



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 12, 2024 – 04:27 AM EDT

PDB ID : 1IHY
Title : GAPDH complexed with ADP-ribose
Authors : Shen, Y.-Q.; Song, S.-Y.; Lin, Z.-J.
Deposited on : 2001-04-20
Resolution : 3.00 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.36.2
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)	:	2.36.2

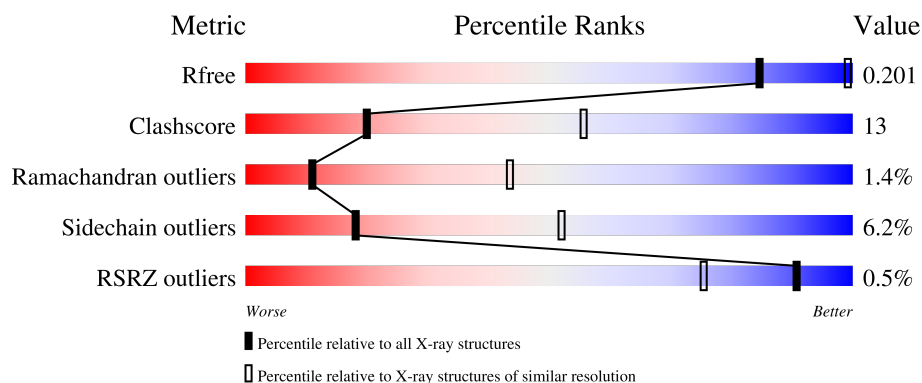
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (3.00-3.00)

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 $\leq 5\%$. 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	333	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 24%, green 73%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 73% 24% </div> </div>
1	B	333	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, orange 1%, yellow 23%, green 74%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 74% 23% </div> </div>
1	C	333	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, orange 1%, yellow 22%, green 74%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 74% 22% </div> </div>
1	D	333	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, orange 1%, yellow 21%, green 75%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 75% 21% </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10317 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 GLYCERALDEHYDE 3-PHOSPHATE DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	333	Total	C	N	O	S	0	0	0
			2512	1593	421	482	16			
1	B	333	Total	C	N	O	S	0	0	0
			2512	1593	421	482	16			
1	C	333	Total	C	N	O	S	0	0	0
			2512	1593	421	482	16			
1	D	333	Total	C	N	O	S	0	0	0
			2512	1593	421	482	16			

There are 12 discrepancies between the modelled and reference sequences:

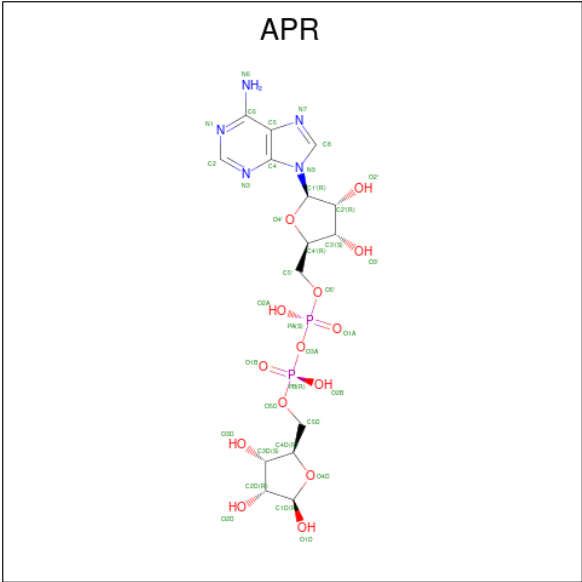
Chain	Residue	Modelled	Actual	Comment	Reference
A	18	THR	ALA	SEE REMARK 999	UNP P56649
A	59	VAL	ALA	SEE REMARK 999	UNP P56649
A	261	THR	ALA	SEE REMARK 999	UNP P56649
B	18	THR	ALA	SEE REMARK 999	UNP P56649
B	59	VAL	ALA	SEE REMARK 999	UNP P56649
B	261	THR	ALA	SEE REMARK 999	UNP P56649
C	18	THR	ALA	SEE REMARK 999	UNP P56649
C	59	VAL	ALA	SEE REMARK 999	UNP P56649
C	261	THR	ALA	SEE REMARK 999	UNP P56649
D	18	THR	ALA	SEE REMARK 999	UNP P56649
D	59	VAL	ALA	SEE REMARK 999	UNP P56649
D	261	THR	ALA	SEE REMARK 999	UNP P56649

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

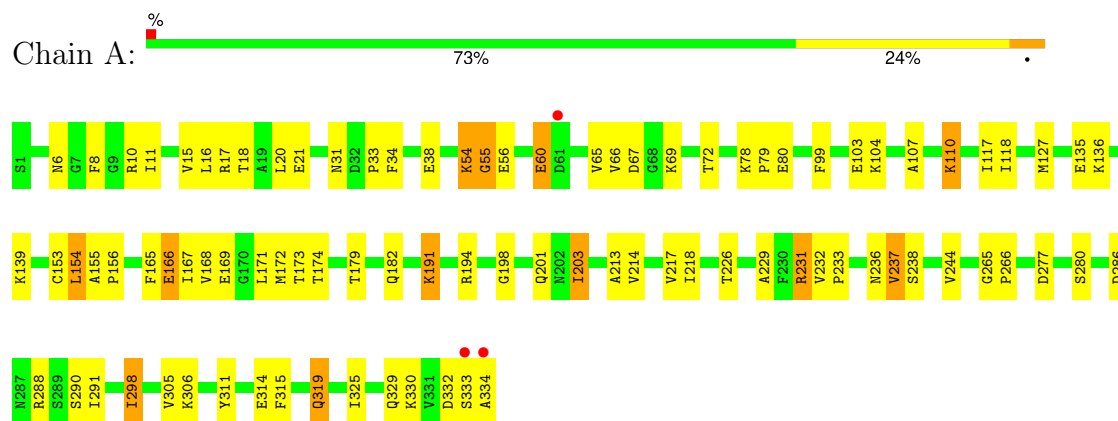
- Molecule 3 is ADENOSINE-5-DIPHOSPHORIBOSE (three-letter code: APR) (formula: $C_{15}H_{23}N_5O_{14}P_2$).



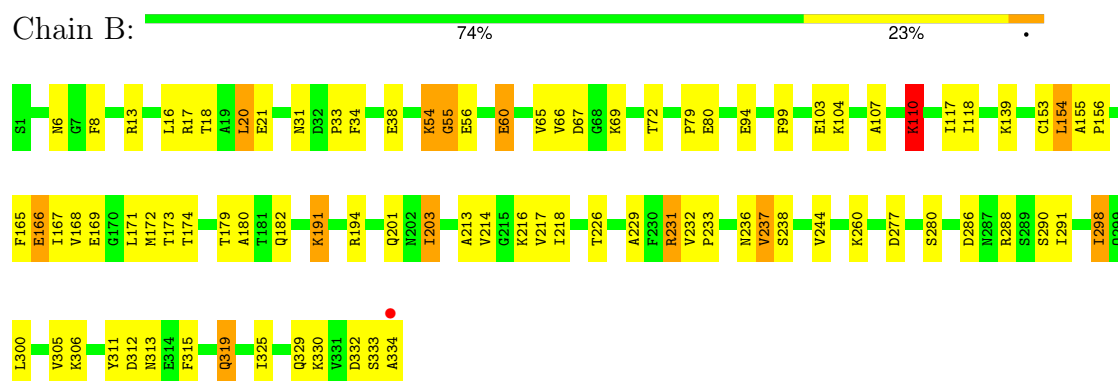
3 Residue-property plots [i](#)

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.

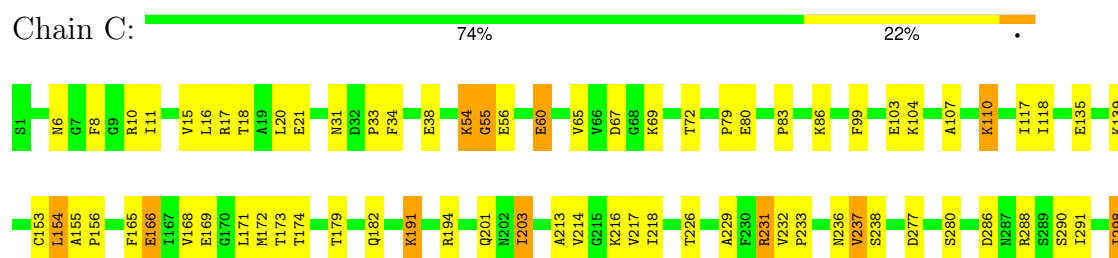
• Molecule 1: GLYCERALDEHYDE 3-PHOSPHATE DEHYDROGENASE

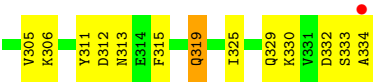


• Molecule 1: GLYCERALDEHYDE 3-PHOSPHATE DEHYDROGENASE



• Molecule 1: GLYCERALDEHYDE 3-PHOSPHATE DEHYDROGENASE





● Molecule 1: GLYCERALDEHYDE 3-PHOSPHATE DEHYDROGENASE



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	152.80Å 100.35Å 128.31Å 90.00° 110.28° 90.00°	Depositor
Resolution (Å)	24.37 – 3.00 24.37 – 3.00	Depositor EDS
% Data completeness (in resolution range)	90.0 (24.37-3.00) 90.1 (24.37-3.00)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.59 (at 2.99Å)	Xtriage
Refinement program	CNS 0.9	Depositor
R, R_{free}	0.176 , 0.207 0.171 , 0.201	Depositor DCC
R_{free} test set	3295 reflections (9.97%)	wwPDB-VP
Wilson B-factor (Å ²)	36.9	Xtriage
Anisotropy	0.272	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 54.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	10317	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 17.12% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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: SO4, APR

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	0.52	0/2557	0.90	5/3455 (0.1%)
1	B	0.50	0/2557	0.69	5/3455 (0.1%)
1	C	0.50	0/2557	0.67	3/3455 (0.1%)
1	D	0.51	0/2557	0.67	3/3455 (0.1%)
All	All	0.51	0/10228	0.74	16/13820 (0.1%)

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	231	ARG	NE-CZ-NH2	-25.65	107.47	120.30
1	A	231	ARG	NE-CZ-NH1	23.67	132.14	120.30
1	A	231	ARG	CD-NE-CZ	10.79	138.70	123.60
1	C	231	ARG	NE-CZ-NH1	-7.74	116.43	120.30
1	B	110	LYS	CB-CG-CD	7.62	131.41	111.60
1	D	231	ARG	NE-CZ-NH1	-7.38	116.61	120.30
1	A	231	ARG	CG-CD-NE	-7.22	96.64	111.80
1	B	231	ARG	NE-CZ-NH1	-6.93	116.83	120.30
1	B	110	LYS	CA-CB-CG	6.88	128.54	113.40
1	D	231	ARG	NE-CZ-NH2	6.21	123.40	120.30
1	C	231	ARG	NE-CZ-NH2	6.17	123.39	120.30
1	B	231	ARG	NE-CZ-NH2	6.15	123.38	120.30
1	C	203	ILE	N-CA-C	-5.75	95.47	111.00
1	A	203	ILE	N-CA-C	-5.55	96.01	111.00
1	B	203	ILE	N-CA-C	-5.45	96.30	111.00
1	D	203	ILE	N-CA-C	-5.34	96.58	111.00

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	2512	0	2537	69	0
1	B	2512	0	2537	70	0
1	C	2512	0	2537	70	0
1	D	2512	0	2537	65	0
2	A	10	0	0	0	0
2	B	10	0	0	0	0
2	C	10	0	0	0	0
2	D	10	0	0	0	0
3	A	36	0	21	3	0
3	B	36	0	21	1	0
3	C	36	0	21	4	0
3	D	36	0	21	3	0
4	A	19	0	0	2	0
4	B	24	0	0	3	0
4	C	18	0	0	1	0
4	D	24	0	0	1	0
All	All	10317	0	10232	257	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 13.

All (257) 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:54:LYS:H	1:B:54:LYS:HD2	1.25	1.00
1:C:54:LYS:H	1:C:54:LYS:HD2	1.25	0.98
1:D:54:LYS:H	1:D:54:LYS:HD2	1.25	0.97
1:A:54:LYS:H	1:A:54:LYS:HD2	1.25	0.96
1:B:154:LEU:HD13	1:B:214:VAL:HG21	1.57	0.87
1:A:154:LEU:HD13	1:A:214:VAL:HG21	1.61	0.82
1:D:154:LEU:HD13	1:D:214:VAL:HG21	1.61	0.82
1:C:154:LEU:HD13	1:C:214:VAL:HG21	1.62	0.81
1:C:83:PRO:HB2	1:C:86:LYS:HD3	1.61	0.81
1:D:179:THR:HB	3:D:337:APR:O2D	1.88	0.74
1:B:182:GLN:OE1	1:B:231:ARG:HD2	1.88	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:191:LYS:HB2	1:C:191:LYS:NZ	2.03	0.74
1:D:191:LYS:NZ	1:D:191:LYS:HB2	2.03	0.74
1:B:191:LYS:HB2	1:B:191:LYS:NZ	2.03	0.73
1:C:182:GLN:OE1	1:C:231:ARG:HD2	1.90	0.71
1:A:191:LYS:HB2	1:A:191:LYS:NZ	2.06	0.71
1:B:280:SER:HB2	1:C:203:ILE:H	1.58	0.69
1:D:182:GLN:OE1	1:D:231:ARG:HD2	1.94	0.68
1:B:168:VAL:HG12	1:B:169:GLU:HG2	1.76	0.68
1:B:277:ASP:HB3	1:C:194:ARG:HG2	1.76	0.68
1:A:203:ILE:H	1:D:280:SER:HB2	1.59	0.68
1:A:179:THR:HB	3:A:335:APR:O2D	1.93	0.67
1:A:277:ASP:HB3	1:D:194:ARG:HG2	1.76	0.67
1:A:194:ARG:HG2	1:D:277:ASP:HB3	1.77	0.66
1:C:153:CYS:HA	1:C:290:SER:HB3	1.76	0.66
1:A:17:ARG:O	1:A:21:GLU:HG2	1.96	0.66
1:C:17:ARG:O	1:C:21:GLU:HG2	1.96	0.65
1:A:280:SER:HB2	1:D:203:ILE:H	1.62	0.65
1:B:203:ILE:H	1:C:280:SER:HB2	1.62	0.65
1:C:54:LYS:H	1:C:54:LYS:CD	1.98	0.65
1:C:286:ASP:OD1	1:C:288:ARG:HD3	1.96	0.65
1:A:182:GLN:HB3	4:A:519:HOH:O	1.97	0.65
1:A:332:ASP:C	1:A:334:ALA:H	2.01	0.64
1:C:80:GLU:HB2	1:C:110:LYS:HG2	1.78	0.64
1:C:86:LYS:NZ	1:C:86:LYS:HB3	2.13	0.64
1:B:153:CYS:HA	1:B:290:SER:HB3	1.78	0.64
1:A:168:VAL:HG12	1:A:169:GLU:HG2	1.79	0.64
1:B:286:ASP:OD1	1:B:288:ARG:HD3	1.97	0.64
1:B:179:THR:HB	3:B:336:APR:O2D	1.98	0.64
1:A:153:CYS:HA	1:A:290:SER:HB3	1.79	0.64
1:D:286:ASP:OD1	1:D:288:ARG:HD3	1.98	0.64
1:D:80:GLU:HB2	1:D:110:LYS:HG2	1.80	0.63
1:D:168:VAL:HG12	1:D:169:GLU:HG2	1.79	0.63
1:B:194:ARG:HG2	1:C:277:ASP:HB3	1.80	0.63
1:B:332:ASP:C	1:B:334:ALA:H	2.02	0.63
1:D:17:ARG:O	1:D:21:GLU:HG2	1.98	0.63
1:D:110:LYS:HE3	1:D:110:LYS:O	1.99	0.63
1:A:99:PHE:HA	1:A:104:LYS:HG3	1.80	0.62
1:D:153:CYS:HA	1:D:290:SER:HB3	1.81	0.62
1:A:286:ASP:OD1	1:A:288:ARG:HD3	1.99	0.62
1:C:179:THR:HB	3:C:338:APR:O2D	2.00	0.62
1:A:80:GLU:HB2	1:A:110:LYS:HG2	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:191:LYS:HB2	1:B:191:LYS:HZ1	1.63	0.61
1:D:99:PHE:HA	1:D:104:LYS:HG3	1.80	0.61
1:D:332:ASP:C	1:D:334:ALA:H	2.03	0.61
1:C:83:PRO:CB	1:C:86:LYS:HD3	2.31	0.61
1:C:332:ASP:C	1:C:334:ALA:H	2.04	0.61
1:C:168:VAL:HG12	1:C:169:GLU:HG2	1.83	0.61
1:B:54:LYS:HD2	1:B:54:LYS:N	2.07	0.61
1:B:233:PRO:HG2	1:C:233:PRO:HG2	1.83	0.61
1:B:17:ARG:O	1:B:21:GLU:HG2	2.01	0.60
1:B:99:PHE:HA	1:B:104:LYS:HG3	1.83	0.60
1:C:99:PHE:HA	1:C:104:LYS:HG3	1.82	0.60
1:B:155:ALA:HB3	1:B:156:PRO:HD3	1.84	0.60
1:B:56:GLU:O	1:B:56:GLU:HG3	2.01	0.60
1:C:110:LYS:HE3	1:C:110:LYS:O	2.01	0.59
1:A:110:LYS:HE3	1:A:110:LYS:O	2.03	0.58
1:C:330:LYS:O	1:C:334:ALA:HB2	2.03	0.58
1:B:330:LYS:O	1:B:334:ALA:HB2	2.04	0.57
1:D:54:LYS:HD2	1:D:54:LYS:N	2.08	0.57
1:D:54:LYS:H	1:D:54:LYS:CD	1.98	0.56
1:B:260:LYS:NZ	4:B:521:HOH:O	2.33	0.56
1:A:233:PRO:HG2	1:D:233:PRO:HG2	1.87	0.56
1:C:54:LYS:HD2	1:C:54:LYS:N	2.08	0.56
1:D:330:LYS:O	1:D:334:ALA:HB2	2.05	0.56
1:D:191:LYS:HB2	1:D:191:LYS:HZ1	1.71	0.56
1:C:56:GLU:O	1:C:56:GLU:HG3	2.05	0.56
1:A:330:LYS:O	1:A:334:ALA:HB2	2.06	0.56
1:A:54:LYS:HD2	1:A:54:LYS:N	2.07	0.55
1:A:155:ALA:HB3	1:A:156:PRO:HD3	1.89	0.55
1:C:191:LYS:HB2	1:C:191:LYS:HZ1	1.72	0.55
1:D:155:ALA:HB3	1:D:156:PRO:HD3	1.88	0.55
1:A:56:GLU:O	1:A:56:GLU:HG3	2.06	0.55
1:B:280:SER:CB	1:C:203:ILE:H	2.18	0.55
1:C:80:GLU:HB2	1:C:110:LYS:CG	2.36	0.55
1:A:54:LYS:H	1:A:54:LYS:CD	1.96	0.55
1:C:86:LYS:HB3	1:C:86:LYS:HZ3	1.73	0.54
1:A:280:SER:CB	1:D:203:ILE:H	2.20	0.54
1:B:80:GLU:HB2	1:B:110:LYS:HD3	1.89	0.54
1:C:155:ALA:HB3	1:C:156:PRO:HD3	1.89	0.54
1:B:54:LYS:H	1:B:54:LYS:CD	1.96	0.54
1:C:291:ILE:N	1:C:291:ILE:HD12	2.23	0.53
1:C:80:GLU:O	1:C:110:LYS:HD3	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:118:ILE:HD12	1:D:118:ILE:N	2.23	0.53
1:A:165:PHE:O	1:A:166:GLU:HB3	2.09	0.52
1:B:80:GLU:O	1:B:110:LYS:HD3	2.10	0.52
1:B:298:ILE:CD1	1:B:306:LYS:HD2	2.38	0.52
1:D:80:GLU:O	1:D:110:LYS:HD3	2.09	0.52
1:D:325:ILE:O	1:D:329:GLN:HG3	2.09	0.52
1:B:118:ILE:HD12	1:B:118:ILE:N	2.25	0.52
1:A:118:ILE:N	1:A:118:ILE:HD12	2.25	0.52
1:D:56:GLU:O	1:D:56:GLU:HG3	2.09	0.52
1:C:298:ILE:CD1	1:C:306:LYS:HD2	2.40	0.52
1:B:171:LEU:HD23	1:B:226:THR:HG23	1.91	0.52
1:D:80:GLU:HB2	1:D:110:LYS:CG	2.39	0.51
1:C:33:PRO:HG2	1:C:34:PHE:CD1	2.44	0.51
1:C:118:ILE:N	1:C:118:ILE:HD12	2.25	0.51
1:D:165:PHE:O	1:D:166:GLU:HB3	2.10	0.51
1:A:127:MET:HE1	1:A:213:ALA:HB1	1.93	0.51
1:A:80:GLU:HB2	1:A:110:LYS:CG	2.41	0.51
1:B:298:ILE:HD11	1:B:306:LYS:HD2	1.92	0.51
1:C:325:ILE:O	1:C:329:GLN:HG3	2.11	0.51
1:B:16:LEU:HD23	1:B:16:LEU:O	2.10	0.51
1:D:291:ILE:N	1:D:291:ILE:HD12	2.26	0.51
1:B:80:GLU:O	1:B:110:LYS:CD	2.58	0.50
1:C:165:PHE:O	1:C:166:GLU:HB3	2.10	0.50
1:D:33:PRO:HG2	1:D:34:PHE:CD1	2.46	0.50
1:B:203:ILE:H	1:C:280:SER:CB	2.23	0.50
1:D:298:ILE:CD1	1:D:306:LYS:HD2	2.42	0.50
1:B:80:GLU:HB3	1:B:107:ALA:HB1	1.94	0.50
1:C:153:CYS:SG	1:C:290:SER:HB2	2.52	0.50
1:C:11:ILE:O	1:C:15:VAL:HG23	2.11	0.50
1:A:332:ASP:C	1:A:334:ALA:N	2.65	0.50
1:B:325:ILE:O	1:B:329:GLN:HG3	2.12	0.50
1:D:172:MET:HG2	1:D:173:THR:N	2.26	0.50
1:D:171:LEU:HD23	1:D:226:THR:HG23	1.93	0.50
1:A:60:GLU:HB2	1:A:65:VAL:HG23	1.93	0.49
1:C:172:MET:HG2	1:C:173:THR:N	2.27	0.49
1:B:33:PRO:HG2	1:B:34:PHE:CD1	2.48	0.49
1:B:165:PHE:O	1:B:166:GLU:HB3	2.11	0.49
1:C:298:ILE:HD11	1:C:306:LYS:HD2	1.93	0.49
1:A:172:MET:HG2	1:A:173:THR:N	2.26	0.49
1:D:60:GLU:HB2	1:D:65:VAL:HG23	1.94	0.49
1:A:291:ILE:N	1:A:291:ILE:HD12	2.27	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:110:LYS:O	1:B:110:LYS:HE3	2.13	0.49
1:A:33:PRO:HG2	1:A:34:PHE:CD1	2.48	0.49
1:D:6:ASN:HD21	3:D:337:APR:H2	1.77	0.49
1:A:136:LYS:HG2	1:A:136:LYS:O	2.12	0.49
1:B:172:MET:HG2	1:B:173:THR:N	2.27	0.49
1:C:10:ARG:HG3	3:C:338:APR:O1D	2.12	0.49
1:D:298:ILE:HD11	1:D:306:LYS:HD2	1.95	0.49
1:D:55:GLY:HA3	1:D:67:ASP:OD2	2.13	0.49
1:A:203:ILE:H	1:D:280:SER:CB	2.22	0.49
1:B:291:ILE:HD12	1:B:291:ILE:N	2.27	0.49
1:C:171:LEU:HD23	1:C:226:THR:HG23	1.94	0.49
1:B:332:ASP:C	1:B:334:ALA:N	2.66	0.48
1:C:80:GLU:HB3	1:C:107:ALA:HB1	1.94	0.48
1:C:117:ILE:HD12	1:C:117:ILE:N	2.28	0.48
1:D:117:ILE:N	1:D:117:ILE:HD12	2.27	0.48
1:D:16:LEU:O	1:D:16:LEU:HD23	2.13	0.48
1:C:60:GLU:HB2	1:C:65:VAL:HG23	1.95	0.48
1:C:86:LYS:NZ	1:C:86:LYS:CB	2.76	0.48
1:D:179:THR:CB	3:D:337:APR:O2D	2.60	0.48
1:A:17:ARG:HD3	4:A:510:HOH:O	2.13	0.48
1:D:80:GLU:HB3	1:D:107:ALA:HB1	1.96	0.48
1:C:305:VAL:HG22	1:C:306:LYS:N	2.28	0.48
1:A:80:GLU:HB3	1:A:107:ALA:HB1	1.96	0.48
1:A:153:CYS:SG	1:A:290:SER:HB2	2.54	0.48
1:D:127:MET:HE1	1:D:213:ALA:HB1	1.95	0.48
1:A:11:ILE:O	1:A:15:VAL:HG23	2.14	0.48
1:A:171:LEU:HD23	1:A:226:THR:HG23	1.95	0.48
1:D:332:ASP:C	1:D:334:ALA:N	2.67	0.48
1:A:232:VAL:HG11	1:D:232:VAL:HG11	1.95	0.47
1:C:6:ASN:ND2	1:C:31:ASN:HD22	2.12	0.47
1:A:298:ILE:CD1	1:A:306:LYS:HD2	2.44	0.47
1:C:217:VAL:HG23	1:C:218:ILE:HG13	1.96	0.47
1:C:332:ASP:C	1:C:334:ALA:N	2.68	0.47
1:D:237:VAL:HG11	1:D:280:SER:O	2.15	0.47
1:B:232:VAL:HG11	1:C:232:VAL:HG11	1.96	0.47
3:C:338:APR:H5'2	4:C:511:HOH:O	2.15	0.46
1:A:6:ASN:ND2	1:A:31:ASN:HD22	2.12	0.46
1:A:325:ILE:O	1:A:329:GLN:HG3	2.13	0.46
1:B:6:ASN:ND2	1:B:31:ASN:HD22	2.11	0.46
1:A:236:ASN:O	1:A:237:VAL:HB	2.15	0.46
1:B:60:GLU:HB2	1:B:65:VAL:HG23	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:171:LEU:HD23	1:B:226:THR:CG2	2.45	0.46
1:B:153:CYS:SG	1:B:290:SER:HB2	2.55	0.46
1:A:16:LEU:O	1:A:16:LEU:HD23	2.14	0.46
1:A:298:ILE:HD13	1:A:306:LYS:HB3	1.98	0.46
1:C:16:LEU:HD23	1:C:16:LEU:O	2.16	0.46
1:B:55:GLY:HA3	1:B:67:ASP:OD2	2.16	0.46
1:B:182:GLN:HB3	4:B:520:HOH:O	2.15	0.46
1:B:305:VAL:HG22	1:B:306:LYS:N	2.31	0.46
1:A:298:ILE:HD11	1:A:306:LYS:HD2	1.97	0.45
1:C:237:VAL:HG11	1:C:280:SER:O	2.17	0.45
1:D:60:GLU:HB2	1:D:65:VAL:CG2	2.47	0.45
1:A:80:GLU:O	1:A:110:LYS:HD3	2.16	0.45
1:B:332:ASP:O	1:B:334:ALA:N	2.49	0.45
1:D:217:VAL:HG23	1:D:218:ILE:HG13	1.98	0.45
1:A:117:ILE:HD12	1:A:117:ILE:N	2.32	0.45
1:C:55:GLY:HA3	1:C:67:ASP:OD2	2.17	0.45
1:B:80:GLU:HB2	1:B:110:LYS:CD	2.47	0.45
1:B:298:ILE:HD13	1:B:306:LYS:HB3	1.99	0.45
1:A:60:GLU:HB2	1:A:65:VAL:CG2	2.47	0.44
1:A:237:VAL:HG11	1:A:280:SER:O	2.17	0.44
1:A:10:ARG:HG3	3:A:335:APR:O1D	2.18	0.44
1:C:135:GLU:OE2	1:C:135:GLU:N	2.44	0.44
1:C:171:LEU:HD23	1:C:226:THR:CG2	2.48	0.44
1:D:6:ASN:ND2	1:D:31:ASN:HD22	2.16	0.44
1:A:217:VAL:HG23	1:A:218:ILE:HG13	1.99	0.44
1:D:171:LEU:HD23	1:D:226:THR:CG2	2.47	0.44
1:A:78:LYS:HE2	1:A:80:GLU:OE1	2.18	0.43
1:D:79:PRO:HD3	1:D:99:PHE:CZ	2.53	0.43
1:B:20:LEU:CD2	1:B:66:VAL:HG11	2.48	0.43
1:B:117:ILE:N	1:B:117:ILE:HD12	2.32	0.43
1:B:238:SER:OG	1:B:313:ASN:ND2	2.51	0.43
1:C:298:ILE:HD13	1:C:306:LYS:HB3	2.01	0.43
1:A:191:LYS:HB2	1:A:191:LYS:HZ1	1.82	0.43
1:B:238:SER:HB2	1:B:311:TYR:CZ	2.53	0.43
1:A:10:ARG:NH2	1:A:314:GLU:OE1	2.50	0.43
1:A:6:ASN:HD21	3:A:335:APR:H2	1.84	0.43
1:A:18:THR:HG21	1:A:319:GLN:HA	2.01	0.43
1:A:332:ASP:O	1:A:334:ALA:N	2.51	0.43
1:B:180:ALA:HA	1:D:185:VAL:HG22	2.01	0.43
1:B:217:VAL:HG23	1:B:218:ILE:HG13	2.00	0.43
1:B:236:ASN:HD21	1:B:312:ASP:CG	2.22	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:153:CYS:SG	1:D:290:SER:HB2	2.59	0.42
1:D:18:THR:HG21	1:D:319:GLN:HA	2.01	0.42
1:B:213:ALA:HA	1:B:216:LYS:HE2	2.02	0.42
1:A:238:SER:HB2	1:A:311:TYR:CZ	2.55	0.42
1:B:18:THR:HG21	1:B:319:GLN:HA	2.02	0.42
1:B:236:ASN:O	1:B:237:VAL:HB	2.19	0.42
1:D:20:LEU:CD2	1:D:66:VAL:HG11	2.49	0.42
1:B:60:GLU:HB2	1:B:65:VAL:CG2	2.50	0.42
1:C:60:GLU:HB2	1:C:65:VAL:CG2	2.49	0.42
1:C:79:PRO:HD3	1:C:99:PHE:CZ	2.54	0.42
1:D:305:VAL:HG22	1:D:306:LYS:N	2.34	0.42
1:C:18:THR:HG21	1:C:319:GLN:HA	2.01	0.42
1:D:271:LEU:HD12	1:D:290:SER:O	2.20	0.42
1:C:6:ASN:HD21	3:C:338:APR:H2	1.85	0.42
1:C:174:THR:HG23	1:C:229:ALA:CB	2.50	0.42
1:A:79:PRO:HD3	1:A:99:PHE:CZ	2.55	0.42
1:A:167:ILE:HG23	1:A:244:VAL:CG2	2.50	0.42
1:B:174:THR:HG23	1:B:229:ALA:CB	2.49	0.42
1:C:238:SER:HB2	1:C:311:TYR:CZ	2.54	0.42
1:D:56:GLU:OE1	1:D:58:LYS:HE3	2.20	0.42
1:A:171:LEU:HD23	1:A:226:THR:CG2	2.50	0.41
1:B:237:VAL:HG11	1:B:280:SER:O	2.20	0.41
1:A:174:THR:HG23	1:A:229:ALA:CB	2.50	0.41
1:D:13:ARG:NH2	4:D:508:HOH:O	2.53	0.41
1:A:20:LEU:CD2	1:A:66:VAL:HG11	2.50	0.41
1:A:55:GLY:HA3	1:A:67:ASP:OD2	2.19	0.41
1:A:305:VAL:HG22	1:A:306:LYS:N	2.35	0.41
1:B:13:ARG:NH2	4:B:510:HOH:O	2.53	0.41
1:C:332:ASP:O	1:C:334:ALA:N	2.54	0.41
1:D:332:ASP:O	1:D:334:ALA:N	2.53	0.41
1:B:79:PRO:HD3	1:B:99:PHE:CZ	2.56	0.41
1:A:265:GLY:HA3	1:A:266:PRO:HD3	1.88	0.41
1:B:20:LEU:HD21	1:B:66:VAL:HG11	2.02	0.41
1:C:236:ASN:HD21	1:C:312:ASP:CG	2.23	0.41
1:D:213:ALA:HA	1:D:216:LYS:HE2	2.02	0.41
1:D:236:ASN:O	1:D:237:VAL:HB	2.21	0.41
1:C:213:ALA:HA	1:C:216:LYS:HE2	2.03	0.41
1:B:167:ILE:HG23	1:B:244:VAL:CG2	2.51	0.40
1:C:83:PRO:HB2	1:C:86:LYS:CD	2.42	0.40
1:B:300:LEU:HD23	1:C:169:GLU:HB2	2.04	0.40
1:D:238:SER:HB2	1:D:311:TYR:CZ	2.56	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	331/333 (99%)	307 (93%)	19 (6%)	5 (2%)	10	42
1	B	331/333 (99%)	305 (92%)	22 (7%)	4 (1%)	13	48
1	C	331/333 (99%)	305 (92%)	21 (6%)	5 (2%)	10	42
1	D	331/333 (99%)	307 (93%)	20 (6%)	4 (1%)	13	48
All	All	1324/1332 (99%)	1224 (92%)	82 (6%)	18 (1%)	11	43

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	55	GLY
1	B	55	GLY
1	C	55	GLY
1	A	237	VAL
1	A	333	SER
1	B	237	VAL
1	B	333	SER
1	C	237	VAL
1	D	55	GLY
1	D	237	VAL
1	C	313	ASN
1	C	333	SER
1	D	333	SER
1	D	166	GLU
1	A	166	GLU
1	B	166	GLU
1	C	166	GLU
1	A	198	GLY

5.3.2 Protein sidechains ⓘ

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	272/272 (100%)	255 (94%)	17 (6%)	18	51
1	B	272/272 (100%)	255 (94%)	17 (6%)	18	51
1	C	272/272 (100%)	256 (94%)	16 (6%)	19	54
1	D	272/272 (100%)	255 (94%)	17 (6%)	18	51
All	All	1088/1088 (100%)	1021 (94%)	67 (6%)	18	52

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	PHE
1	A	38	GLU
1	A	54	LYS
1	A	60	GLU
1	A	69	LYS
1	A	72	THR
1	A	103	GLU
1	A	110	LYS
1	A	135	GLU
1	A	139	LYS
1	A	154	LEU
1	A	191	LYS
1	A	201	GLN
1	A	231	ARG
1	A	298	ILE
1	A	315	PHE
1	A	319	GLN
1	B	8	PHE
1	B	20	LEU
1	B	38	GLU
1	B	54	LYS
1	B	60	GLU
1	B	69	LYS
1	B	72	THR

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Mol	Chain	Res	Type
1	B	94	GLU
1	B	103	GLU
1	B	110	LYS
1	B	139	LYS
1	B	154	LEU
1	B	191	LYS
1	B	201	GLN
1	B	298	ILE
1	B	315	PHE
1	B	319	GLN
1	C	8	PHE
1	C	20	LEU
1	C	38	GLU
1	C	54	LYS
1	C	60	GLU
1	C	69	LYS
1	C	72	THR
1	C	103	GLU
1	C	110	LYS
1	C	139	LYS
1	C	154	LEU
1	C	191	LYS
1	C	201	GLN
1	C	298	ILE
1	C	315	PHE
1	C	319	GLN
1	D	8	PHE
1	D	20	LEU
1	D	38	GLU
1	D	54	LYS
1	D	60	GLU
1	D	69	LYS
1	D	72	THR
1	D	103	GLU
1	D	110	LYS
1	D	135	GLU
1	D	139	LYS
1	D	154	LEU
1	D	191	LYS
1	D	201	GLN
1	D	298	ILE
1	D	315	PHE

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Mol	Chain	Res	Type
1	D	319	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	ASN
1	A	146	ASN
1	B	6	ASN
1	B	81	ASN
1	B	146	ASN
1	B	201	GLN
1	B	313	ASN
1	C	6	ASN
1	C	146	ASN
1	D	6	ASN
1	D	146	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](#)

12 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	APR	D	337	-	35,39,39	1.63	4 (11%)	42,60,60	1.11	3 (7%)
2	SO4	A	501	-	4,4,4	0.79	0	6,6,6	0.59	0
2	SO4	D	507	-	4,4,4	0.84	0	6,6,6	0.46	0
3	APR	B	336	-	35,39,39	1.65	5 (14%)	42,60,60	1.07	3 (7%)
2	SO4	C	508	-	4,4,4	0.93	0	6,6,6	0.47	0
2	SO4	D	503	-	4,4,4	0.89	0	6,6,6	0.48	0
2	SO4	B	502	-	4,4,4	0.82	0	6,6,6	0.58	0
3	APR	A	335	-	35,39,39	1.74	4 (11%)	42,60,60	1.04	2 (4%)
3	APR	C	338	-	35,39,39	1.71	4 (11%)	42,60,60	1.05	3 (7%)
2	SO4	C	504	-	4,4,4	0.80	0	6,6,6	0.50	0
2	SO4	B	506	-	4,4,4	0.92	0	6,6,6	0.47	0
2	SO4	A	505	-	4,4,4	0.88	0	6,6,6	0.42	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	APR	B	336	-	-	6/18/54/54	0/4/4/4
3	APR	D	337	-	-	7/18/54/54	0/4/4/4
3	APR	A	335	-	-	7/18/54/54	0/4/4/4
3	APR	C	338	-	-	8/18/54/54	0/4/4/4

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	335	APR	PB-O3A	8.26	1.68	1.59
3	C	338	APR	PB-O3A	7.97	1.68	1.59
3	D	337	APR	PB-O3A	7.31	1.67	1.59
3	B	336	APR	PB-O3A	7.05	1.67	1.59
3	B	336	APR	C2-N1	3.03	1.39	1.33
3	A	335	APR	O1D-C1D	2.74	1.48	1.39
3	D	337	APR	O1D-C1D	2.72	1.48	1.39
3	D	337	APR	C2-N1	2.65	1.38	1.33
3	A	335	APR	C2-N1	2.61	1.38	1.33
3	C	338	APR	O4D-C1D	-2.56	1.40	1.43
3	C	338	APR	C2-N1	2.55	1.38	1.33
3	B	336	APR	O1D-C1D	2.44	1.47	1.39
3	C	338	APR	O1D-C1D	2.42	1.47	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	336	APR	PA-O3A	-2.40	1.56	1.59
3	B	336	APR	C4-N3	2.36	1.38	1.35
3	D	337	APR	PA-O3A	-2.24	1.57	1.59
3	A	335	APR	O4D-C1D	-2.17	1.40	1.43

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	336	APR	C1'-N9-C4	-2.64	122.00	126.64
3	C	338	APR	O4'-C1'-N9	2.59	112.19	108.75
3	A	335	APR	O4'-C1'-N9	2.52	112.09	108.75
3	D	337	APR	C1'-N9-C4	-2.34	122.53	126.64
3	B	336	APR	O4'-C1'-N9	2.32	111.82	108.75
3	D	337	APR	C4'-O4'-C1'	2.28	112.01	109.92
3	B	336	APR	C5-C6-N6	2.26	123.75	120.31
3	D	337	APR	C1D-C2D-C3D	2.22	105.03	102.29
3	C	338	APR	C1'-N9-C4	-2.19	122.80	126.64
3	C	338	APR	C5-C6-N6	2.11	123.53	120.31
3	A	335	APR	C1'-N9-C4	-2.10	122.96	126.64

There are no chirality outliers.

All (28) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	335	APR	C5'-O5'-PA-O1A
3	A	335	APR	C5'-O5'-PA-O2A
3	A	335	APR	C5'-O5'-PA-O3A
3	B	336	APR	C5'-O5'-PA-O2A
3	B	336	APR	C5'-O5'-PA-O3A
3	C	338	APR	C5'-O5'-PA-O1A
3	C	338	APR	C5'-O5'-PA-O2A
3	C	338	APR	C5'-O5'-PA-O3A
3	D	337	APR	C5'-O5'-PA-O2A
3	D	337	APR	C5'-O5'-PA-O3A
3	D	337	APR	PB-O3A-PA-O5'
3	A	335	APR	C3'-C4'-C5'-O5'
3	A	335	APR	O4'-C4'-C5'-O5'
3	B	336	APR	O4'-C4'-C5'-O5'
3	D	337	APR	O4'-C4'-C5'-O5'
3	B	336	APR	C3'-C4'-C5'-O5'
3	C	338	APR	O4'-C4'-C5'-O5'
3	D	337	APR	C3'-C4'-C5'-O5'

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Mol	Chain	Res	Type	Atoms
3	C	338	APR	C3'-C4'-C5'-O5'
3	A	335	APR	PB-O3A-PA-O5'
3	B	336	APR	PB-O3A-PA-O5'
3	C	338	APR	PB-O3A-PA-O5'
3	A	335	APR	PB-O3A-PA-O1A
3	B	336	APR	PB-O3A-PA-O1A
3	C	338	APR	PB-O3A-PA-O1A
3	D	337	APR	PB-O3A-PA-O1A
3	C	338	APR	PB-O3A-PA-O2A
3	D	337	APR	PB-O3A-PA-O2A

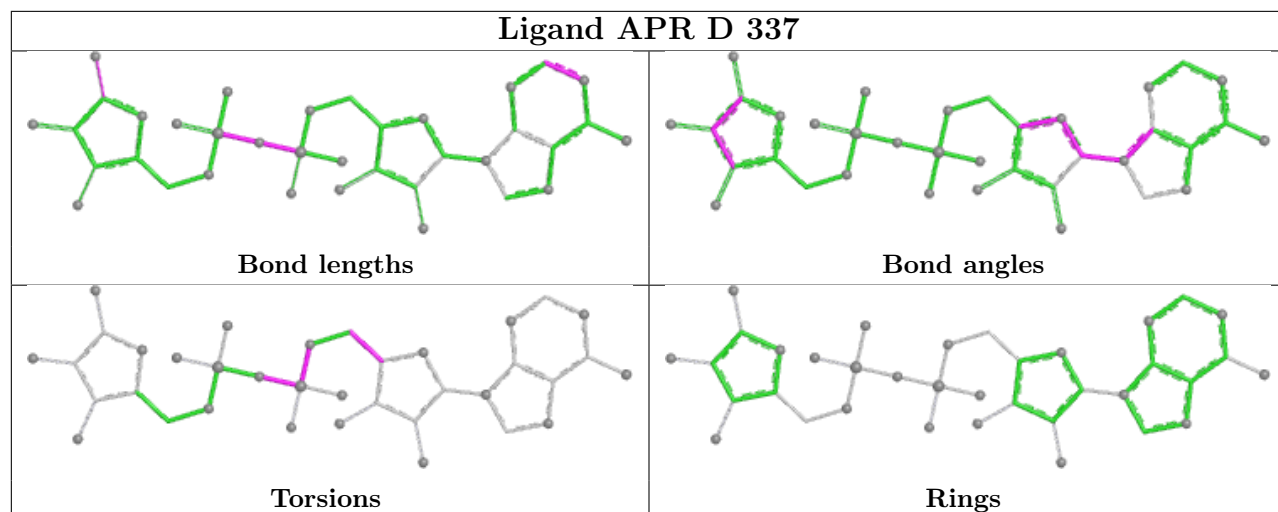
There are no ring outliers.

4 monomers are involved in 11 short contacts:

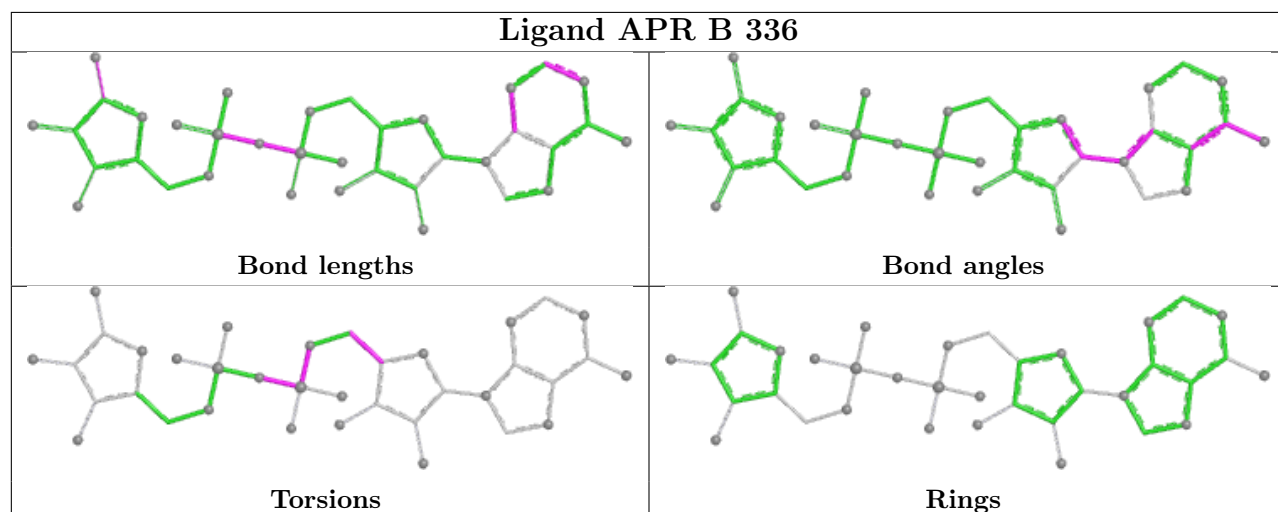
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	337	APR	3	0
3	B	336	APR	1	0
3	A	335	APR	3	0
3	C	338	APR	4	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less 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.

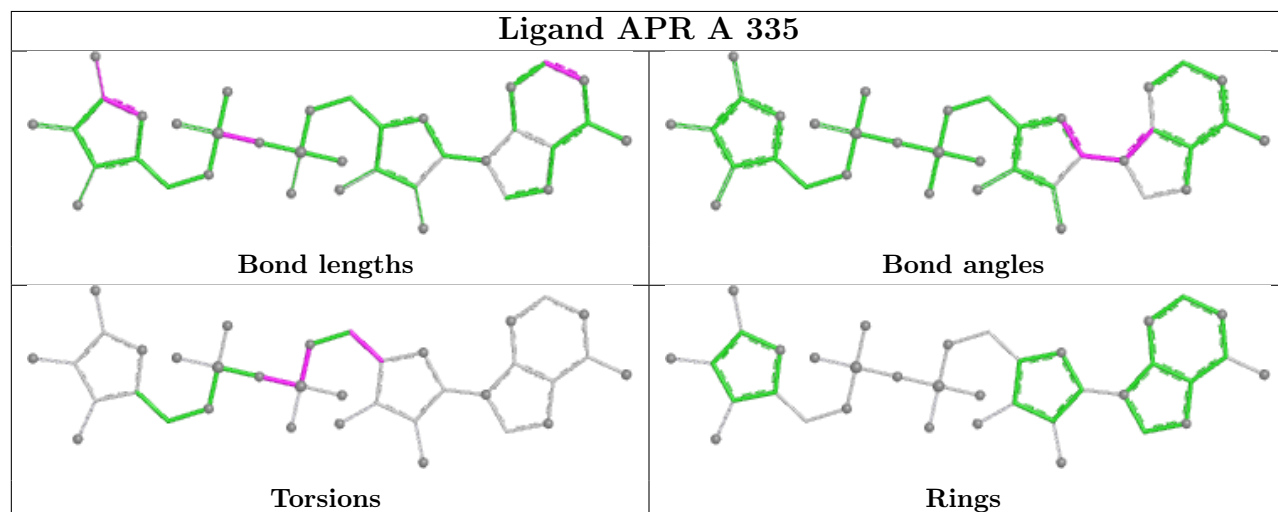
Ligand APR D 337

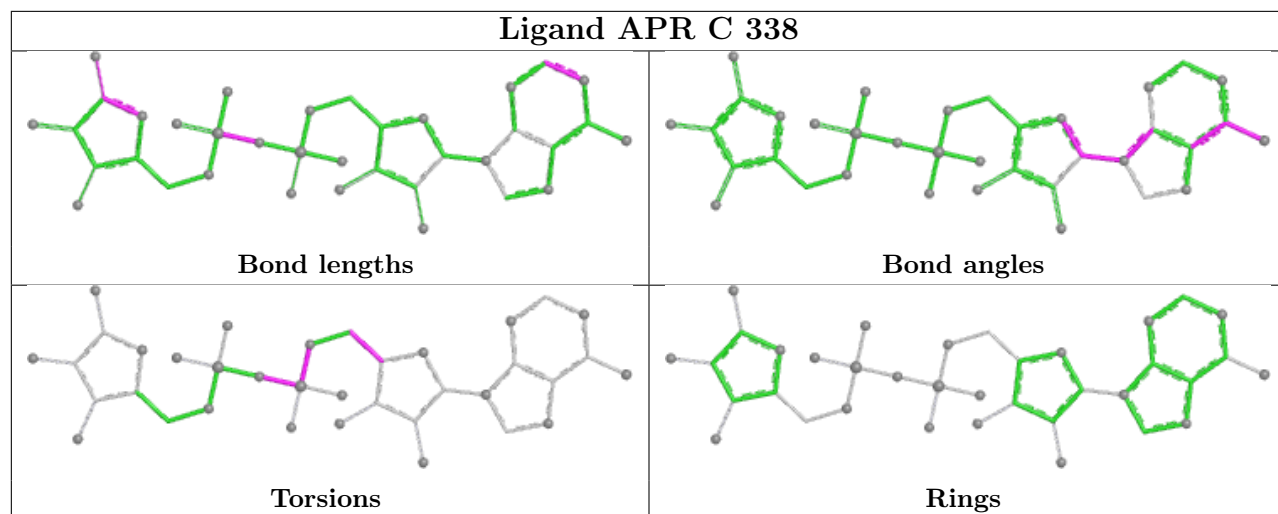


Ligand APR B 336



Ligand APR A 335





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	333/333 (100%)	-0.74	3 (0%) 84 63	8, 23, 46, 64	0
1	B	333/333 (100%)	-0.77	1 (0%) 94 84	7, 23, 42, 61	0
1	C	333/333 (100%)	-0.65	1 (0%) 94 84	13, 31, 56, 72	0
1	D	333/333 (100%)	-0.66	1 (0%) 94 84	8, 28, 58, 71	0
All	All	1332/1332 (100%)	-0.71	6 (0%) 91 75	7, 25, 52, 72	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	334	ALA	5.3
1	C	334	ALA	3.9
1	B	334	ALA	3.1
1	A	334	ALA	2.9
1	A	61	ASP	2.4
1	A	333	SER	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

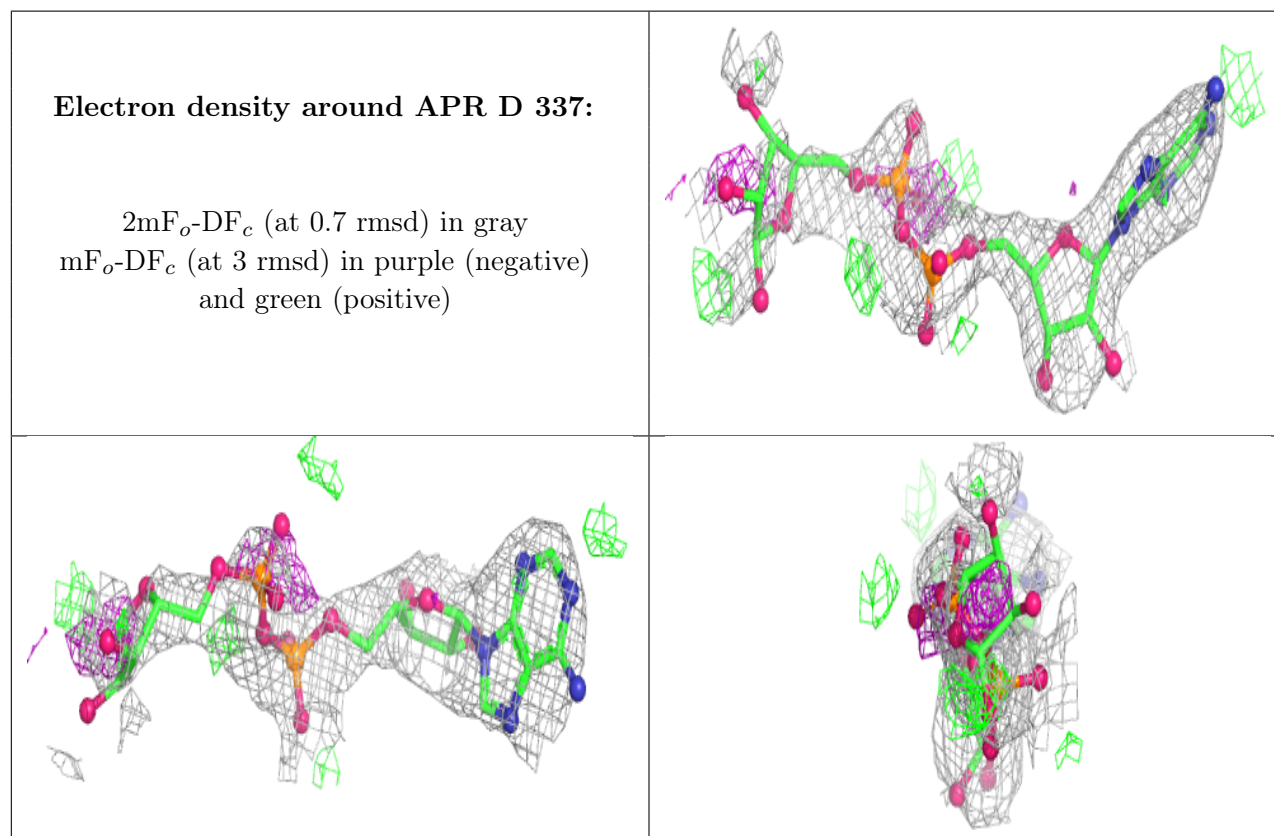
6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

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

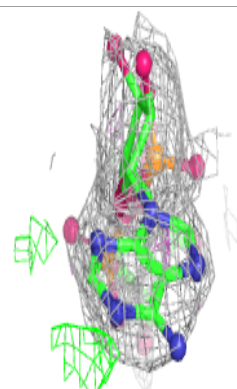
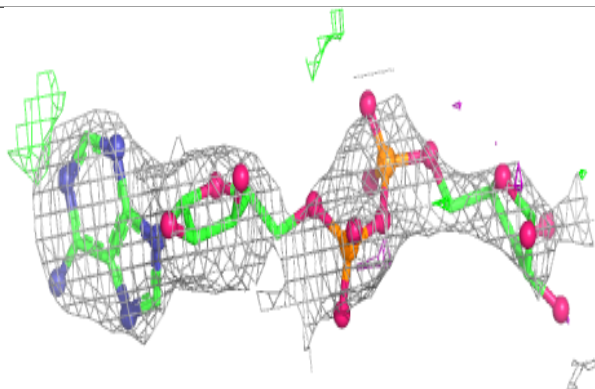
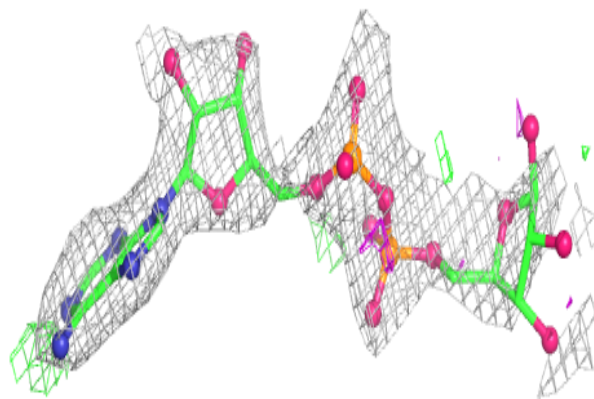
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	APR	D	337	36/36	0.81	0.28	79,94,103,103	0
2	SO4	C	508	5/5	0.85	0.33	103,103,103,103	0
3	APR	C	338	36/36	0.86	0.29	67,87,103,103	0
2	SO4	B	506	5/5	0.88	0.24	102,103,103,103	0
3	APR	A	335	36/36	0.89	0.23	52,76,103,103	0
3	APR	B	336	36/36	0.89	0.23	45,74,103,103	0
2	SO4	C	504	5/5	0.90	0.23	89,89,90,90	0
2	SO4	B	502	5/5	0.91	0.17	81,82,82,82	0
2	SO4	A	505	5/5	0.91	0.18	100,100,100,100	0
2	SO4	D	503	5/5	0.92	0.19	72,73,73,74	0
2	SO4	D	507	5/5	0.95	0.20	101,102,102,102	0
2	SO4	A	501	5/5	0.96	0.14	74,75,75,75	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.

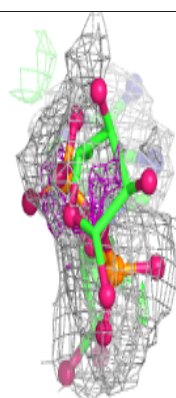
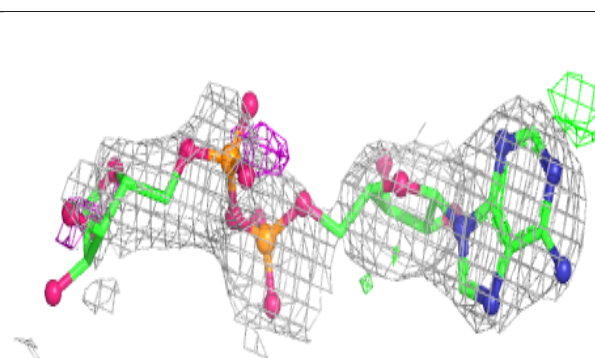
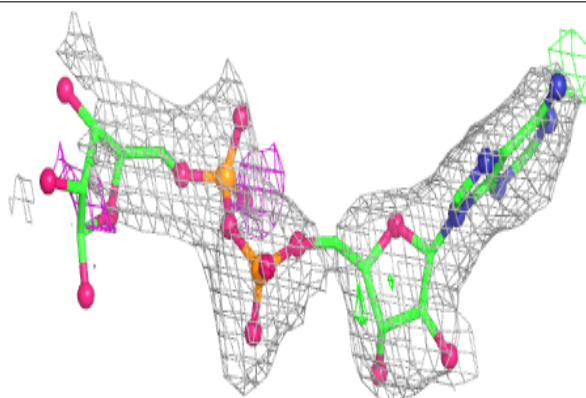


Electron density around APR C 338:

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

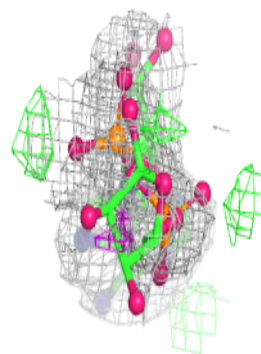
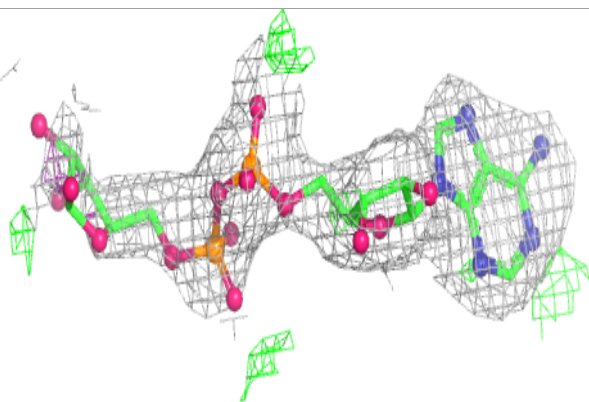
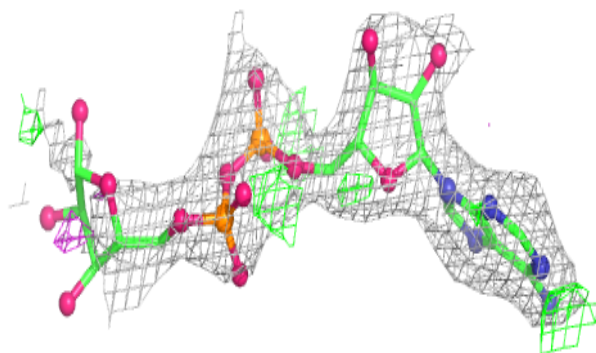
**Electron density around APR A 335:**

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



Electron density around APR B 336:

$2mF_o-DF_c$ (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.