



# Full wwPDB X-ray Structure Validation Report ⓘ

Sep 2, 2021 – 04:06 PM EDT

PDB ID : 7S0Z  
Title : Structures of TcdB in complex with R-Ras  
Authors : Zheng, L.; Rongsheng, J.; Peng, C.  
Deposited on : 2021-08-31  
Resolution : 2.34 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](mailto:validation@mail.wwpdb.org)

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) 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 : 2.23.1  
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) : 2.23.1

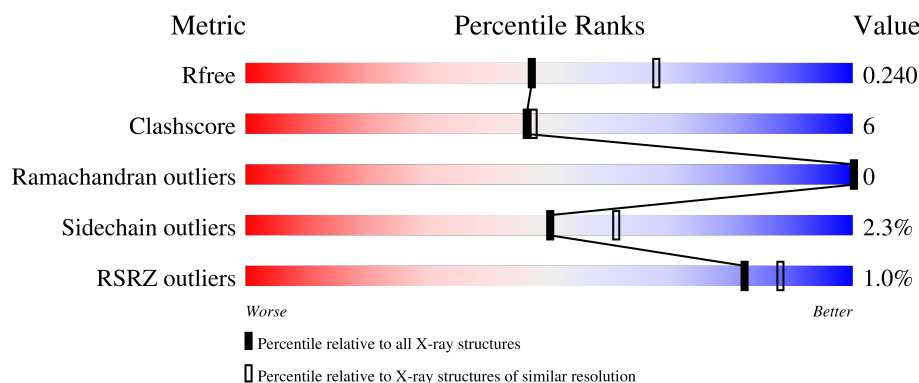
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.34 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	2096 (2.36-2.32)
Clashscore	141614	2193 (2.36-2.32)
Ramachandran outliers	138981	2159 (2.36-2.32)
Sidechain outliers	138945	2160 (2.36-2.32)
RSRZ outliers	127900	2067 (2.36-2.32)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	541	<div> <div>%</div> <div> <div></div> <div>85%</div> <div>14%</div> <div>.</div> </div> </div>
1	B	541	<div> <div>2%</div> <div> <div></div> <div>89%</div> <div>10%</div> <div>.</div> </div> </div>
2	C	179	<div> <div></div> <div> <div></div> <div>88%</div> <div>12%</div> <div>.</div> </div> </div>
2	D	179	<div> <div>%</div> <div> <div></div> <div>85%</div> <div>15%</div> <div>.</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	NH4	A	602	-	-	X	-
6	ACT	A	614	-	-	X	-
6	ACT	C	304	-	-	X	-
7	PEG	B	612	-	-	X	-
9	GLC	A	623	-	-	X	-
9	GLC	B	616	-	-	X	-

## 2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 12426 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 Toxin B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	541	Total	C	N	O	S	0	0	0
			4404	2808	713	873	10			
1	A	541	Total	C	N	O	S	0	0	0
			4402	2807	713	872	10			

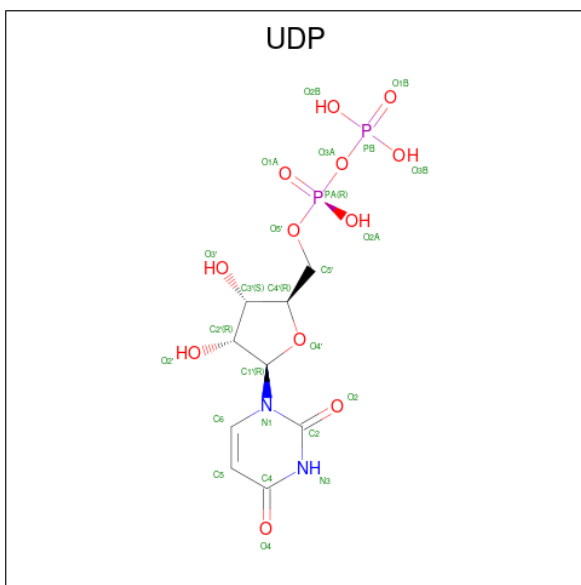
- Molecule 2 is a protein called Ras-related protein R-Ras.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	179	Total	C	N	O	S	0	1	0
			1414	892	245	274	3			
2	D	179	Total	C	N	O	S	0	0	0
			1402	885	243	271	3			

There are 2 discrepancies between the modelled and reference sequences:

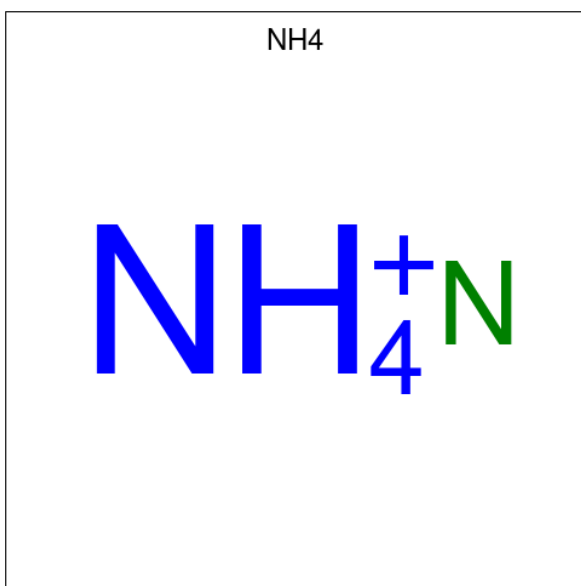
Chain	Residue	Modelled	Actual	Comment	Reference
C	61	ASN	THR	conflict	UNP P10301
D	61	ASN	THR	conflict	UNP P10301

- Molecule 3 is URIDINE-5'-DIPHOSPHATE (three-letter code: UDP) (formula: C<sub>9</sub>H<sub>14</sub>N<sub>2</sub>O<sub>12</sub>P<sub>2</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	B	1	Total 25	C 9	N 2	O 12	P 2	0	0
3	A	1	Total 25	C 9	N 2	O 12	P 2	0	0

- Molecule 4 is AMMONIUM ION (three-letter code: NH4) (formula:  $\text{H}_4\text{N}$ ).



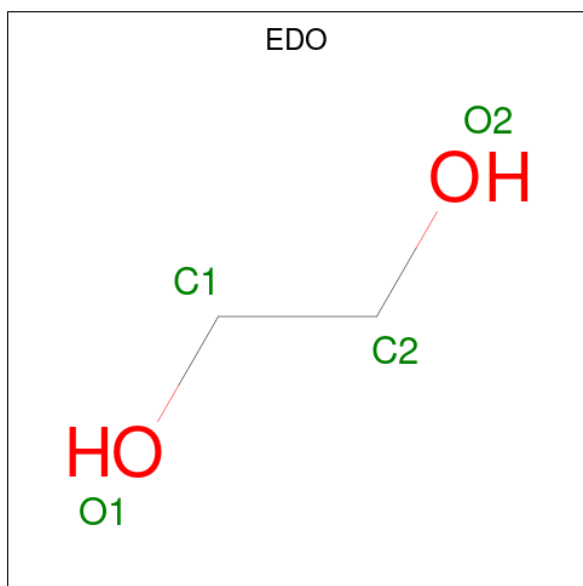
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total N 1 1	0	0
4	B	1	Total N 1 1	0	0

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

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula:  $C_2H_6O_2$ ).



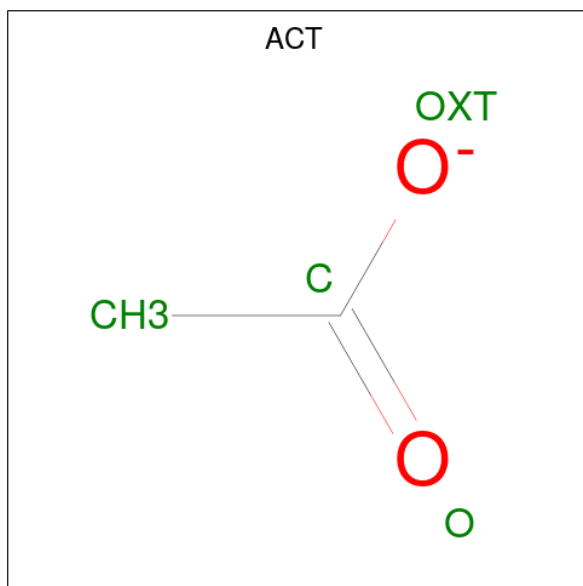
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total C O 4 2 2	0	0
5	B	1	Total C O 4 2 2	0	0
5	B	1	Total C O 4 2 2	0	0
5	B	1	Total C O 4 2 2	0	0

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

- Molecule 6 is ACETATE ION (three-letter code: ACT) (formula:  $C_2H_3O_2$ ).



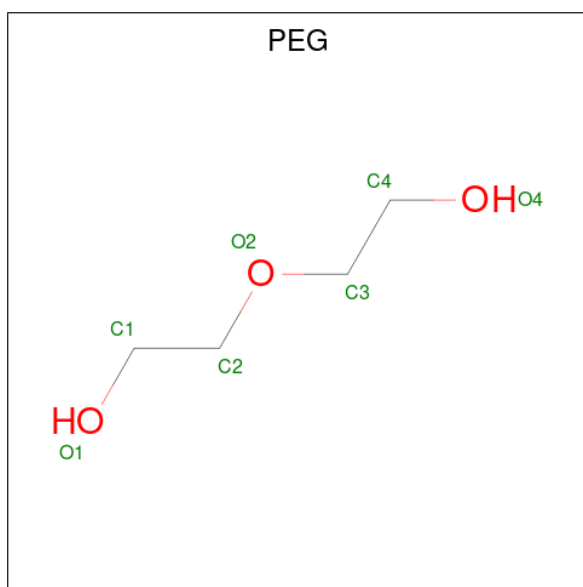
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total	C	O	0	0
			4	2	2		
6	B	1	Total	C	O	0	0
			4	2	2		

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

- Molecule 7 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C<sub>4</sub>H<sub>10</sub>O<sub>3</sub>).

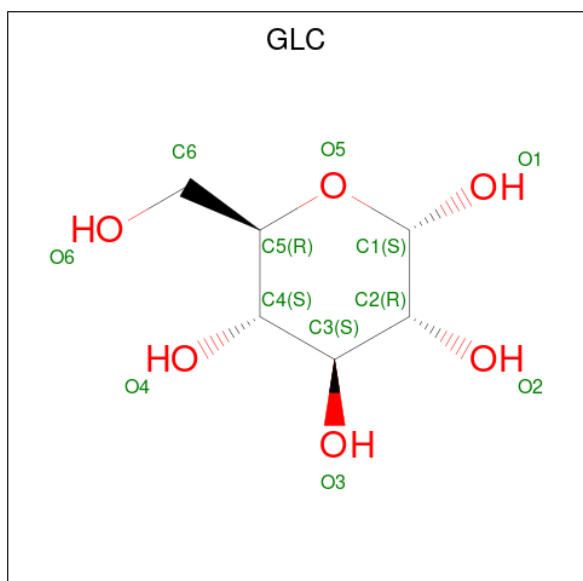


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	C	O	0	0
			7	4	3		

- Molecule 8 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	B	1	Total	Mn	0	0
			1	1		
8	A	1	Total	Mn	0	0
			1	1		

- Molecule 9 is alpha-D-glucopyranose (three-letter code: GLC) (formula: C<sub>6</sub>H<sub>12</sub>O<sub>6</sub>).

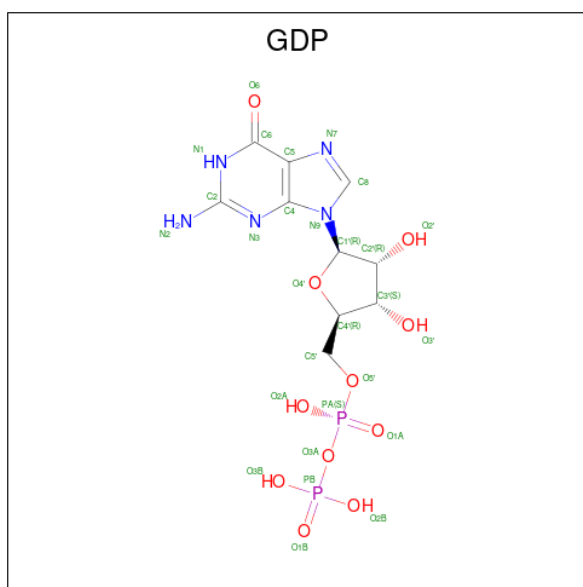


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	B	1	Total	C	O	0	0
			12	6	6		
9	A	1	Total	C	O	0	0
			12	6	6		

- Molecule 10 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	4	Total	Mg	0	0
			4	4		
10	D	1	Total	Mg	0	0
			1	1		

- Molecule 11 is GUANOSINE-5'-DIPHOSPHATE (three-letter code: GDP) (formula: C<sub>10</sub>H<sub>15</sub>N<sub>5</sub>O<sub>11</sub>P<sub>2</sub>).



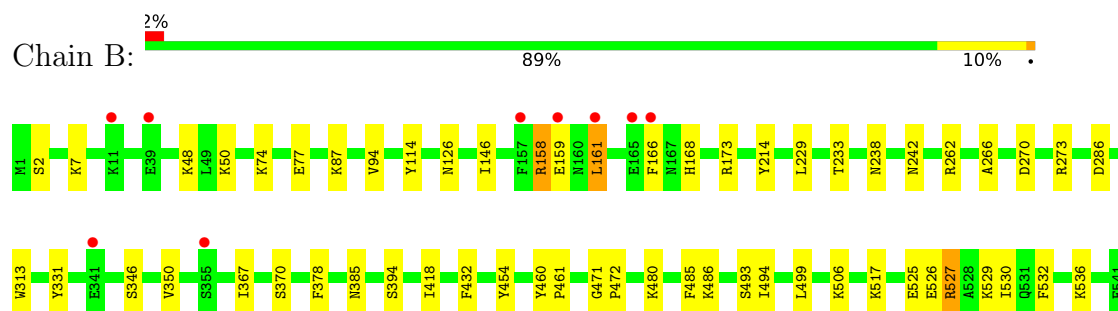
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	A	208	Total 208	O 208	0	0
12	C	63	Total 63	O 63	0	0
12	D	62	Total 62	O 62	0	0

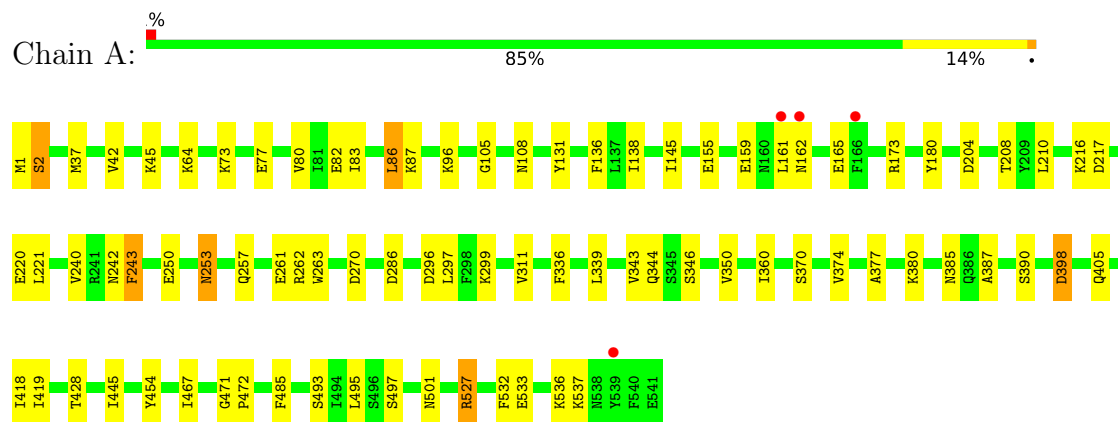
### 3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

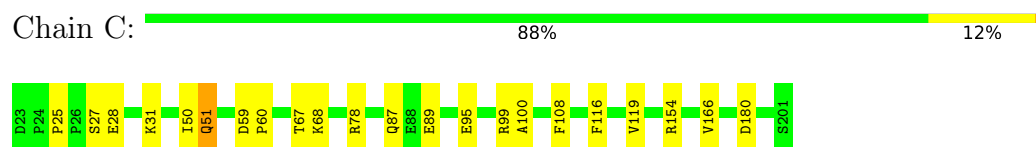
#### • Molecule 1: Toxin B



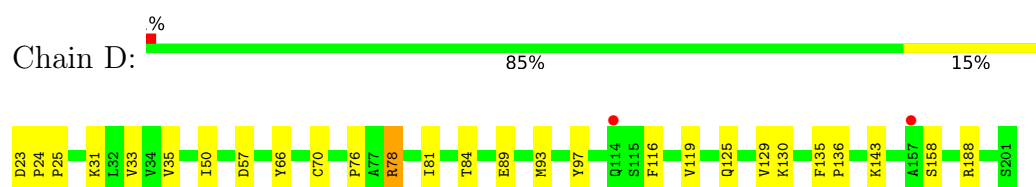
#### • Molecule 1: Toxin B



#### • Molecule 2: Ras-related protein R-Ras



#### • Molecule 2: Ras-related protein R-Ras



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	76.31Å 112.11Å 101.57Å 90.00° 95.89° 90.00°	Depositor
Resolution (Å)	76.02 – 2.34 75.91 – 2.34	Depositor EDS
% Data completeness (in resolution range)	98.8 (76.02-2.34) 98.8 (75.91-2.34)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.78 (at 2.34Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, $R_{free}$	0.201 , 0.241 0.200 , 0.240	Depositor DCC
$R_{free}$ test set	3516 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	36.9	Xtriage
Anisotropy	0.311	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 33.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	12426	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.36% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: EDO, GLC, ACT, PEG, MN, MG, UDP, NH<sub>4</sub>, GDP

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	0.62	0/4485	0.64	0/6062
1	B	0.62	0/4487	0.64	0/6066
2	C	0.61	0/1443	0.67	0/1953
2	D	0.62	0/1431	0.69	0/1937
All	All	0.62	0/11846	0.65	0/16018

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	4402	0	4313	64	0
1	B	4404	0	4315	49	0
2	C	1414	0	1374	22	0
2	D	1402	0	1359	21	0
3	A	25	0	11	0	0
3	B	25	0	11	0	0
4	A	4	0	0	2	0
4	B	4	0	0	1	0
4	C	1	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	20	0	30	5	0
5	B	24	0	36	5	0
5	C	8	0	12	2	0
6	A	36	0	27	4	0
6	B	16	0	12	1	0
6	C	4	0	3	2	0
6	D	4	0	3	0	0
7	B	7	0	10	5	0
8	A	1	0	0	0	0
8	B	1	0	0	0	0
9	A	12	0	11	6	0
9	B	12	0	11	9	0
10	A	4	0	0	0	0
10	D	1	0	0	0	0
11	C	28	0	12	0	0
11	D	28	0	12	1	0
12	A	208	0	0	4	0
12	B	206	0	0	6	0
12	C	63	0	0	2	0
12	D	62	0	0	1	0
All	All	12426	0	11562	147	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 6.

All (147) 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:385:ASN:OD1	9:A:623:GLC:H4	1.59	1.02
1:B:242:ASN:O	7:B:612:PEG:H22	1.65	0.94
2:C:166:VAL:HA	5:C:301:EDO:H11	1.49	0.94
1:B:77:GLU:OE1	2:C:25:PRO:HG2	1.68	0.93
1:B:385:ASN:OD1	9:B:616:GLC:H4	1.70	0.91
1:B:266:ALA:HB1	9:B:616:GLC:H1	1.56	0.85
1:B:471:GLY:HA3	9:B:616:GLC:O3	1.77	0.85
1:B:242:ASN:O	7:B:612:PEG:C2	2.27	0.82
1:A:210:LEU:HD23	1:A:216:LYS:HD3	1.62	0.80
1:A:262:ARG:HD3	4:A:602:NH4:N	1.97	0.80
1:B:161:LEU:H	1:B:161:LEU:HD23	1.48	0.79
2:C:50:ILE:HD12	2:C:68:LYS:HB2	1.65	0.78
1:A:471:GLY:HA3	9:A:623:GLC:O3	1.88	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:472:PRO:HD3	9:A:623:GLC:O3	1.87	0.73
1:A:80:VAL:HG11	6:A:614:ACT:H2	1.69	0.73
1:A:533:GLU:CD	1:A:536:LYS:HZ1	1.91	0.73
1:A:533:GLU:OE1	1:A:536:LYS:NZ	2.21	0.73
1:A:532:PHE:CE2	1:A:536:LYS:HE2	2.26	0.70
1:A:270:ASP:OD2	9:A:623:GLC:H2	1.91	0.70
1:A:253:ASN:HD22	1:A:253:ASN:H	1.39	0.69
2:D:23:ASP:N	2:D:24:PRO:HD2	2.07	0.69
2:D:50:ILE:HD11	2:D:81:ILE:HD12	1.74	0.69
1:A:398:ASP:HB2	12:A:730:HOH:O	1.94	0.68
1:B:161:LEU:HD23	1:B:161:LEU:N	2.10	0.66
7:B:612:PEG:H21	12:B:800:HOH:O	1.96	0.65
1:A:471:GLY:CA	9:A:623:GLC:O3	2.47	0.63
1:A:243:PHE:HA	5:A:604:EDO:H22	1.78	0.62
1:A:418:ILE:HG22	1:A:428:THR:HG23	1.81	0.62
2:D:23:ASP:N	2:D:24:PRO:CD	2.63	0.61
2:D:70:CYS:O	2:D:76:PRO:HA	2.01	0.61
5:A:625:EDO:H11	2:D:188:ARG:HD3	1.82	0.61
1:B:378:PHE:O	1:B:506:LYS:NZ	2.34	0.61
1:B:262:ARG:HD3	12:B:723:HOH:O	2.01	0.60
1:B:480:LYS:HE3	5:B:609:EDO:O1	2.01	0.60
2:C:180:ASP:OD2	6:C:304:ACT:H1	2.02	0.60
1:B:168:HIS:CE1	1:B:529:LYS:HD3	2.39	0.57
1:A:533:GLU:HA	1:A:536:LYS:HZ1	1.70	0.56
1:A:83:ILE:HD11	1:A:485:PHE:CZ	2.40	0.56
1:A:204:ASP:O	1:A:208:THR:HG23	2.05	0.56
1:B:472:PRO:HD3	9:B:616:GLC:O3	2.05	0.55
1:A:374:VAL:HG22	1:A:390:SER:HB2	1.89	0.55
1:A:533:GLU:HA	1:A:536:LYS:NZ	2.22	0.55
1:B:527:ARG:HG3	12:B:818:HOH:O	2.06	0.55
1:B:262:ARG:HD2	1:B:454:TYR:CZ	2.42	0.55
1:B:161:LEU:H	1:B:161:LEU:CD2	2.16	0.54
1:B:493:SER:OG	2:D:78:ARG:NH2	2.40	0.54
1:A:533:GLU:CD	1:A:536:LYS:NZ	2.58	0.54
1:A:216:LYS:HE2	1:A:221:LEU:HD21	1.89	0.53
2:D:23:ASP:N	12:D:2403:HOH:O	2.39	0.53
1:B:418:ILE:HD13	1:B:432:PHE:HA	1.90	0.53
2:D:125:GLN:O	2:D:129:VAL:HG23	2.08	0.53
1:A:253:ASN:HD22	1:A:253:ASN:N	2.06	0.53
1:A:311:VAL:HG22	5:A:607:EDO:H11	1.90	0.53
1:A:495:LEU:HD23	2:C:67:THR:HG22	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:50:LYS:HG3	1:B:485:PHE:HB3	1.91	0.52
1:B:173:ARG:CZ	2:D:89:GLU:HB2	2.40	0.51
2:C:59[A]:ASP:OD2	5:C:302:EDO:H22	2.10	0.51
1:B:346:SER:O	1:B:350:VAL:HG23	2.11	0.51
1:A:262:ARG:HD2	1:A:454:TYR:CE2	2.46	0.50
2:D:84:THR:HG22	2:D:97:TYR:CG	2.45	0.50
1:A:173:ARG:CZ	2:C:89:GLU:HB2	2.42	0.50
1:B:270:ASP:OD2	9:B:616:GLC:H2	2.12	0.49
1:B:532:PHE:CZ	1:B:536:LYS:HD3	2.47	0.49
1:A:87:LYS:O	1:A:370:SER:HB2	2.12	0.49
2:D:116:PHE:O	2:D:119:VAL:HG22	2.13	0.49
2:D:35:VAL:HG12	2:D:84:THR:HG21	1.93	0.49
5:B:611:EDO:H22	12:B:708:HOH:O	2.12	0.49
1:A:336:PHE:CZ	1:A:344:GLN:HG2	2.47	0.49
2:C:154:ARG:NH1	12:C:404:HOH:O	2.46	0.49
1:B:262:ARG:HD2	1:B:454:TYR:CE2	2.48	0.49
1:A:262:ARG:CD	4:A:602:NH4:N	2.71	0.48
1:A:87:LYS:HE2	1:A:501:ASN:HB3	1.94	0.48
1:A:2:SER:HB2	1:A:82:GLU:OE1	2.12	0.48
1:A:380:LYS:HE2	2:C:60:PRO:HD2	1.96	0.48
1:B:471:GLY:CA	9:B:616:GLC:O3	2.55	0.48
2:C:116:PHE:O	2:C:119:VAL:HG22	2.14	0.48
1:A:80:VAL:HG11	6:A:614:ACT:CH3	2.41	0.48
2:D:84:THR:HG22	2:D:97:TYR:CD1	2.49	0.47
1:A:286:ASP:OD2	9:A:623:GLC:H61	2.14	0.47
1:A:2:SER:HB2	12:A:758:HOH:O	2.14	0.47
1:A:445:ILE:HG13	2:C:100:ALA:HB2	1.96	0.47
2:D:143:LYS:HG2	11:D:2302:GDP:C6	2.50	0.47
5:B:611:EDO:C2	12:B:708:HOH:O	2.63	0.47
1:A:262:ARG:HD2	1:A:454:TYR:CZ	2.50	0.47
5:A:604:EDO:H21	12:A:720:HOH:O	2.15	0.46
2:D:130:LYS:HE3	2:D:135:PHE:CZ	2.50	0.46
2:D:50:ILE:HG13	2:D:66:TYR:HB3	1.97	0.46
1:A:296:ASP:O	1:A:299:LYS:HG2	2.16	0.46
1:B:166:PHE:CD1	1:B:166:PHE:N	2.85	0.45
1:B:173:ARG:NH1	2:D:89:GLU:HB2	2.31	0.45
1:A:297:LEU:HD22	1:A:360:ILE:HA	1.97	0.45
1:B:77:GLU:OE1	2:C:25:PRO:CG	2.53	0.45
1:A:339:LEU:HB3	1:A:343:VAL:CG2	2.46	0.45
1:B:126:ASN:ND2	5:B:603:EDO:H22	2.31	0.45
1:B:94:VAL:HG21	1:B:367:ILE:HG22	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:217:ASP:OD2	1:A:220:GLU:HG3	2.16	0.44
2:C:59[B]:ASP:OD1	4:C:303:NH4:N	2.50	0.44
1:B:526:GLU:O	1:B:530:ILE:HG13	2.16	0.44
1:A:42:VAL:HG13	1:A:86:LEU:HD22	2.00	0.44
1:B:238:ASN:OD1	7:B:612:PEG:H42	2.18	0.44
1:B:87:LYS:O	1:B:370:SER:HB2	2.18	0.44
1:A:527:ARG:HD3	1:A:527:ARG:HA	1.64	0.44
2:C:28:GLU:HB2	12:C:456:HOH:O	2.16	0.44
1:A:77:GLU:HG3	2:D:25:PRO:HG2	1.99	0.43
1:A:242:ASN:HD22	5:A:603:EDO:H11	1.83	0.43
1:A:250:GLU:HB3	1:A:405:GLN:OE1	2.19	0.43
1:A:336:PHE:CE2	1:A:344:GLN:HG2	2.53	0.43
1:B:229:LEU:O	1:B:233:THR:HG23	2.19	0.43
1:A:532:PHE:CZ	1:A:536:LYS:HE2	2.53	0.43
1:B:460:TYR:CG	1:B:461:PRO:HD2	2.54	0.43
1:A:467:ILE:HG23	1:A:472:PRO:HD2	2.00	0.43
1:B:214:TYR:OH	12:B:701:HOH:O	2.21	0.43
1:A:131:TYR:O	1:A:240:VAL:HG13	2.19	0.43
1:A:346:SER:O	1:A:350:VAL:HG23	2.18	0.43
1:A:377:ALA:HB2	1:A:387:ALA:HB3	2.01	0.43
1:B:286:ASP:OD2	9:B:616:GLC:H61	2.19	0.42
2:C:50:ILE:HD12	2:C:68:LYS:CB	2.44	0.42
1:A:73:LYS:NZ	12:A:722:HOH:O	2.53	0.42
1:A:257:GLN:O	1:A:261:GLU:HB2	2.18	0.42
1:B:74:LYS:HE2	5:B:617:EDO:O2	2.19	0.42
2:D:135:PHE:CD2	2:D:136:PRO:HD2	2.54	0.42
1:A:419:ILE:HG22	6:A:608:ACT:H2	2.01	0.42
1:B:114:TYR:HH	1:B:331:TYR:HE2	1.64	0.42
1:B:494:ILE:HG23	1:B:499:LEU:HD11	2.01	0.42
1:B:517:LYS:HD3	2:D:57:ASP:O	2.19	0.42
1:A:180:TYR:CE2	6:A:610:ACT:H3	2.54	0.42
1:B:146:ILE:HD13	6:B:608:ACT:H3	2.02	0.42
1:B:394:SER:HB3	4:B:602:NH4:N	2.34	0.42
1:A:493:SER:OG	2:C:78:ARG:NH2	2.53	0.42
2:C:51:GLN:HE21	2:C:51:GLN:HB3	1.50	0.42
2:C:180:ASP:OD2	6:C:304:ACT:CH3	2.68	0.42
1:A:105:GLY:HA2	1:A:136:PHE:O	2.20	0.42
1:A:159:GLU:C	1:A:161:LEU:H	2.22	0.41
1:B:158:ARG:HB3	1:B:159:GLU:HG3	2.01	0.41
2:C:95:GLU:O	2:C:99:ARG:HG3	2.20	0.41
1:B:526:GLU:OE1	1:B:526:GLU:HA	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:138:ILE:HD13	1:A:138:ILE:HA	1.92	0.41
1:B:313:TRP:HA	1:B:313:TRP:CE3	2.56	0.41
2:C:31:LYS:HB3	2:C:100:ALA:O	2.21	0.41
1:A:173:ARG:NH2	2:C:87:GLN:HG3	2.36	0.41
2:C:108:PHE:HB3	2:C:119:VAL:HG11	2.02	0.41
2:D:33:VAL:HG12	2:D:35:VAL:HG13	2.03	0.41
1:B:266:ALA:HB1	9:B:616:GLC:C1	2.39	0.41
1:B:242:ASN:O	7:B:612:PEG:H21	2.15	0.40
1:B:273:ARG:HH12	9:B:616:GLC:H62	1.86	0.40
1:A:145:ILE:HD12	1:A:263:TRP:CH2	2.57	0.40
1:A:37:MET:O	1:A:45:LYS:HE2	2.22	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	539/541 (100%)	524 (97%)	15 (3%)	0	100	100
1	B	539/541 (100%)	526 (98%)	13 (2%)	0	100	100
2	C	178/179 (99%)	170 (96%)	8 (4%)	0	100	100
2	D	177/179 (99%)	168 (95%)	9 (5%)	0	100	100
All	All	1433/1440 (100%)	1388 (97%)	45 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 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	490/494 (99%)	475 (97%)	15 (3%)	40	49
1	B	491/494 (99%)	483 (98%)	8 (2%)	62	74
2	C	152/151 (101%)	150 (99%)	2 (1%)	69	79
2	D	150/151 (99%)	146 (97%)	4 (3%)	44	55
All	All	1283/1290 (100%)	1254 (98%)	29 (2%)	50	61

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	2	SER
1	B	7	LYS
1	B	48	LYS
1	B	158	ARG
1	B	161	LEU
1	B	486	LYS
1	B	525	GLU
1	B	527	ARG
1	A	1	MET
1	A	2	SER
1	A	64	LYS
1	A	86	LEU
1	A	96	LYS
1	A	108	ASN
1	A	155	GLU
1	A	162	ASN
1	A	165	GLU
1	A	243	PHE
1	A	253	ASN
1	A	398	ASP
1	A	497	SER
1	A	527	ARG
1	A	537	LYS
2	C	27	SER
2	C	51	GLN
2	D	31	LYS
2	D	78	ARG
2	D	93	MET
2	D	158	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	106	GLN
1	B	126	ASN
1	B	508	ASN
1	B	522	GLN
1	B	538	ASN
1	A	242	ASN
1	A	253	ASN
2	C	51	GLN
2	D	51	GLN
2	D	196	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

## 5.6 Ligand geometry ⓘ

Of 51 ligands modelled in this entry, 9 are modelled with single atom and 7 are monoatomic - leaving 35 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	$\# Z  > 2$
5	EDO	B	609	-	3,3,3	0.13	0	2,2,2	0.21	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
5	EDO	B	611	-	3,3,3	0.04	0	2,2,2	0.31	0
7	PEG	B	612	-	6,6,6	0.39	0	5,5,5	0.39	0
6	ACT	A	614	-	1,3,3	3.37	1 (100%)	0,3,3	0.00	-
5	EDO	A	603	-	3,3,3	0.11	0	2,2,2	0.38	0
5	EDO	C	302	-	3,3,3	0.20	0	2,2,2	0.45	0
6	ACT	B	610	-	1,3,3	3.90	1 (100%)	0,3,3	0.00	-
6	ACT	B	608	-	1,3,3	3.17	1 (100%)	0,3,3	0.00	-
5	EDO	A	607	-	3,3,3	0.08	0	2,2,2	0.25	0
5	EDO	B	603	-	3,3,3	0.11	0	2,2,2	0.38	0
6	ACT	B	607	-	1,3,3	3.70	1 (100%)	0,3,3	0.00	-
3	UDP	B	601	8	20,26,26	0.84	1 (5%)	25,40,40	1.08	1 (4%)
6	ACT	A	616	-	1,3,3	3.49	1 (100%)	0,3,3	0.00	-
6	ACT	A	610	-	1,3,3	3.44	1 (100%)	0,3,3	0.00	-
6	ACT	B	618	-	1,3,3	3.61	1 (100%)	0,3,3	0.00	-
6	ACT	A	608	-	1,3,3	3.47	1 (100%)	0,3,3	0.00	-
5	EDO	A	605	-	3,3,3	0.03	0	2,2,2	0.17	0
5	EDO	B	604	-	3,3,3	0.12	0	2,2,2	0.15	0
5	EDO	A	625	-	3,3,3	0.34	0	2,2,2	0.75	0
6	ACT	A	611	-	1,3,3	3.52	1 (100%)	0,3,3	0.00	-
9	GLC	A	623	-	12,12,12	1.33	2 (16%)	17,17,17	2.51	7 (41%)
9	GLC	B	616	-	12,12,12	1.33	2 (16%)	17,17,17	2.64	6 (35%)
6	ACT	A	622	-	1,3,3	3.38	1 (100%)	0,3,3	0.00	-
3	UDP	A	601	8,10	20,26,26	0.86	0	25,40,40	1.07	2 (8%)
11	GDP	D	2302	10	24,30,30	1.14	3 (12%)	31,47,47	1.79	6 (19%)
6	ACT	A	613	-	1,3,3	3.92	1 (100%)	0,3,3	0.00	-
6	ACT	A	606	-	1,3,3	3.75	1 (100%)	0,3,3	0.00	-
5	EDO	B	617	-	3,3,3	0.13	0	2,2,2	0.22	0
5	EDO	C	301	-	3,3,3	0.13	0	2,2,2	0.47	0
6	ACT	C	304	-	1,3,3	3.55	1 (100%)	0,3,3	0.00	-
6	ACT	D	2301	-	1,3,3	3.69	1 (100%)	0,3,3	0.00	-
11	GDP	C	305	-	24,30,30	1.10	2 (8%)	31,47,47	1.91	7 (22%)
5	EDO	B	606	-	3,3,3	0.07	0	2,2,2	0.33	0
5	EDO	A	604	-	3,3,3	0.07	0	2,2,2	0.21	0
6	ACT	A	615	-	1,3,3	3.30	1 (100%)	0,3,3	0.00	-

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
5	EDO	B	609	-	-	1/1/1/1	-
5	EDO	B	611	-	-	1/1/1/1	-
7	PEG	B	612	-	-	2/4/4/4	-
5	EDO	A	603	-	-	1/1/1/1	-
5	EDO	C	302	-	-	1/1/1/1	-
5	EDO	A	607	-	-	1/1/1/1	-
5	EDO	B	603	-	-	0/1/1/1	-
3	UDP	B	601	8	-	0/14/32/32	0/2/2/2
5	EDO	A	605	-	-	1/1/1/1	-
5	EDO	B	604	-	-	0/1/1/1	-
5	EDO	A	625	-	-	0/1/1/1	-
9	GLC	A	623	-	-	1/2/22/22	0/1/1/1
9	GLC	B	616	-	-	2/2/22/22	0/1/1/1
3	UDP	A	601	8,10	-	0/14/32/32	0/2/2/2
11	GDP	D	2302	10	-	0/12/32/32	0/3/3/3
5	EDO	B	617	-	-	0/1/1/1	-
5	EDO	C	301	-	-	0/1/1/1	-
11	GDP	C	305	-	-	1/12/32/32	0/3/3/3
5	EDO	B	606	-	-	0/1/1/1	-
5	EDO	A	604	-	-	1/1/1/1	-

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	613	ACT	CH3-C	3.92	1.53	1.48
6	B	610	ACT	CH3-C	3.90	1.53	1.48
11	C	305	GDP	C6-C5	3.80	1.47	1.41
6	A	606	ACT	CH3-C	3.75	1.53	1.48
6	B	607	ACT	CH3-C	3.70	1.53	1.48
6	D	2301	ACT	CH3-C	3.69	1.53	1.48
6	B	618	ACT	CH3-C	3.61	1.53	1.48
6	C	304	ACT	CH3-C	3.55	1.53	1.48
6	A	611	ACT	CH3-C	3.52	1.53	1.48
11	D	2302	GDP	C6-C5	3.51	1.47	1.41
6	A	616	ACT	CH3-C	3.49	1.53	1.48
6	A	608	ACT	CH3-C	3.47	1.53	1.48
6	A	610	ACT	CH3-C	3.44	1.53	1.48
6	A	622	ACT	CH3-C	3.38	1.53	1.48
6	A	614	ACT	CH3-C	3.37	1.53	1.48
9	A	623	GLC	O5-C1	3.35	1.51	1.42
6	A	615	ACT	CH3-C	3.30	1.52	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	608	ACT	CH3-C	3.17	1.52	1.48
9	B	616	GLC	O5-C1	3.03	1.50	1.42
11	D	2302	GDP	C5-C4	2.34	1.47	1.40
11	C	305	GDP	C5-C4	2.33	1.47	1.40
11	D	2302	GDP	O4'-C1'	2.15	1.44	1.41
9	B	616	GLC	C3-C2	-2.06	1.47	1.52
9	A	623	GLC	C3-C2	-2.03	1.47	1.52
3	B	601	UDP	C2-N3	-2.00	1.34	1.38

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	A	623	GLC	C4-C3-C2	6.33	121.87	110.82
9	B	616	GLC	O3-C3-C2	-5.39	97.88	110.35
9	B	616	GLC	C4-C3-C2	5.28	120.03	110.82
9	B	616	GLC	C1-O5-C5	-4.77	104.66	113.66
11	C	305	GDP	C6-N1-C2	4.55	123.16	115.93
9	A	623	GLC	O3-C3-C2	-4.54	99.85	110.35
11	D	2302	GDP	C5-C6-N1	-4.34	117.50	123.43
11	D	2302	GDP	C6-N1-C2	4.34	122.82	115.93
11	C	305	GDP	C6-C5-C4	-4.29	116.70	120.80
11	C	305	GDP	C5-C6-N1	-4.20	117.69	123.43
11	D	2302	GDP	C2-N3-C4	4.03	119.96	115.36
9	A	623	GLC	C1-O5-C5	-3.99	106.14	113.66
11	C	305	GDP	C2-N3-C4	3.93	119.84	115.36
11	C	305	GDP	N3-C2-N1	-3.43	122.65	127.22
11	D	2302	GDP	C6-C5-C4	-3.40	117.55	120.80
11	D	2302	GDP	N3-C2-N1	-3.28	122.85	127.22
9	B	616	GLC	O5-C5-C6	3.17	114.31	106.44
3	A	601	UDP	PA-O3A-PB	-3.14	122.06	132.83
3	B	601	UDP	PA-O3A-PB	-2.96	122.69	132.83
9	B	616	GLC	C3-C4-C5	2.76	115.17	110.24
9	A	623	GLC	O2-C2-C3	-2.76	103.97	110.35
9	A	623	GLC	C1-C2-C3	2.68	115.88	110.31
9	A	623	GLC	O5-C5-C6	2.63	112.97	106.44
11	C	305	GDP	PA-O3A-PB	-2.55	124.07	132.83
11	D	2302	GDP	PA-O3A-PB	-2.44	124.47	132.83
9	B	616	GLC	O2-C2-C3	-2.37	104.88	110.35
3	A	601	UDP	O3B-PB-O2B	2.23	116.15	107.64
11	C	305	GDP	C4-C5-N7	-2.22	107.09	109.40
9	A	623	GLC	C3-C4-C5	2.01	113.82	110.24

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
9	B	616	GLC	O5-C5-C6-O6
9	A	623	GLC	O5-C5-C6-O6
5	C	302	EDO	O1-C1-C2-O2
7	B	612	PEG	O1-C1-C2-O2
5	B	611	EDO	O1-C1-C2-O2
11	C	305	GDP	PA-O3A-PB-O1B
5	A	605	EDO	O1-C1-C2-O2
5	B	609	EDO	O1-C1-C2-O2
7	B	612	PEG	O2-C3-C4-O4
5	A	603	EDO	O1-C1-C2-O2
5	A	604	EDO	O1-C1-C2-O2
5	A	607	EDO	O1-C1-C2-O2
9	B	616	GLC	C4-C5-C6-O6

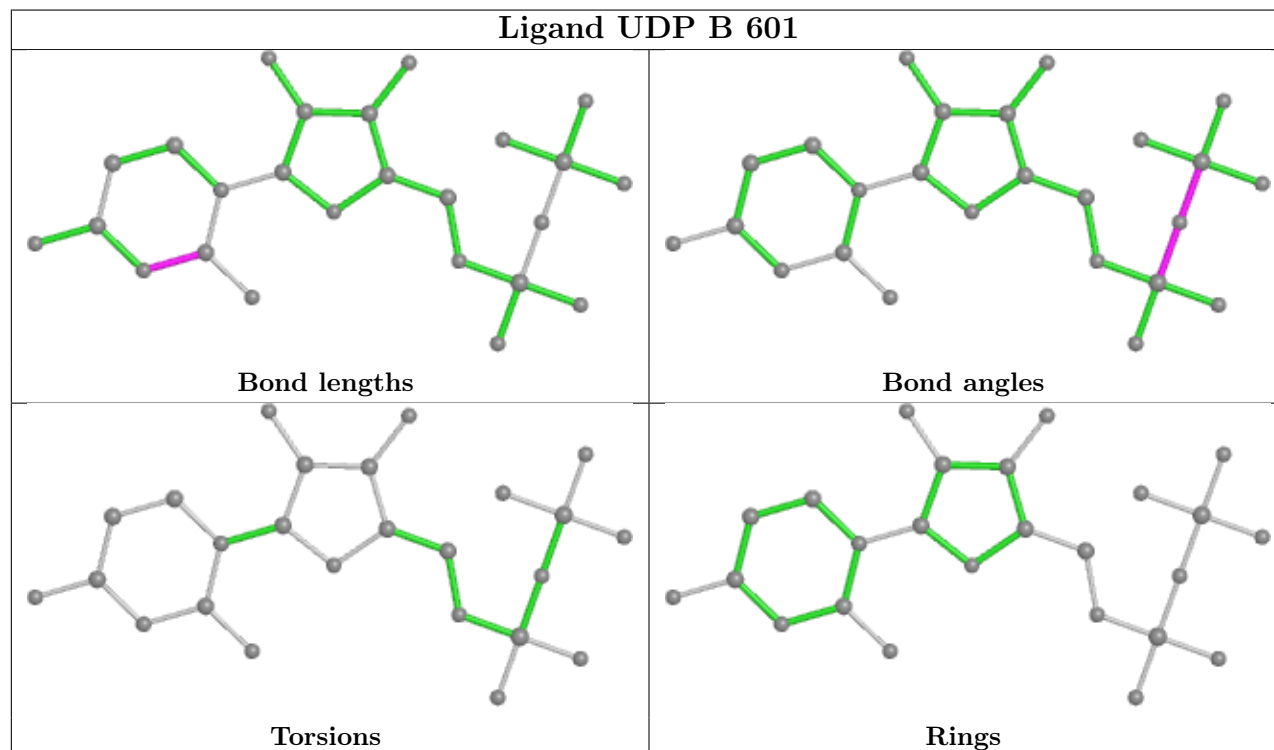
There are no ring outliers.

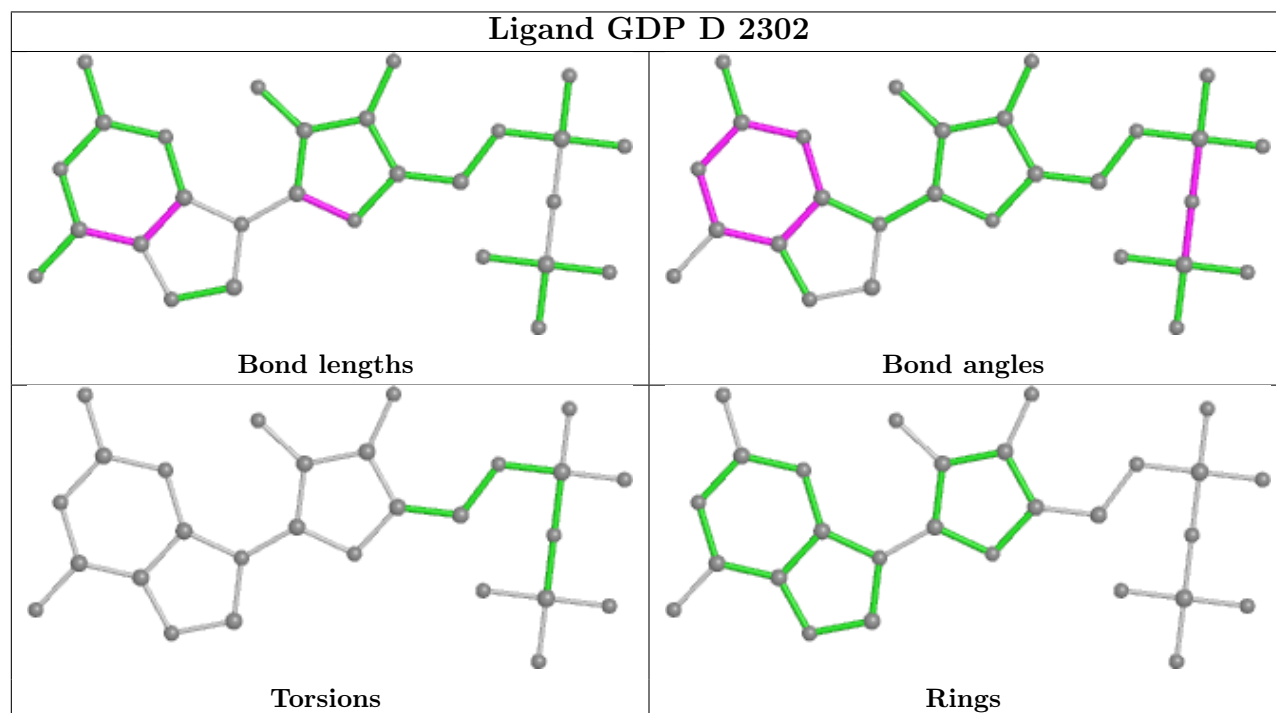
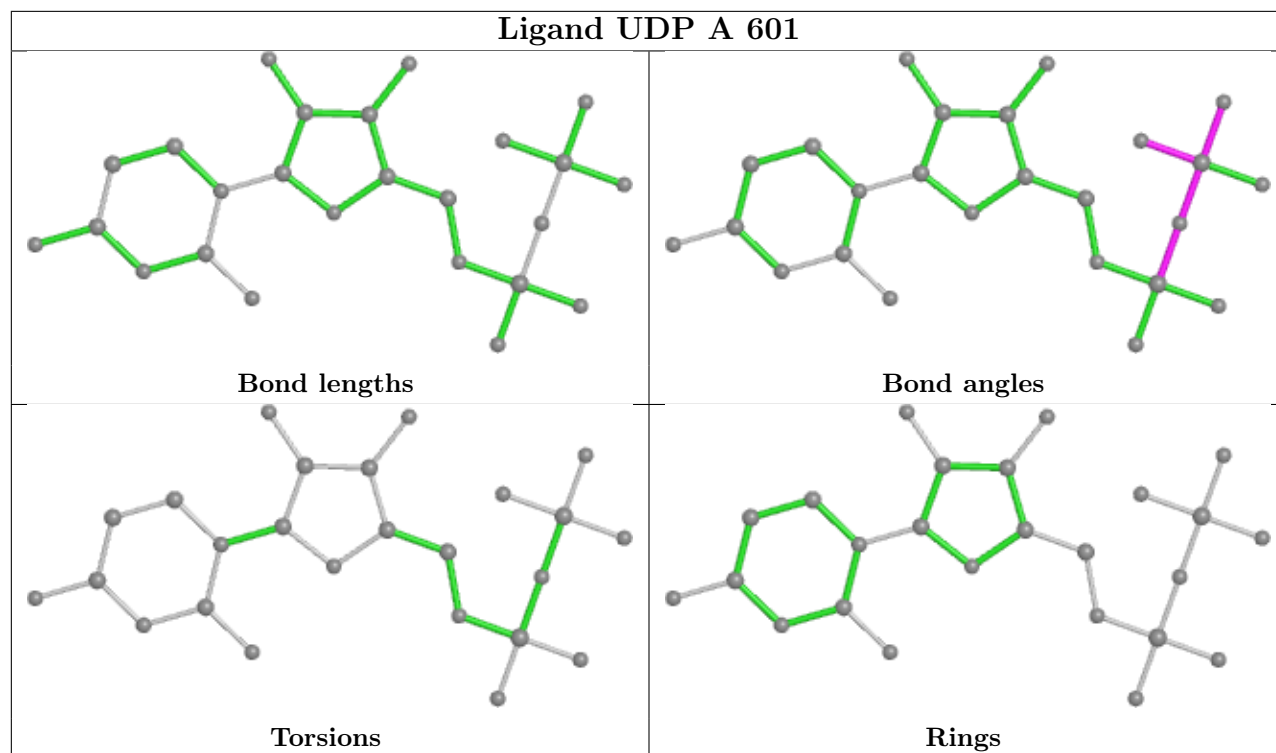
19 monomers are involved in 40 short contacts:

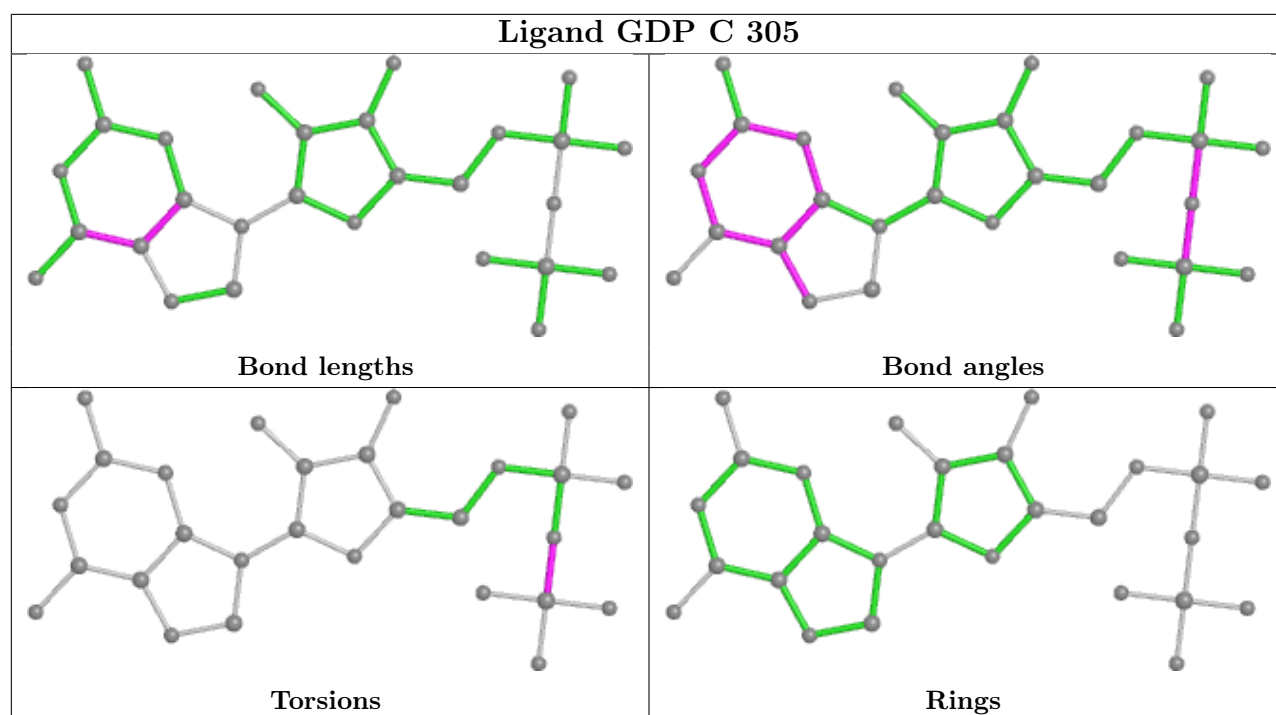
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	609	EDO	1	0
5	B	611	EDO	2	0
7	B	612	PEG	5	0
6	A	614	ACT	2	0
5	A	603	EDO	1	0
5	C	302	EDO	1	0
6	B	608	ACT	1	0
5	A	607	EDO	1	0
5	B	603	EDO	1	0
6	A	610	ACT	1	0
6	A	608	ACT	1	0
5	A	625	EDO	1	0
9	A	623	GLC	6	0
9	B	616	GLC	9	0
11	D	2302	GDP	1	0
5	B	617	EDO	1	0
5	C	301	EDO	1	0
6	C	304	ACT	2	0
5	A	604	EDO	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 than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.







## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	541/541 (100%)	-0.05	4 (0%) 87 92	26, 38, 60, 81	0
1	B	541/541 (100%)	-0.11	9 (1%) 70 78	28, 39, 58, 90	0
2	C	179/179 (100%)	-0.03	0 100 100	28, 38, 54, 77	0
2	D	179/179 (100%)	0.01	2 (1%) 80 86	29, 41, 57, 72	0
All	All	1440/1440 (100%)	-0.06	15 (1%) 82 88	26, 39, 58, 90	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	161	LEU	3.6
1	A	161	LEU	3.3
1	B	166	PHE	2.8
1	A	166	PHE	2.7
1	B	39	GLU	2.7
1	B	159	GLU	2.7
1	B	341	GLU	2.7
1	B	165	GLU	2.3
2	D	157	ALA	2.3
1	A	162	ASN	2.3
1	B	157	PHE	2.3
1	A	539	TYR	2.2
1	B	11	LYS	2.2
1	B	355	SER	2.2
2	D	114	GLN	2.1

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

## 6.4 Ligands ⓘ

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
5	EDO	A	625	4/4	0.32	0.35	50,51,52,52	0
6	ACT	B	618	4/4	0.64	0.31	49,54,54,57	0
10	MG	A	624	1/1	0.68	0.16	65,65,65,65	0
5	EDO	C	301	4/4	0.70	0.36	49,50,51,52	0
6	ACT	A	611	4/4	0.71	0.23	61,62,63,63	0
7	PEG	B	612	7/7	0.73	0.23	54,56,57,58	0
6	ACT	C	304	4/4	0.73	0.22	60,62,64,65	0
10	MG	D	2303	1/1	0.74	0.08	60,60,60,60	0
5	EDO	C	302	4/4	0.75	0.22	49,51,51,51	0
5	EDO	B	603	4/4	0.76	0.26	54,54,55,57	0
6	ACT	A	615	4/4	0.76	0.34	50,54,57,57	0
5	EDO	A	604	4/4	0.79	0.16	49,49,50,52	0
9	GLC	B	616	12/12	0.81	0.31	25,26,26,27	12
5	EDO	B	609	4/4	0.82	0.30	54,54,55,57	0
9	GLC	A	623	12/12	0.82	0.32	25,26,26,27	12
6	ACT	B	610	4/4	0.84	0.36	60,62,63,65	0
6	ACT	A	622	4/4	0.84	0.31	31,31,31,32	0
5	EDO	B	606	4/4	0.84	0.32	45,51,53,55	0
4	NH4	A	602	1/1	0.84	0.30	28,28,28,28	0
5	EDO	A	603	4/4	0.85	0.25	46,50,50,51	0
4	NH4	C	303	1/1	0.85	0.33	38,38,38,38	0
6	ACT	A	613	4/4	0.85	0.16	68,68,68,68	0
6	ACT	A	614	4/4	0.85	0.30	57,58,59,59	0
4	NH4	B	614	1/1	0.88	0.21	33,33,33,33	0
5	EDO	B	617	4/4	0.89	0.24	53,53,54,55	0
6	ACT	A	608	4/4	0.89	0.17	57,58,59,59	0
6	ACT	D	2301	4/4	0.89	0.16	63,64,65,66	0
6	ACT	A	616	4/4	0.89	0.26	66,67,67,68	0
5	EDO	B	611	4/4	0.90	0.33	51,53,53,54	0
6	ACT	A	606	4/4	0.90	0.22	58,60,60,62	0
6	ACT	B	607	4/4	0.91	0.31	49,50,51,53	0
4	NH4	B	615	1/1	0.91	0.33	30,30,30,30	0

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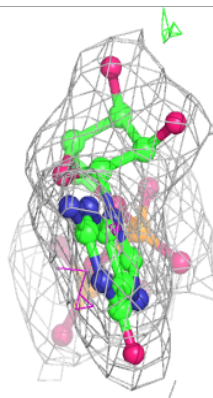
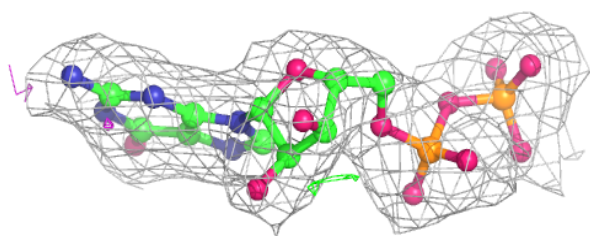
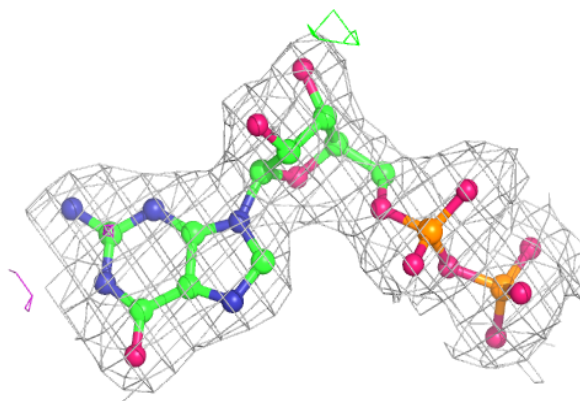
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
4	NH4	A	612	1/1	0.92	0.46	42,42,42,42	0
10	MG	A	621	1/1	0.93	0.20	50,50,50,50	0
6	ACT	A	610	4/4	0.93	0.24	61,61,62,64	0
4	NH4	B	602	1/1	0.93	0.30	28,28,28,28	0
5	EDO	A	607	4/4	0.94	0.13	48,48,49,50	0
4	NH4	A	617	1/1	0.94	0.24	42,42,42,42	0
4	NH4	B	605	1/1	0.95	0.21	20,20,20,20	0
5	EDO	B	604	4/4	0.95	0.18	43,43,44,44	0
5	EDO	A	605	4/4	0.95	0.15	50,51,52,55	0
6	ACT	B	608	4/4	0.96	0.16	51,52,53,55	0
11	GDP	D	2302	28/28	0.97	0.12	35,36,38,38	0
4	NH4	A	609	1/1	0.98	0.36	19,19,19,19	0
3	UDP	A	601	25/25	0.98	0.12	27,29,34,36	0
11	GDP	C	305	28/28	0.98	0.12	34,34,35,36	0
3	UDP	B	601	25/25	0.98	0.11	28,31,33,35	0
8	MN	A	618	1/1	0.99	0.14	29,29,29,29	0
10	MG	A	619	1/1	0.99	0.12	39,39,39,39	0
10	MG	A	620	1/1	0.99	0.19	39,39,39,39	0
8	MN	B	613	1/1	0.99	0.12	33,33,33,33	0

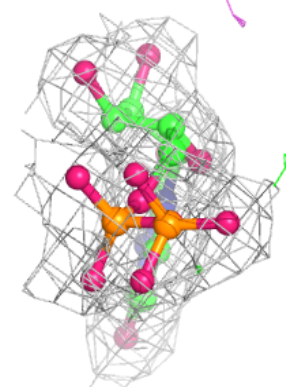
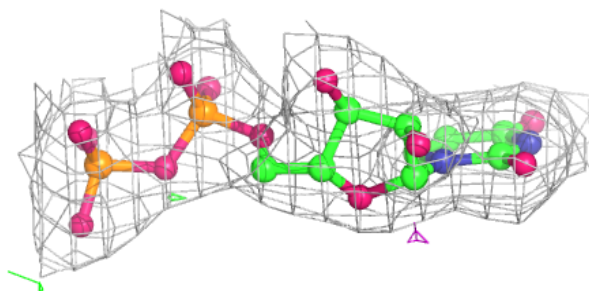
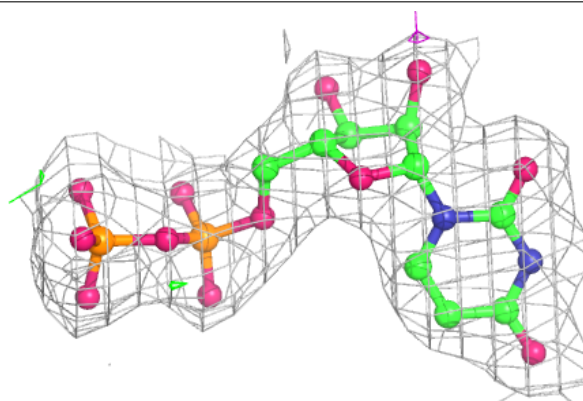
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around GDP D 2302:**

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

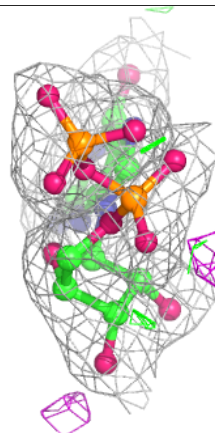
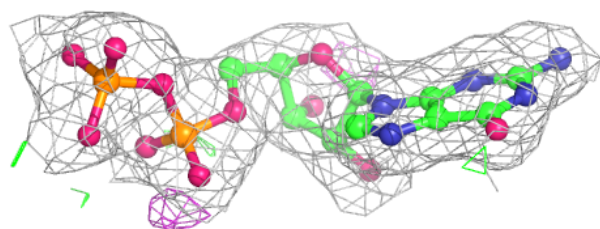
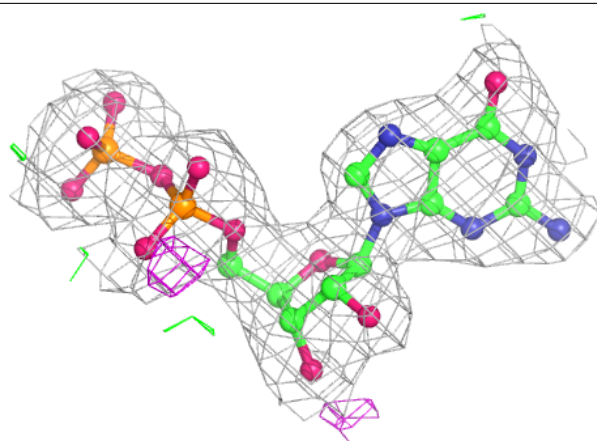
**Electron density around UDP A 601:**

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

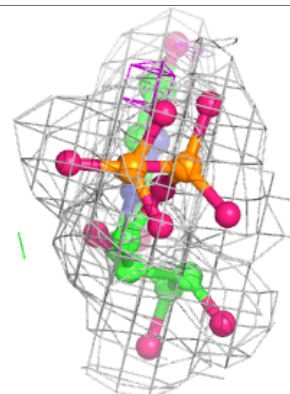
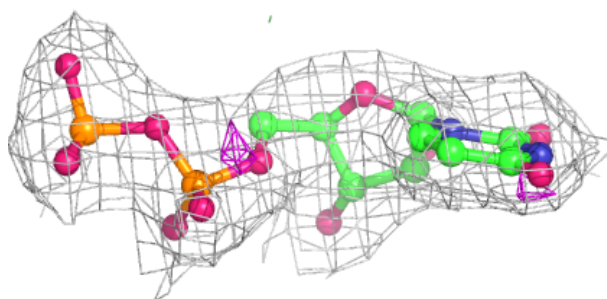
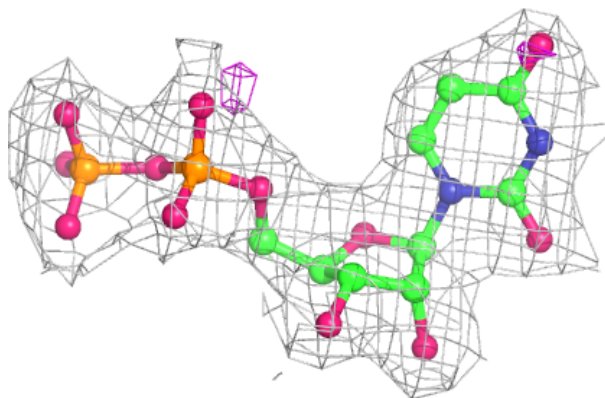


**Electron density around GDP C 305:**

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

**Electron density around UDP B 601:**

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



## 6.5 Other polymers ⓘ

There are no such residues in this entry.