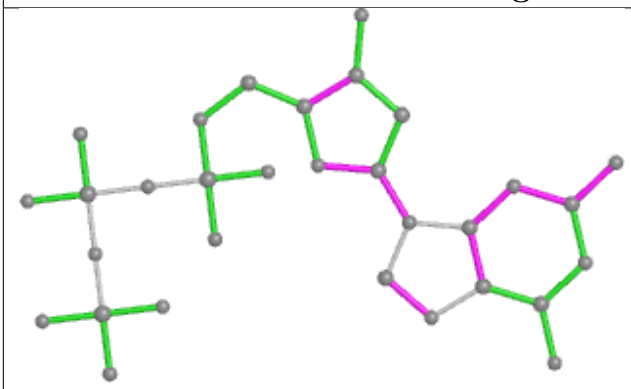


Torsions

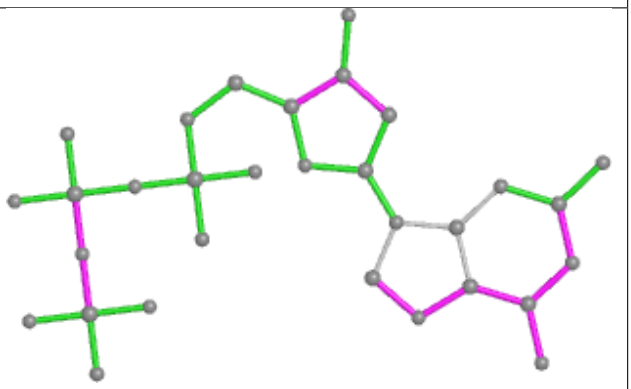


Rings

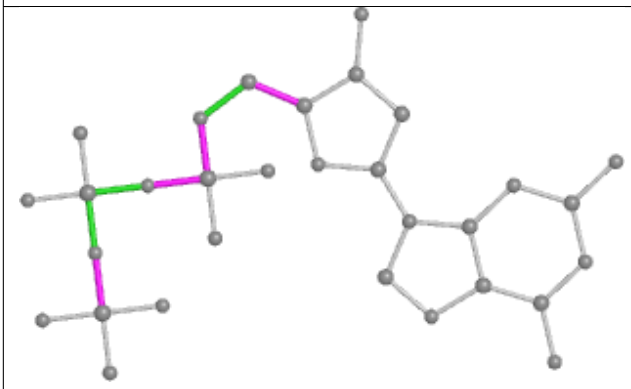
Ligand DGT C 1583



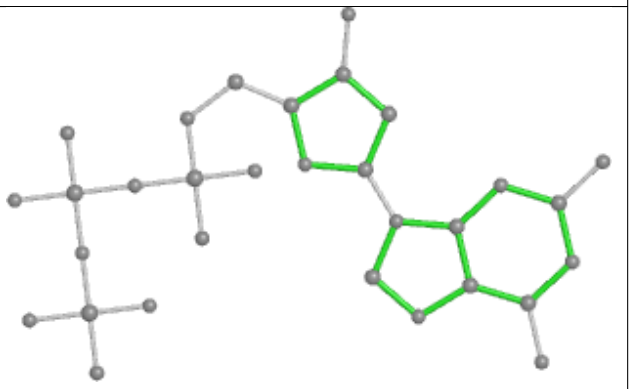
Bond lengths



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	436/491 (88%)	-0.20	12 (2%)	53	36	30, 58, 98, 176	1 (0%)
1	B	419/491 (85%)	-0.20	8 (1%)	66	49	29, 55, 108, 147	1 (0%)
1	C	438/491 (89%)	-0.02	17 (3%)	39	25	33, 65, 117, 151	0
1	D	419/491 (85%)	-0.10	13 (3%)	49	32	31, 65, 115, 163	0
All	All	1712/1964 (87%)	-0.13	50 (2%)	51	35	29, 61, 113, 176	2 (0%)

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	487	VAL	5.8
1	C	507	TYR	5.0
1	C	466	ILE	4.1
1	C	464	GLY	3.9
1	D	343	VAL	3.7
1	A	487	VAL	3.6
1	D	490	ASP	3.5
1	A	524	THR	3.5
1	B	487	VAL	3.4
1	D	345	ASN	3.4
1	B	343	VAL	3.3
1	C	491	VAL	3.3
1	A	517	HIS	3.3
1	A	284	LEU	3.2
1	C	524	THR	3.1
1	D	488	LEU	3.0
1	C	563	TYR	3.0
1	C	403	GLY	2.9
1	B	190	GLN	2.9
1	C	490	ASP	2.9
1	A	526	PRO	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	403	GLY	2.7
1	A	525	ALA	2.7
1	D	505	MET	2.7
1	C	463	THR	2.6
1	B	344	ASP	2.6
1	D	293	ASN	2.5
1	A	485	PRO	2.5
1	A	490	ASP	2.4
1	A	563	TYR	2.4
1	C	401	GLY	2.4
1	C	530	ILE	2.3
1	C	514	PRO	2.3
1	B	489	LEU	2.3
1	D	484	LYS	2.3
1	A	400	THR	2.3
1	D	344	ASP	2.3
1	D	581	PRO	2.3
1	B	355	GLU	2.2
1	D	276	LEU	2.1
1	C	465	GLN	2.1
1	A	344	ASP	2.1
1	D	346	GLU	2.1
1	C	506	ASP	2.1
1	B	341	CYS	2.1
1	C	525	ALA	2.1
1	B	337	PHE	2.1
1	C	562	LEU	2.0
1	D	547	GLU	2.0
1	C	467	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)

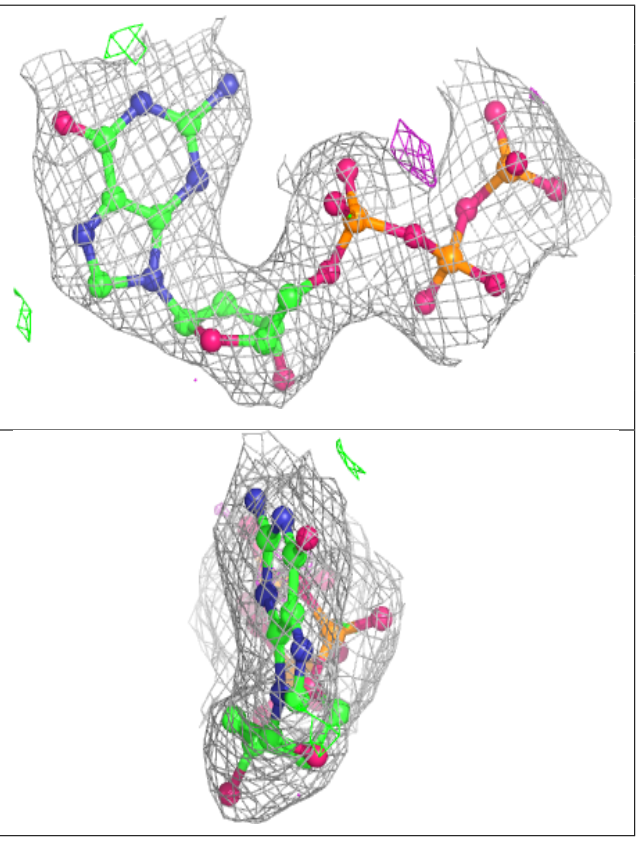
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	DGT	C	1583	31/31	0.91	0.18	47,81,125,186	0
3	DGT	C	1584	31/31	0.94	0.14	54,83,131,161	0
3	DGT	B	1583	31/31	0.95	0.15	41,76,129,146	0
3	DGT	A	1583	31/31	0.96	0.12	46,76,107,114	0
2	FE	D	1001	1/1	0.98	0.20	69,69,69,69	0
2	FE	C	1001	1/1	0.99	0.22	60,60,60,60	0
2	FE	A	1001	1/1	0.99	0.17	58,58,58,58	0
2	FE	B	1001	1/1	0.99	0.17	63,63,63,63	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.

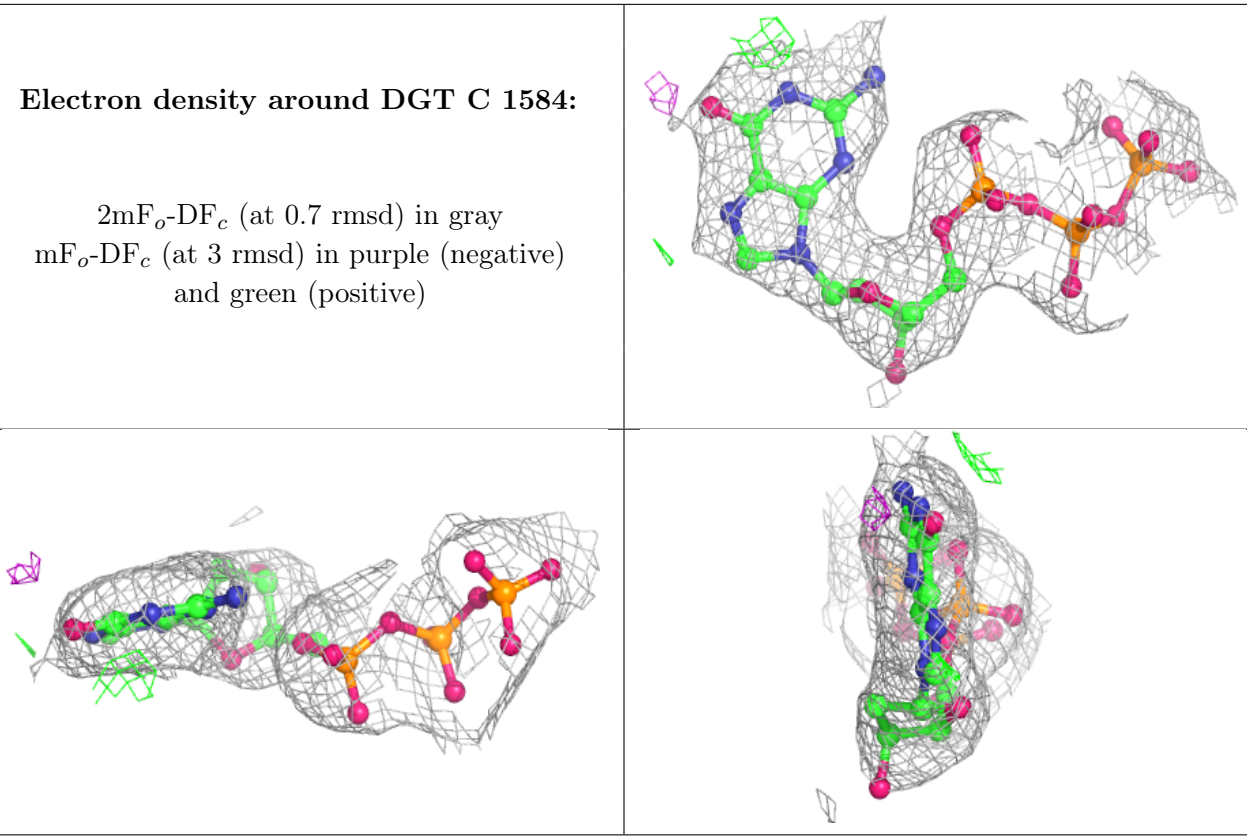
Electron density around DGT C 1583:

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



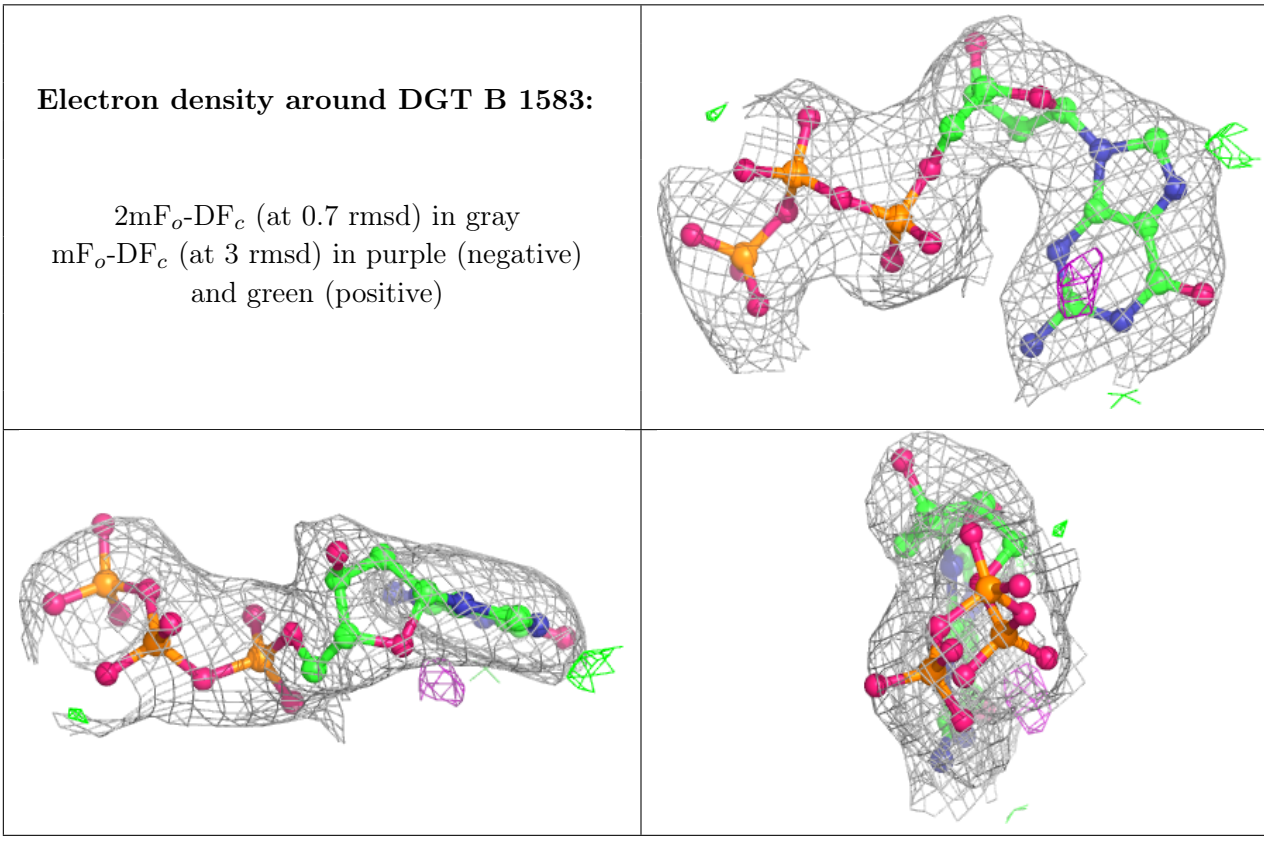
Electron density around DGT C 1584:

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



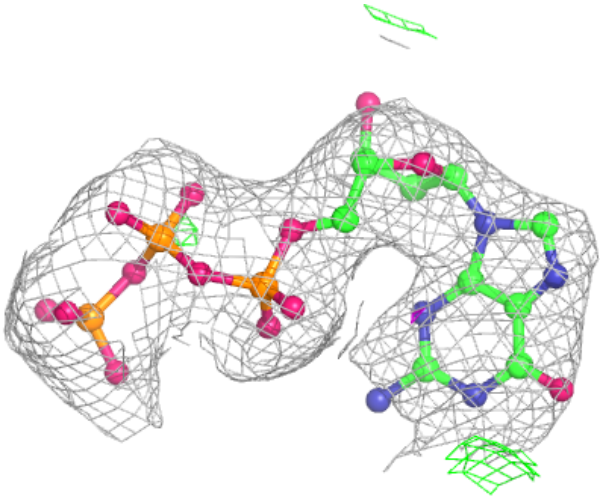
Electron density around DGT B 1583:

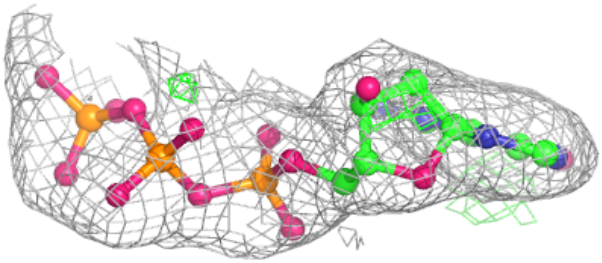
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

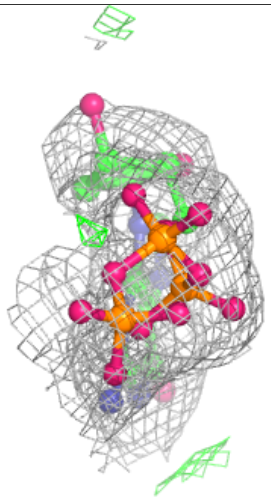


Electron density around DGT A 1583:

$2mF_o$-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.