



Full wwPDB X-ray Structure Validation Report

May 13, 2020 – 08:06 pm BST

PDB ID : 4F97
Title : Crystal Structure of VldE, the pseudo-glycosyltransferase, in complex with GDP and validoxylamine A 7'-phosphate
Authors : Cavalier, M.C.; Yim, Y.-S.; Asamizu, S.; Neau, D.; Almabruk, K.H.; Mahmud, T.; Lee, Y.-H.
Deposited on : 2012-05-18
Resolution : 2.11 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

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

with specific help available everywhere you see the  symbol.

The 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 : 2.11
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) : 2.11

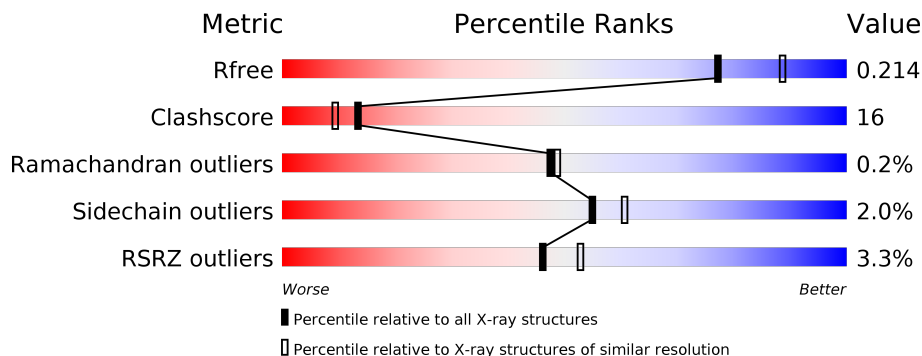
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.11 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	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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

Mol	Chain	Length	Quality of chain
1	A	497	 2% 74% 20% • 5%
1	B	497	 4% 71% 22% • 5%

2 Entry composition [i](#)

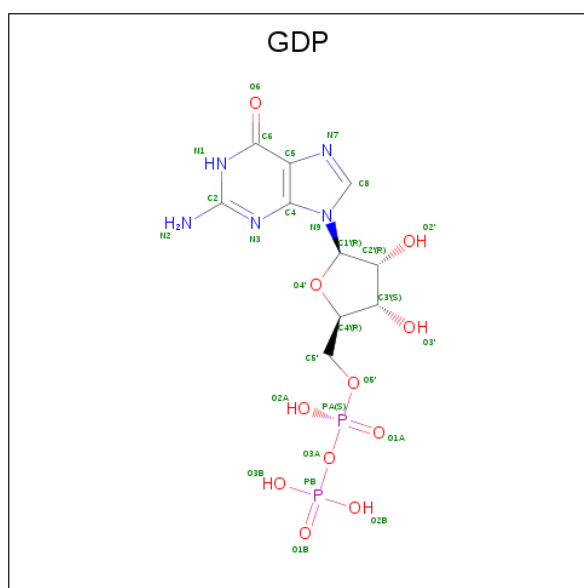
There are 6 unique types of molecules in this entry. The entry contains 7969 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 VldE.

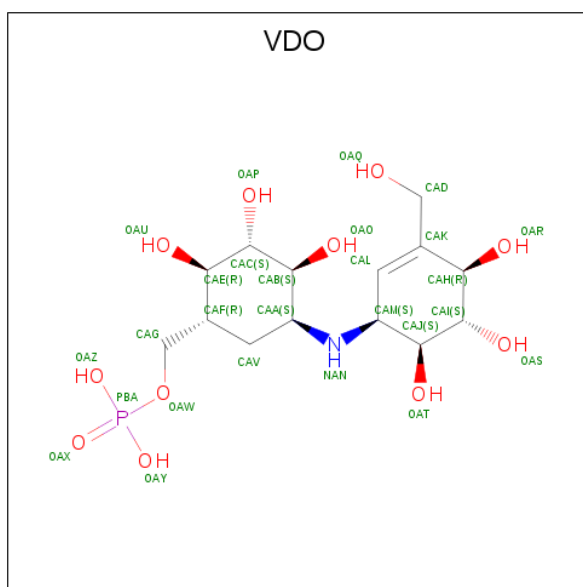
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	472	Total	C	N	O	S	0	0	0
			3688	2307	680	689	12			
1	B	470	Total	C	N	O	S	0	0	0
			3673	2298	676	687	12			

- Molecule 2 is GUANOSINE-5'-DIPHOSPHATE (three-letter code: GDP) (formula: C₁₀H₁₅N₅O₁₁P₂).



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

- Molecule 3 is [(1R,2R,3S,4S,5S)-2,3,4-TRIHYDROXY-5-([(1S,4R,5S,6S)-4,5,6-TRIHYDROXY-3-(HYDROXYMETHYL)CYCLOHEX-2-EN-1-YL]AMINO)CYCLOHEXYL]METHYL DIHYDROGEN PHOSPHATE (three-letter code: VDO) (formula: C₁₄H₂₆N₂O₁₁P).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
			Total	C	N	O			P
3	A	1	Total	C	N	O	P	0	0
			27	14	1	11	1		
3	B	1	Total	C	N	O	P	0	0
			27	14	1	11	1		

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



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

- Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total Cl 1 1	0	0
5	A	1	Total Cl 1 1	0	0

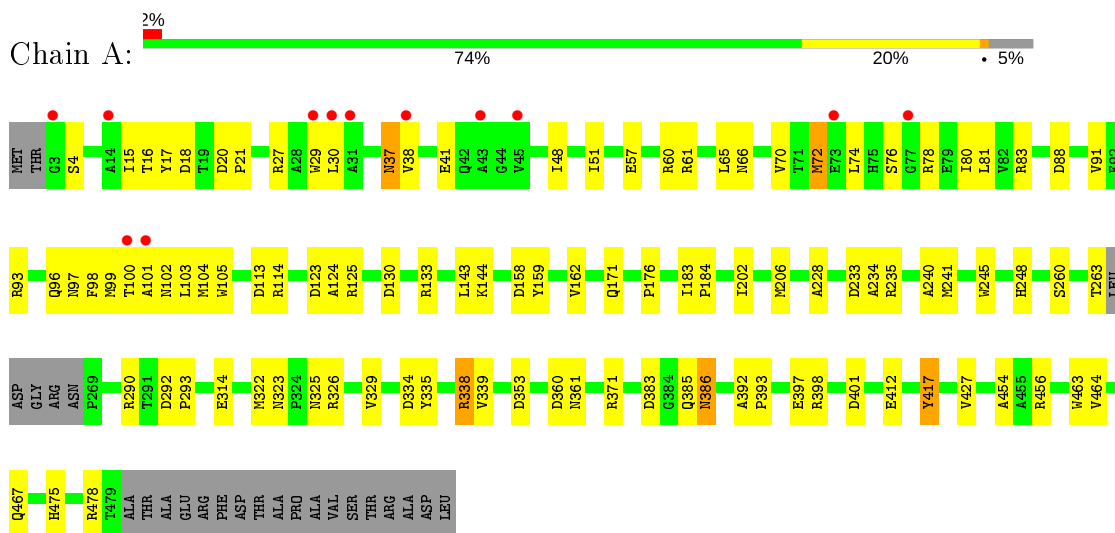
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	267	Total O 267 267	0	0
6	B	221	Total O 221 221	0	0

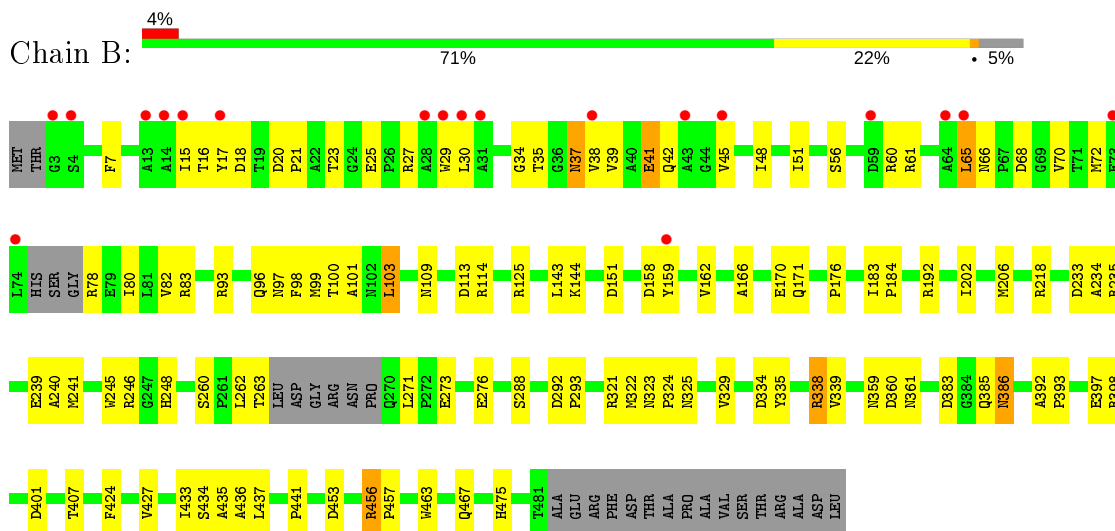
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: VldE



- Molecule 1: VldE



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	47.96Å 120.71Å 84.24Å 90.00° 91.63° 90.00°	Depositor
Resolution (Å)	47.94 - 2.11 47.94 - 2.11	Depositor EDS
% Data completeness (in resolution range)	99.7 (47.94-2.11) 95.0 (47.94-2.11)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.64 (at 2.10Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.191 , 0.222 0.183 , 0.214	Depositor DCC
R_{free} test set	2791 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	30.5	Xtrriage
Anisotropy	0.118	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 32.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	0.217 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7969	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.60% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² 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: GDP, EDO, VDO, CL

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	0.71	5/3774 (0.1%)	0.77	8/5143 (0.2%)
1	B	0.70	6/3756 (0.2%)	0.78	9/5118 (0.2%)
All	All	0.71	11/7530 (0.1%)	0.77	17/10261 (0.2%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	417	TYR	CE2-CZ	-10.73	1.24	1.38
1	B	456	ARG	CZ-NH1	-9.24	1.21	1.33
1	A	456	ARG	CZ-NH1	-9.03	1.21	1.33
1	B	7	PHE	CG-CD1	-8.83	1.25	1.38
1	B	7	PHE	CG-CD2	-8.81	1.25	1.38
1	A	417	TYR	CE1-CZ	-8.77	1.27	1.38
1	A	417	TYR	CG-CD1	-8.50	1.28	1.39
1	A	417	TYR	CG-CD2	-8.30	1.28	1.39
1	B	7	PHE	CE1-CZ	-7.31	1.23	1.37
1	B	7	PHE	CE2-CZ	-7.04	1.24	1.37
1	B	109	ASN	CG-OD1	-5.15	1.12	1.24

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	456	ARG	NE-CZ-NH2	18.82	129.71	120.30
1	A	456	ARG	NE-CZ-NH2	18.68	129.64	120.30
1	A	93	ARG	CB-CG-CD	-7.68	91.63	111.60
1	A	338	ARG	NE-CZ-NH2	-7.64	116.48	120.30
1	B	338	ARG	NE-CZ-NH2	-7.63	116.48	120.30
1	A	93	ARG	NE-CZ-NH1	-6.82	116.89	120.30
1	B	456	ARG	NE-CZ-NH1	-6.68	116.96	120.30
1	B	65	LEU	CB-CG-CD2	-6.23	100.41	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	93	ARG	CB-CG-CD	-6.02	95.95	111.60
1	A	456	ARG	NH1-CZ-NH2	-5.97	112.83	119.40
1	A	93	ARG	CD-NE-CZ	-5.85	115.42	123.60
1	A	456	ARG	NE-CZ-NH1	-5.55	117.52	120.30
1	B	456	ARG	NH1-CZ-NH2	-5.52	113.32	119.40
1	B	41	GLU	O-C-N	-5.46	113.97	122.70
1	A	98	PHE	C-N-CA	5.33	135.04	121.70
1	B	93	ARG	NE-CZ-NH1	-5.24	117.68	120.30
1	B	338	ARG	NE-CZ-NH1	5.01	122.81	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3688	0	3577	116	0
1	B	3673	0	3565	119	0
2	A	28	0	12	0	0
2	B	28	0	12	0	0
3	A	27	0	24	3	0
3	B	27	0	24	3	0
4	A	4	0	6	1	0
4	B	4	0	6	2	0
5	A	1	0	0	1	0
5	B	1	0	0	1	0
6	A	267	0	0	19	0
6	B	221	0	0	20	0
All	All	7969	0	7226	236	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 16.

All (236) 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:B:48:ILE:HG23	1:B:80:ILE:CD1	1.51	1.38
1:B:48:ILE:CG2	1:B:80:ILE:HD13	1.66	1.26
1:A:83:ARG:HH22	1:A:144:LYS:HD3	1.03	1.12
1:A:61:ARG:O	1:A:65:LEU:HD13	1.49	1.10
1:B:83:ARG:HH22	1:B:144:LYS:HD3	0.96	1.06
1:A:176:PRO:HG3	1:A:475:HIS:ND1	1.72	1.04
1:A:88:ASP:CG	6:A:843:HOH:O	1.97	1.03
1:A:96:GLN:O	1:A:100:THR:HG23	1.58	1.03
1:B:48:ILE:HG23	1:B:80:ILE:HD13	1.04	1.03
1:B:83:ARG:NH2	1:B:144:LYS:HD3	1.74	1.01
1:B:48:ILE:HG23	1:B:80:ILE:HD12	1.44	1.00
1:A:478:ARG:HD3	6:A:852:HOH:O	1.60	1.00
1:A:314:GLU:OE2	6:A:857:HOH:O	1.80	0.98
1:B:56:SER:O	1:B:60:ARG:HG2	1.62	0.97
1:A:125:ARG:O	6:A:809:HOH:O	1.83	0.95
1:A:101:ALA:O	6:A:863:HOH:O	1.83	0.95
1:B:176:PRO:HG3	1:B:475:HIS:ND1	1.81	0.95
1:A:83:ARG:NH2	1:A:144:LYS:HD3	1.81	0.94
1:B:48:ILE:HG12	1:B:80:ILE:HD11	1.47	0.94
1:B:48:ILE:O	1:B:80:ILE:HD12	1.68	0.93
1:B:96:GLN:O	1:B:100:THR:HG23	1.69	0.92
1:B:48:ILE:CG2	1:B:80:ILE:CD1	2.36	0.92
1:A:37:ASN:H	1:A:37:ASN:ND2	1.60	0.91
1:A:70:VAL:CG1	1:A:72:MET:CE	2.52	0.88
1:A:96:GLN:O	1:A:100:THR:CG2	2.22	0.88
1:A:37:ASN:N	1:A:37:ASN:HD22	1.72	0.87
1:A:99:MET:HE1	1:A:104:MET:HG3	1.57	0.87
1:B:48:ILE:HG12	1:B:80:ILE:CD1	2.04	0.87
1:A:37:ASN:H	1:A:37:ASN:HD22	0.88	0.86
1:B:436:ALA:N	6:B:773:HOH:O	2.08	0.86
1:B:83:ARG:HH22	1:B:144:LYS:CD	1.86	0.85
1:B:433:ILE:O	6:B:773:HOH:O	1.96	0.83
1:A:113:ASP:HB3	1:B:113:ASP:HB3	1.60	0.82
1:A:48:ILE:HG23	1:A:80:ILE:HD13	1.59	0.81
1:B:325:ASN:ND2	3:B:501:VDO:OAY	2.14	0.81
1:A:70:VAL:CG1	1:A:72:MET:HE2	2.12	0.80
1:A:70:VAL:HG11	1:A:72:MET:CE	2.11	0.78
1:B:401:ASP:OD2	6:B:783:HOH:O	2.03	0.77
1:A:70:VAL:HG11	1:A:72:MET:HE1	1.65	0.77
1:A:176:PRO:HG3	1:A:475:HIS:CE1	2.20	0.77
1:A:88:ASP:OD1	6:A:843:HOH:O	1.96	0.77
1:A:74:LEU:HB2	1:A:78:ARG:O	1.87	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:143:LEU:HD13	1:B:171:GLN:HG3	1.69	0.75
1:B:66:ASN:CB	1:B:70:VAL:HG12	2.17	0.74
1:A:70:VAL:HG12	1:A:72:MET:CE	2.17	0.74
1:A:412:GLU:HG3	4:A:503:EDO:H22	1.70	0.73
1:A:143:LEU:HD13	1:A:171:GLN:HG3	1.70	0.73
1:B:176:PRO:HG3	1:B:475:HIS:CE1	2.24	0.72
1:A:57:GLU:OE2	1:A:60:ARG:NH2	2.20	0.71
1:B:434:SER:C	6:B:773:HOH:O	2.29	0.71
1:A:16:THR:HG21	1:A:29:TRP:CZ2	2.25	0.71
1:A:228:ALA:O	6:A:816:HOH:O	2.08	0.70
1:B:66:ASN:HB2	1:B:70:VAL:HG12	1.71	0.70
1:A:401:ASP:OD2	6:A:806:HOH:O	2.08	0.70
1:A:83:ARG:HH22	1:A:144:LYS:CD	1.95	0.69
1:B:361:ASN:HB2	6:B:791:HOH:O	1.91	0.69
1:A:37:ASN:N	1:A:37:ASN:ND2	2.33	0.69
1:B:437:LEU:O	6:B:805:HOH:O	2.09	0.69
3:A:502:VDO:HAM	3:A:502:VDO:OAO	1.93	0.68
1:B:453:ASP:OD1	1:B:456:ARG:NH2	2.26	0.68
3:B:501:VDO:OAO	3:B:501:VDO:HAM	1.93	0.68
1:A:20:ASP:OD1	1:A:21:PRO:HD2	1.93	0.68
1:A:41:GLU:HG2	1:A:78:ARG:NH1	2.09	0.67
1:A:101:ALA:HA	1:A:105:TRP:HB3	1.77	0.67
1:B:151:ASP:OD1	6:B:819:HOH:O	2.12	0.67
1:A:290:ARG:HD2	1:A:325:ASN:ND2	2.10	0.66
1:A:417:TYR:CE1	1:A:454:ALA:HB2	2.31	0.65
1:A:360:ASP:OD2	1:A:361:ASN:N	2.29	0.65
1:A:51:ILE:HD12	1:A:51:ILE:N	2.11	0.65
1:A:417:TYR:CD1	1:A:454:ALA:CB	2.80	0.65
1:B:241:MET:HG3	6:B:818:HOH:O	1.96	0.64
1:B:246:ARG:NH1	6:B:776:HOH:O	1.98	0.64
1:B:18:ASP:O	1:B:27:ARG:N	2.23	0.64
1:A:233:ASP:OD1	6:A:829:HOH:O	2.15	0.64
1:B:37:ASN:OD1	1:B:37:ASN:N	2.30	0.63
1:B:98:PHE:CE1	1:B:103:LEU:HD13	2.33	0.63
1:B:16:THR:HG21	1:B:29:TRP:CZ2	2.33	0.63
1:A:30:LEU:HD12	1:A:80:ILE:HG21	1.80	0.63
1:B:218:ARG:HD2	4:B:503:EDO:H12	1.80	0.63
5:A:504:CL:CL	5:B:504:CL:CL	2.90	0.63
1:B:51:ILE:N	1:B:51:ILE:HD12	2.14	0.62
1:A:99:MET:CE	1:A:104:MET:HG3	2.29	0.61
1:B:48:ILE:C	1:B:80:ILE:HD12	2.21	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:360:ASP:OD2	1:B:361:ASN:N	2.34	0.61
1:A:97:ASN:HA	1:A:100:THR:HG23	1.83	0.61
1:A:417:TYR:CE1	1:A:454:ALA:CB	2.85	0.60
1:A:101:ALA:C	6:A:863:HOH:O	2.34	0.59
1:B:18:ASP:OD1	1:B:27:ARG:HB2	2.02	0.59
1:A:99:MET:CE	1:A:104:MET:CG	2.81	0.59
1:B:56:SER:O	1:B:60:ARG:CG	2.46	0.59
1:A:292:ASP:OD1	1:A:293:PRO:HD2	2.03	0.58
1:B:143:LEU:CD1	1:B:171:GLN:HG3	2.34	0.58
1:A:4:SER:HA	6:A:732:HOH:O	2.03	0.58
1:B:434:SER:O	6:B:773:HOH:O	2.16	0.58
1:A:66:ASN:CB	1:A:70:VAL:HG23	2.34	0.58
1:A:143:LEU:CD1	1:A:171:GLN:HG3	2.34	0.57
1:A:263:THR:HG21	6:A:748:HOH:O	2.03	0.57
1:B:437:LEU:N	6:B:773:HOH:O	1.88	0.57
1:B:48:ILE:HG21	1:B:80:ILE:HD13	1.75	0.56
1:B:335:TYR:O	1:B:339:VAL:HG23	2.05	0.56
1:A:97:ASN:HA	1:A:100:THR:CG2	2.36	0.56
1:B:41:GLU:O	1:B:42:GLN:C	2.43	0.56
1:A:478:ARG:NH1	6:A:852:HOH:O	1.97	0.56
1:A:99:MET:HE1	1:A:104:MET:CG	2.30	0.55
1:B:98:PHE:HD2	1:B:99:MET:CE	2.20	0.55
1:A:159:TYR:O	1:A:162:VAL:HG22	2.07	0.55
1:A:48:ILE:CG2	1:A:80:ILE:HD13	2.35	0.54
1:A:18:ASP:OD1	1:A:27:ARG:HB2	2.08	0.54
1:B:41:GLU:HA	1:B:78:ARG:NH1	2.21	0.54
3:B:501:VDO:OAZ	6:B:740:HOH:O	2.18	0.54
1:A:335:TYR:O	1:A:339:VAL:HG23	2.08	0.54
1:A:48:ILE:HG23	1:A:80:ILE:CD1	2.32	0.54
1:A:38:VAL:CG1	1:A:464:VAL:HG21	2.38	0.54
1:B:323:ASN:HB2	6:B:684:HOH:O	2.08	0.53
1:A:241:MET:HG3	6:A:606:HOH:O	2.08	0.53
1:B:20:ASP:OD2	1:B:23:THR:HG23	2.09	0.53
1:A:323:ASN:HB2	6:A:706:HOH:O	2.09	0.53
1:B:20:ASP:OD1	1:B:21:PRO:HD2	2.09	0.53
1:A:463:TRP:O	1:A:467:GLN:HG3	2.10	0.52
1:B:324:PRO:HG3	1:B:359:ASN:OD1	2.08	0.52
1:B:48:ILE:CG1	1:B:80:ILE:CD1	2.83	0.52
1:B:292:ASP:OD1	1:B:293:PRO:HD2	2.09	0.52
1:B:240:ALA:O	1:B:241:MET:HB2	2.09	0.52
1:A:65:LEU:HD12	1:A:65:LEU:N	2.25	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:74:LEU:CB	1:A:78:ARG:O	2.58	0.51
1:A:97:ASN:CA	1:A:100:THR:HG23	2.40	0.51
1:A:70:VAL:O	1:A:81:LEU:HD12	2.10	0.51
1:A:48:ILE:O	1:A:80:ILE:HD12	2.11	0.51
1:B:34:GLY:O	1:B:38:VAL:HG23	2.11	0.51
1:B:38:VAL:O	1:B:42:GLN:NE2	2.41	0.51
1:B:61:ARG:O	1:B:65:LEU:HB2	2.11	0.51
1:B:98:PHE:HD2	1:B:99:MET:HE3	1.75	0.51
1:B:159:TYR:O	1:B:162:VAL:HG22	2.11	0.50
1:A:101:ALA:HB3	1:A:329:VAL:HG11	1.92	0.50
1:A:18:ASP:O	1:A:27:ARG:N	2.28	0.50
1:B:441:PRO:HD2	6:B:699:HOH:O	2.10	0.50
1:B:239:GLU:OE1	1:B:239:GLU:N	2.45	0.50
1:B:48:ILE:C	1:B:80:ILE:CD1	2.80	0.50
1:A:293:PRO:HA	1:A:335:TYR:CD1	2.46	0.49
1:B:338:ARG:NH2	6:B:639:HOH:O	2.46	0.49
1:B:15:ILE:HG21	1:B:17:TYR:CZ	2.48	0.49
1:B:35:THR:O	1:B:39:VAL:HG23	2.11	0.49
1:A:392:ALA:N	1:A:393:PRO:HD2	2.27	0.49
1:B:463:TRP:O	1:B:467:GLN:HG3	2.13	0.49
1:B:385:GLN:O	1:B:386:ASN:CB	2.60	0.49
1:B:66:ASN:HB2	1:B:70:VAL:CG1	2.41	0.49
1:A:326:ARG:HH21	3:A:502:VDO:PBA	2.36	0.49
1:A:65:LEU:N	1:A:65:LEU:CD1	2.76	0.48
1:B:407:THR:HA	4:B:503:EDO:O1	2.13	0.48
1:B:97:ASN:O	1:B:97:ASN:OD1	2.32	0.48
1:A:183:ILE:HB	1:A:184:PRO:CD	2.44	0.48
6:A:626:HOH:O	1:B:114:ARG:HD2	2.14	0.47
1:A:385:GLN:O	1:A:386:ASN:CB	2.63	0.47
1:B:260:SER:O	1:B:263:THR:HG23	2.14	0.47
1:A:202:ILE:O	1:A:206:MET:HG3	2.13	0.47
1:A:417:TYR:CD1	1:A:454:ALA:HB2	2.50	0.47
1:A:15:ILE:HG21	1:A:17:TYR:CZ	2.51	0.46
1:B:96:GLN:O	1:B:100:THR:CG2	2.52	0.46
1:B:98:PHE:CD2	1:B:99:MET:CE	2.99	0.46
1:B:202:ILE:O	1:B:206:MET:HG3	2.16	0.46
1:B:16:THR:HB	1:B:29:TRP:CE2	2.51	0.45
1:B:234:ALA:O	1:B:235:ARG:HD3	2.16	0.45
1:A:38:VAL:HG11	1:A:464:VAL:HG21	1.99	0.45
1:B:48:ILE:CB	1:B:80:ILE:CD1	2.94	0.45
1:B:245:TRP:O	1:B:248:HIS:HB2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:51:ILE:CD1	1:A:51:ILE:N	2.79	0.45
1:B:72:MET:HG3	1:B:82:VAL:HG23	1.96	0.45
1:A:475:HIS:HD2	1:A:475:HIS:O	1.99	0.45
1:B:293:PRO:HA	1:B:335:TYR:CD1	2.51	0.45
1:A:325:ASN:C	1:A:325:ASN:OD1	2.55	0.45
1:B:360:ASP:C	1:B:360:ASP:OD2	2.54	0.45
1:B:233:ASP:OD1	6:B:775:HOH:O	2.21	0.45
1:B:262:LEU:HD22	1:B:262:LEU:N	2.32	0.45
1:A:76:SER:HB3	1:A:78:ARG:HD2	1.99	0.44
1:B:65:LEU:HA	1:B:65:LEU:HD23	1.67	0.44
1:B:183:ILE:HB	1:B:184:PRO:CD	2.47	0.44
1:B:437:LEU:C	6:B:805:HOH:O	2.54	0.44
1:A:100:THR:C	1:A:102:ASN:N	2.69	0.44
1:A:91:VAL:HG23	6:A:843:HOH:O	2.17	0.44
1:B:48:ILE:HG12	1:B:80:ILE:HD13	1.94	0.44
1:A:260:SER:O	1:A:263:THR:HG23	2.18	0.44
1:B:475:HIS:O	1:B:475:HIS:HD2	2.01	0.44
1:B:68:ASP:OD1	1:B:144:LYS:NZ	2.49	0.43
1:A:322:MET:CE	1:A:335:TYR:CE2	3.01	0.43
1:A:48:ILE:HG12	1:A:80:ILE:HD13	2.00	0.43
1:B:271:LEU:HD12	1:B:276:GLU:HB2	2.00	0.43
1:A:113:ASP:O	1:A:114:ARG:HB2	2.18	0.43
1:A:290:ARG:HD2	1:A:325:ASN:HD22	1.81	0.43
1:B:322:MET:HE3	1:B:335:TYR:CE2	2.53	0.43
1:A:114:ARG:HG2	1:B:424:PHE:CZ	2.54	0.43
1:B:288:SER:HA	1:B:321:ARG:O	2.18	0.43
1:A:123:ASP:OD1	1:A:124:ALA:N	2.52	0.43
1:B:113:ASP:O	1:B:114:ARG:HB2	2.19	0.43
1:B:23:THR:OG1	1:B:25:GLU:CG	2.66	0.43
1:B:235:ARG:HD3	1:B:235:ARG:HA	1.76	0.43
1:A:245:TRP:O	1:A:248:HIS:HB2	2.19	0.43
1:A:70:VAL:HG12	1:A:72:MET:HE2	1.88	0.43
1:B:263:THR:HG21	6:B:649:HOH:O	2.18	0.42
1:B:97:ASN:C	1:B:97:ASN:OD1	2.58	0.42
1:A:130:ASP:OD1	1:A:133:ARG:NH1	2.52	0.42
1:B:39:VAL:O	1:B:45:VAL:HG21	2.18	0.42
1:B:60:ARG:HG2	1:B:60:ARG:H	1.61	0.42
1:A:234:ALA:O	1:A:235:ARG:HD3	2.19	0.42
1:B:334:ASP:OD2	1:B:338:ARG:HD3	2.20	0.42
1:B:435:ALA:C	6:B:773:HOH:O	2.51	0.42
1:A:360:ASP:OD2	1:A:360:ASP:C	2.58	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:15:ILE:HD11	1:B:82:VAL:HG11	2.00	0.42
1:A:100:THR:O	1:A:102:ASN:N	2.53	0.42
1:A:322:MET:CE	1:A:335:TYR:HE2	2.32	0.42
1:B:322:MET:CE	1:B:335:TYR:HE2	2.33	0.42
1:B:51:ILE:N	1:B:51:ILE:CD1	2.83	0.42
1:B:166:ALA:O	1:B:170:GLU:HG3	2.20	0.42
1:A:334:ASP:OD2	1:A:338:ARG:HD3	2.20	0.41
1:B:101:ALA:HB3	1:B:329:VAL:HG11	2.02	0.41
1:A:38:VAL:HG12	1:A:464:VAL:HG21	2.02	0.41
6:A:681:HOH:O	1:B:192:ARG:HD2	2.20	0.41
1:B:392:ALA:N	1:B:393:PRO:HD2	2.35	0.41
1:A:235:ARG:HA	1:A:235:ARG:HD3	1.80	0.41
1:A:240:ALA:O	1:A:241:MET:HB2	2.21	0.41
1:A:114:ARG:HD2	6:B:609:HOH:O	2.19	0.41
1:B:456:ARG:N	1:B:457:PRO:CD	2.84	0.41
3:A:502:VDO:HAA	3:A:502:VDO:OAT	2.19	0.41
1:A:99:MET:HE2	1:A:104:MET:CG	2.50	0.41
1:B:23:THR:OG1	1:B:25:GLU:HG2	2.20	0.41
1:A:385:GLN:O	1:A:386:ASN:HB3	2.21	0.41
1:B:322:MET:CE	1:B:335:TYR:CE2	3.04	0.41
1:A:100:THR:C	1:A:102:ASN:H	2.24	0.41
1:A:70:VAL:O	1:A:70:VAL:HG12	2.21	0.41
1:A:96:GLN:O	1:A:100:THR:HG22	2.15	0.40
1:A:427:VAL:HG11	1:B:125:ARG:CZ	2.51	0.40
1:B:385:GLN:O	1:B:386:ASN:HB3	2.22	0.40
1:A:475:HIS:C	1:A:475:HIS:CD2	2.94	0.40
1:A:371:ARG:NH2	6:A:850:HOH:O	2.26	0.40
1:A:125:ARG:CZ	1:B:427:VAL:HG11	2.51	0.40
1:B:456:ARG:HH11	1:B:456:ARG:HD3	1.67	0.40
1:B:30:LEU:HA	1:B:30:LEU:HD23	1.89	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	468/497 (94%)	456 (97%)	11 (2%)	1 (0%)	47	49
1	B	464/497 (93%)	450 (97%)	13 (3%)	1 (0%)	47	49
All	All	932/994 (94%)	906 (97%)	24 (3%)	2 (0%)	47	49

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	386	ASN
1	B	386	ASN

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	376/395 (95%)	368 (98%)	8 (2%)	53	59
1	B	374/395 (95%)	367 (98%)	7 (2%)	57	63
All	All	750/790 (95%)	735 (98%)	15 (2%)	55	60

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	ASN
1	A	72	MET
1	A	103	LEU
1	A	158	ASP
1	A	353	ASP
1	A	383	ASP
1	A	397	GLU
1	A	398	ARG
1	B	37	ASN
1	B	103	LEU
1	B	158	ASP

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Mol	Chain	Res	Type
1	B	273	GLU
1	B	383	ASP
1	B	397	GLU
1	B	398	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	37	ASN
1	A	475	HIS
1	B	475	HIS

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 2 are monoatomic - leaving 6 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	$\# Z > 2$
4	EDO	B	503	-	3,3,3	0.52	0	2,2,2	0.44	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	VDO	B	501	-	28,28,28	0.63	0	31,42,42	1.07	2 (6%)
2	GDP	A	501	-	24,30,30	1.10	1 (4%)	31,47,47	2.10	9 (29%)
3	VDO	A	502	-	28,28,28	0.62	0	31,42,42	1.06	1 (3%)
4	EDO	A	503	-	3,3,3	0.21	0	2,2,2	1.16	0
2	GDP	B	502	-	24,30,30	1.15	2 (8%)	31,47,47	2.25	7 (22%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	B	503	-	-	0/1/1/1	-
3	VDO	B	501	-	-	5/12/52/52	0/2/2/2
2	GDP	A	501	-	-	5/12/32/32	0/3/3/3
3	VDO	A	502	-	-	4/12/52/52	0/2/2/2
4	EDO	A	503	-	-	1/1/1/1	-
2	GDP	B	502	-	-	4/12/32/32	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	GDP	C6-C5	3.74	1.47	1.41
2	B	502	GDP	C6-C5	2.80	1.46	1.41
2	B	502	GDP	PB-O1B	2.01	1.57	1.50

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	502	GDP	C6-N1-C2	5.64	124.89	115.93
2	A	501	GDP	C2-N3-C4	5.05	121.12	115.36
2	B	502	GDP	C5-C6-N1	-4.78	116.90	123.43
2	A	501	GDP	C6-N1-C2	4.42	122.95	115.93
2	B	502	GDP	O4'-C1'-C2'	4.26	113.16	106.93
2	A	501	GDP	N3-C2-N1	-4.15	121.68	127.22
2	A	501	GDP	C6-C5-C4	-4.03	116.95	120.80
2	B	502	GDP	N3-C2-N1	-4.02	121.86	127.22
2	B	502	GDP	C6-C5-C4	-3.92	117.05	120.80
2	B	502	GDP	C2-N3-C4	3.90	119.81	115.36
2	A	501	GDP	C5-C6-N1	-3.87	118.14	123.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	502	VDO	CAF-CAV-CAA	3.68	113.11	108.49
3	B	501	VDO	CAF-CAV-CAA	3.60	113.00	108.49
2	B	502	GDP	O3B-PB-O3A	3.47	116.29	104.64
2	A	501	GDP	C4-C5-N7	-3.36	105.90	109.40
2	A	501	GDP	O3B-PB-O3A	2.62	113.42	104.64
3	B	501	VDO	CAM-NAN-CAA	2.37	120.76	116.47
2	A	501	GDP	O3B-PB-O1B	2.34	119.85	110.68
2	A	501	GDP	N2-C2-N1	2.06	120.45	117.25

There are no chirality outliers.

All (19) torsion outliers are listed below:

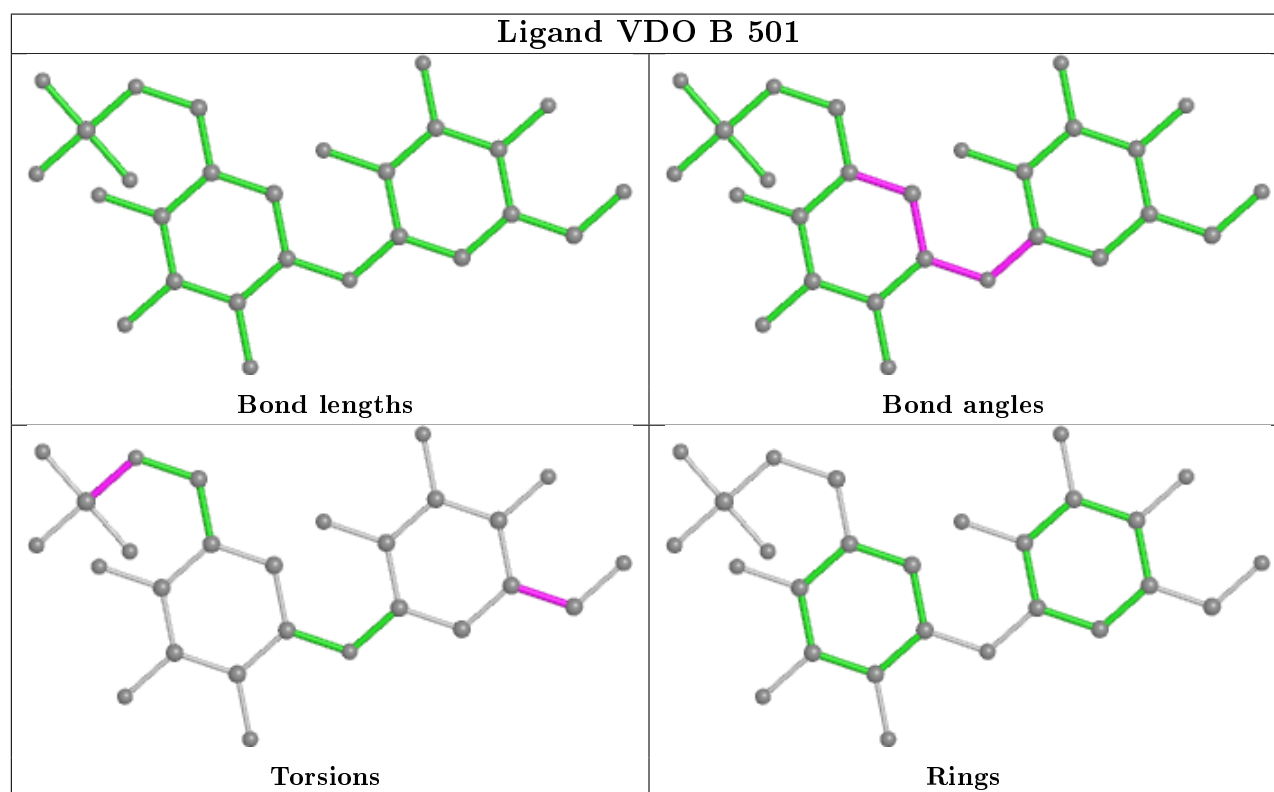
Mol	Chain	Res	Type	Atoms
3	B	501	VDO	CAG-OAW-PBA-OAY
3	B	501	VDO	CAG-OAW-PBA-OAZ
3	B	501	VDO	CAG-OAW-PBA-OAX
3	B	501	VDO	OAQ-CAD-CAK-CAL
3	B	501	VDO	OAQ-CAD-CAK-CAH
2	A	501	GDP	PB-O3A-PA-O5'
2	A	501	GDP	O4'-C4'-C5'-O5'
3	A	502	VDO	CAG-OAW-PBA-OAY
3	A	502	VDO	CAG-OAW-PBA-OAZ
3	A	502	VDO	CAG-OAW-PBA-OAX
2	B	502	GDP	O4'-C4'-C5'-O5'
2	A	501	GDP	C3'-C4'-C5'-O5'
2	B	502	GDP	C3'-C4'-C5'-O5'
4	A	503	EDO	O1-C1-C2-O2
2	B	502	GDP	PA-O3A-PB-O1B
2	B	502	GDP	PB-O3A-PA-O5'
2	A	501	GDP	PA-O3A-PB-O2B
3	A	502	VDO	CAF-CAG-OAW-PBA
2	A	501	GDP	PA-O3A-PB-O1B

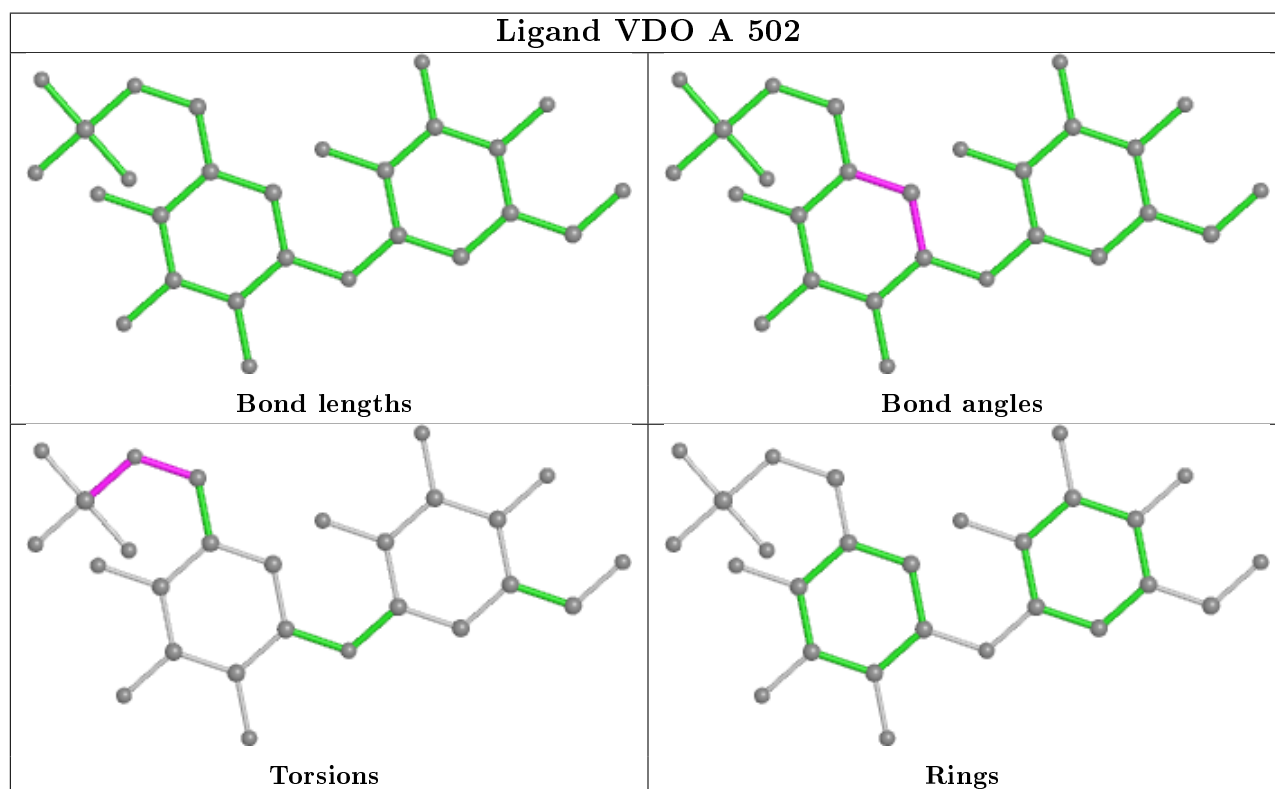
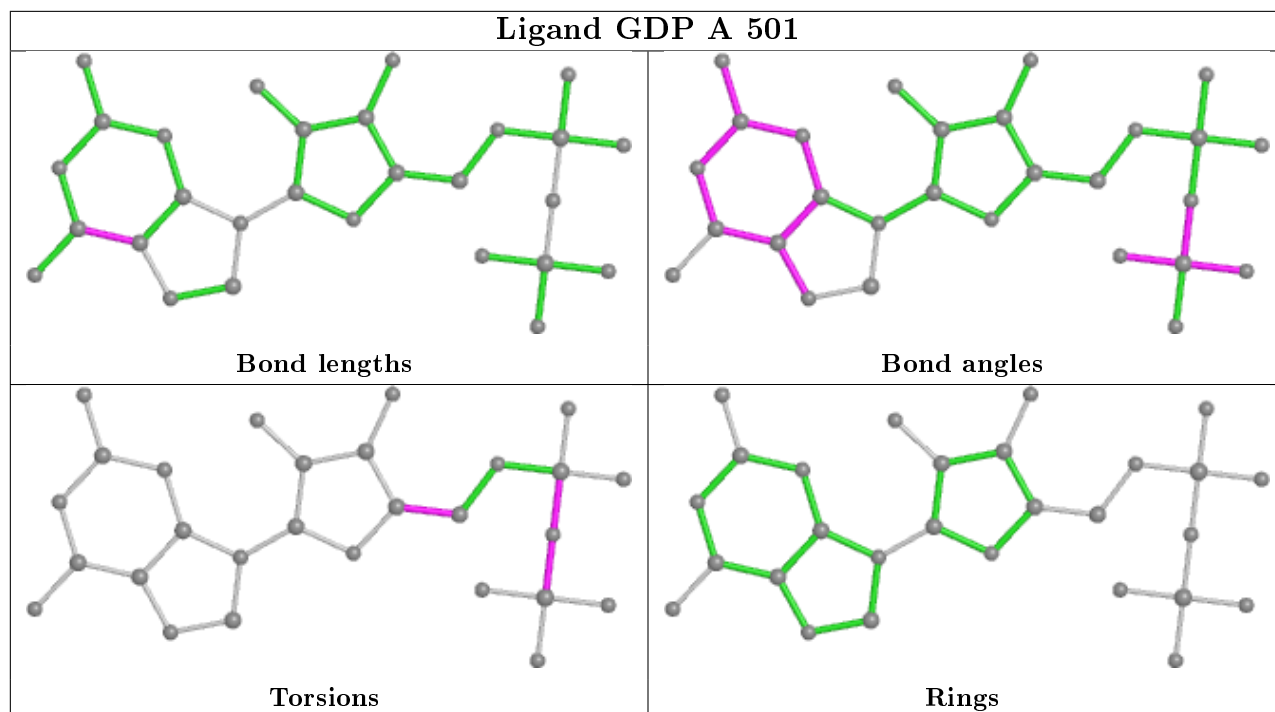
There are no ring outliers.

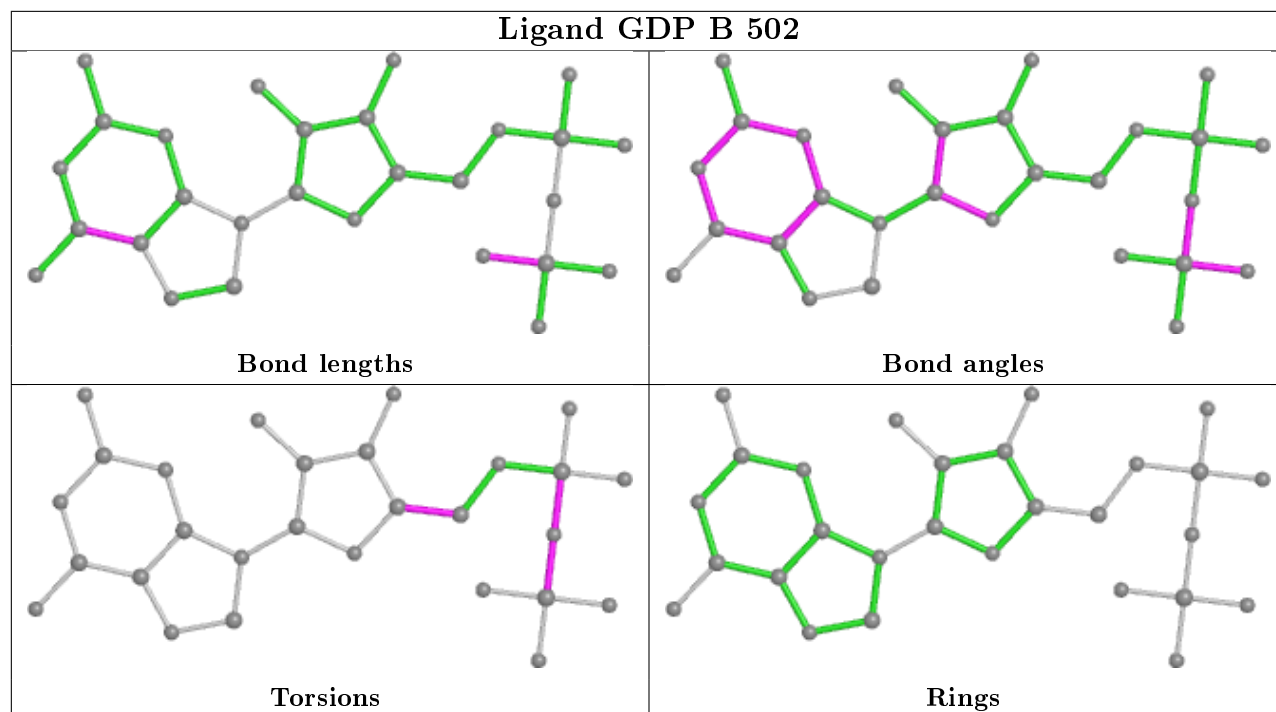
4 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	503	EDO	2	0
3	B	501	VDO	3	0
3	A	502	VDO	3	0
4	A	503	EDO	1	0

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







5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	472/497 (94%)	-0.14	12 (2%) 57 62	19, 35, 72, 95	0
1	B	470/497 (94%)	-0.08	19 (4%) 38 44	20, 36, 72, 87	0
All	All	942/994 (94%)	-0.11	31 (3%) 46 53	19, 35, 72, 95	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	31	ALA	5.2
1	B	43	ALA	5.0
1	B	45	VAL	4.2
1	B	30	LEU	3.7
1	A	101	ALA	3.5
1	A	43	ALA	3.3
1	B	3	GLY	3.3
1	A	45	VAL	3.3
1	A	31	ALA	3.1
1	B	73	GLU	3.1
1	A	3	GLY	3.0
1	A	30	LEU	2.9
1	B	13	ALA	2.8
1	B	29	TRP	2.8
1	B	38	VAL	2.8
1	B	74	LEU	2.6
1	A	73	GLU	2.6
1	A	77	GLY	2.5
1	A	14	ALA	2.5
1	B	14	ALA	2.5
1	B	28	ALA	2.3
1	A	100	THR	2.2
1	A	29	TRP	2.2
1	B	17	TYR	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	64	ALA	2.2
1	B	65	LEU	2.2
1	A	38	VAL	2.1
1	B	4	SER	2.1
1	B	59	ASP	2.0
1	B	15	ILE	2.0
1	B	159	TYR	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 carbohydrates in this entry.

6.4 Ligands [i](#)

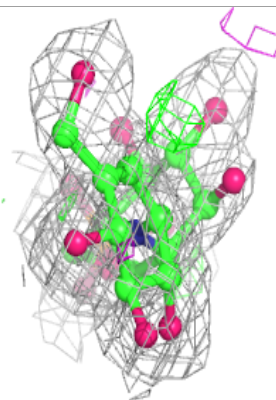
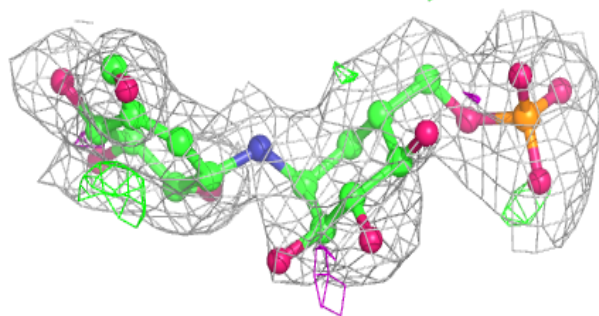
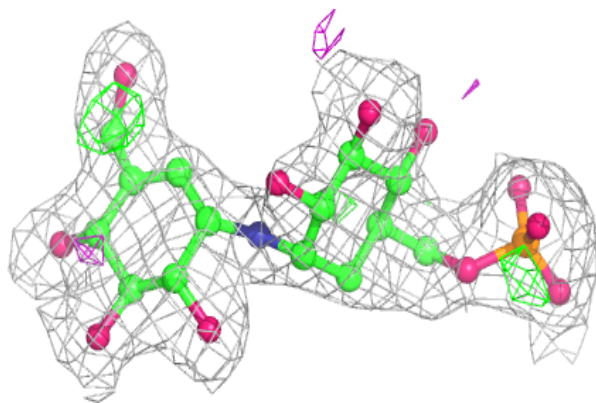
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	VDO	B	501	27/27	0.89	0.17	28,42,71,77	0
3	VDO	A	502	27/27	0.91	0.17	24,40,58,70	0
4	EDO	B	503	4/4	0.94	0.14	23,34,35,44	0
4	EDO	A	503	4/4	0.96	0.17	23,31,32,46	0
2	GDP	A	501	28/28	0.98	0.08	23,29,34,41	0
2	GDP	B	502	28/28	0.98	0.09	25,32,37,41	0
5	CL	A	504	1/1	0.99	0.10	26,26,26,26	0
5	CL	B	504	1/1	1.00	0.11	29,29,29,29	0

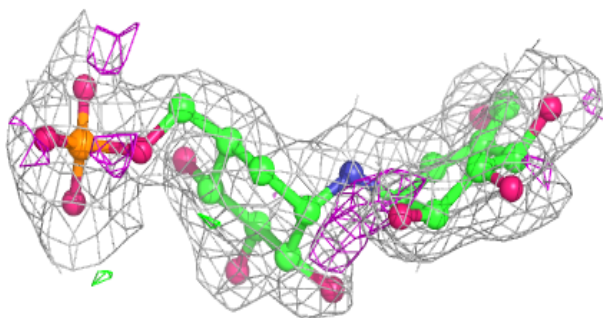
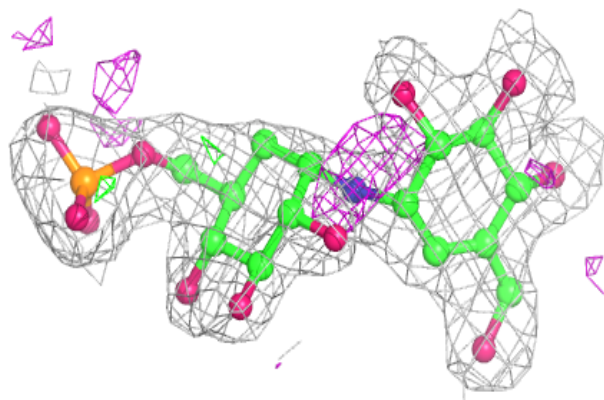
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around VDO B 501:

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

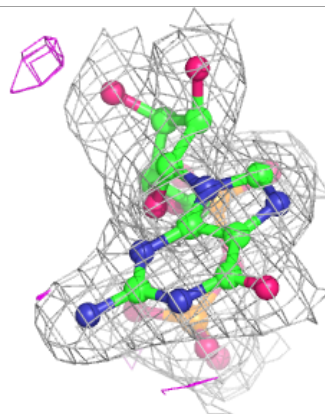
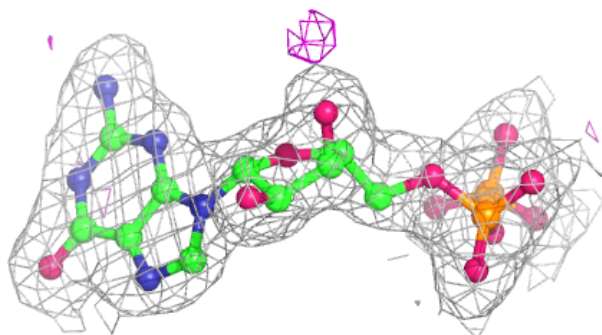
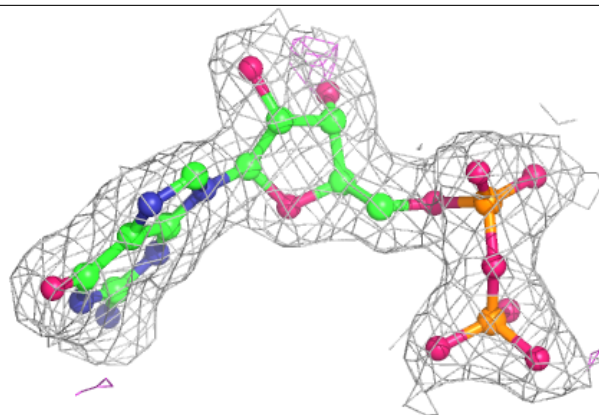
**Electron density around VDO A 502:**

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

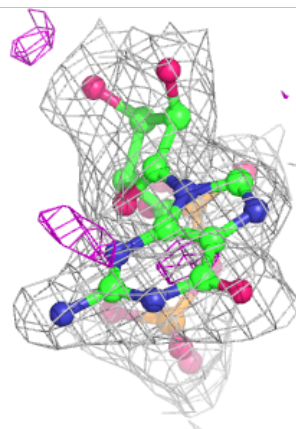
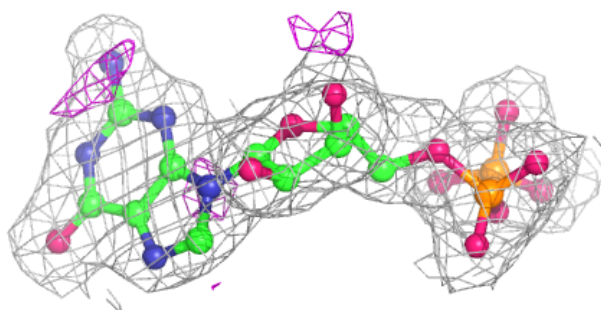
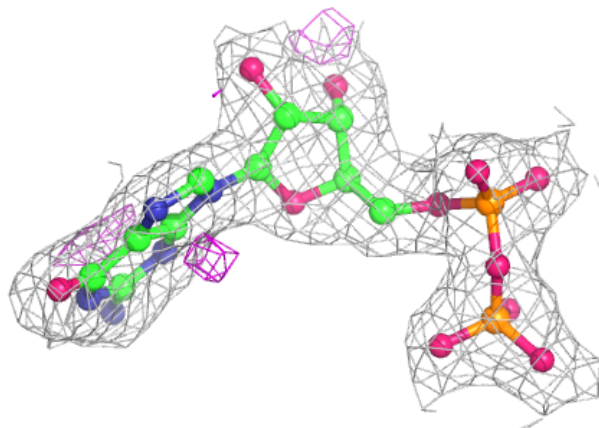


Electron density around GDP A 501:

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

**Electron density around GDP B 502:**

$2mF_o-DF_c$ (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.