Full wwPDB X-ray Structure Validation Report (i)

Jun 10, 2024 – 07:49 PM EDT

PDB ID	:	8SWC
Title	:	RNase H complex with ASO (OOO) and RNA
Authors	:	Cho, Y.-J.; Iwamoto, N.
Deposited on	:	2023-05-18
Resolution	:	2.68 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

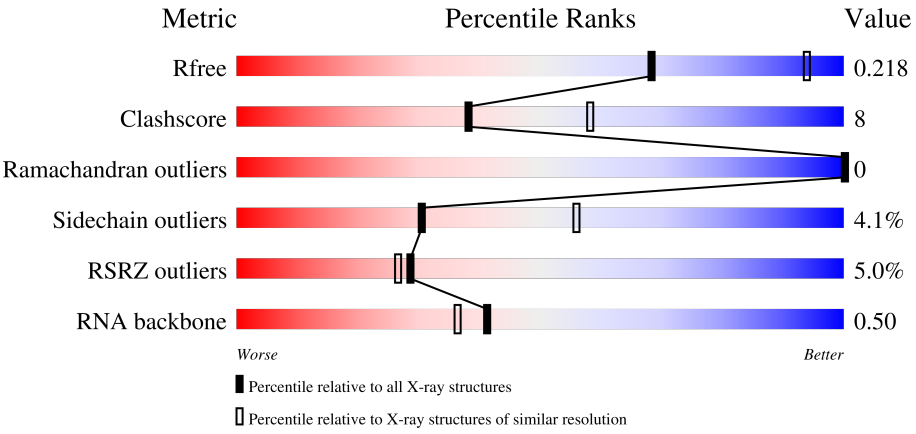
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.68 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	3863 (2.70-2.66)
Clashscore	141614	4210 (2.70-2.66)
Ramachandran outliers	138981	4141 (2.70-2.66)
Sidechain outliers	138945	4141 (2.70-2.66)
RSRZ outliers	127900	3780 (2.70-2.66)
RNA backbone	3102	1007 (2.98-2.38)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	294	3%, 38%, 12%, ·, 49%
2	B	20	75%, 25%
3	C	20	40%, 35%, 20%, 5%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 2052 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 Ribonuclease H1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	149	Total	C	N	O	S	0	0	0
			1168	727	220	215	6			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	HIS	-	expression tag	UNP O60930
A	-6	HIS	-	expression tag	UNP O60930
A	-5	HIS	-	expression tag	UNP O60930
A	-4	HIS	-	expression tag	UNP O60930
A	-3	HIS	-	expression tag	UNP O60930
A	-2	HIS	-	expression tag	UNP O60930
A	-1	HIS	-	expression tag	UNP O60930
A	0	HIS	-	expression tag	UNP O60930
A	210	ASN	ASP	engineered mutation	UNP O60930

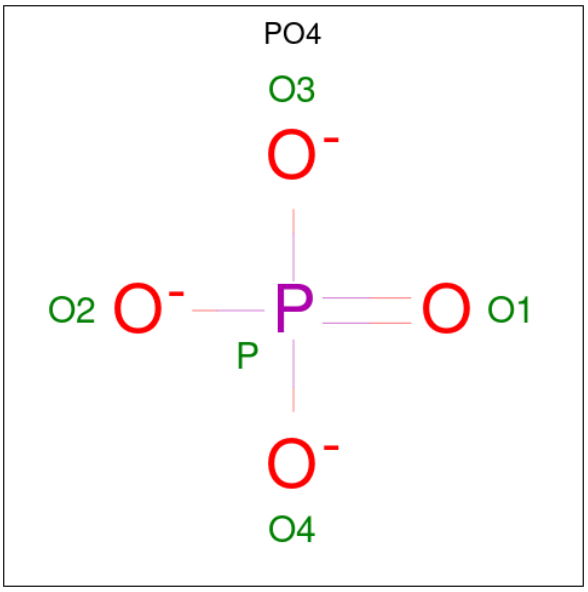
- Molecule 2 is a RNA chain called RNA (5'-R(*UP*GP*GP*CP*GP*AP*GP*UP*GP*GP*GP*UP*GP*AP*GP*UP*GP*AP*GP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	20	Total	C	N	O	P	0	0	0
			439	195	86	139	19			

- Molecule 3 is a DNA chain called PO ASO.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	20	Total	C	N	O	P	0	0	0
			422	206	67	130	19			

- Molecule 4 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).

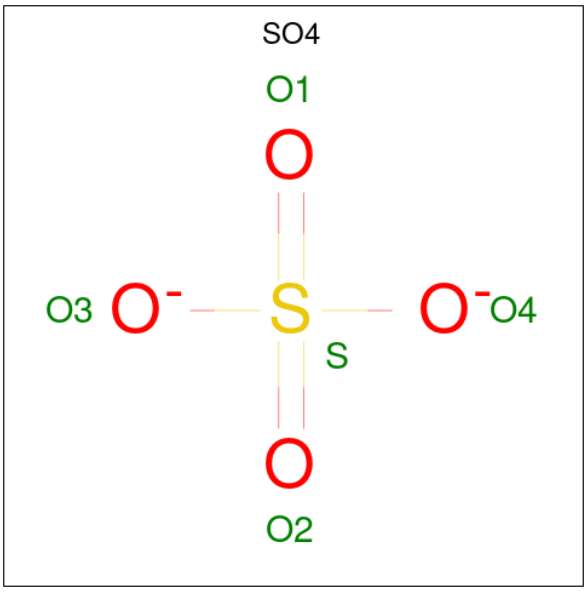


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total O P 5 4 1	0	0
4	B	1	Total O P 5 4 1	0	0

- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total Ca 1 1	0	0
5	B	2	Total Ca 2 2	0	0

- Molecule 6 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	B	1	Total O S 5 4 1	0	0

- Molecule 7 is water.

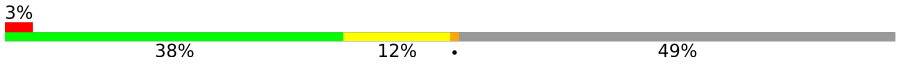
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	2	Total O 2 2	0	0
7	B	2	Total O 2 2	0	0
7	C	1	Total O 1 1	0	0

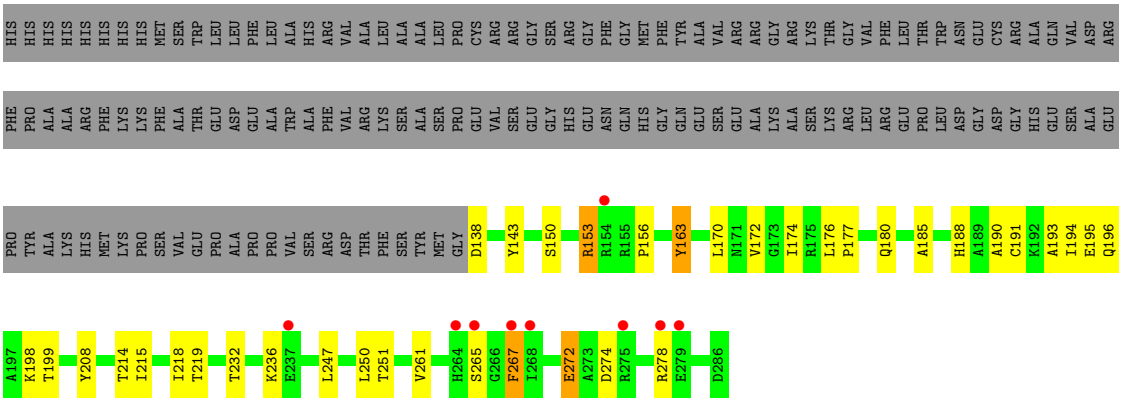
3 Residue-property plots (i)

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

• Molecule 1: Ribonuclease H1

Chain A:





• Molecule 2: RNA (5'-R(*UP*GP*GP*CP*GP*AP*GP*UP*GP*GP*GP*UP*GP*AP*GP*UP *GP*AP*GP*G)-3')

Chain B:

• Molecule 3: PO ASO

Chain C:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α, β, γ	100.62Å 100.62Å 82.09Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	38.48 – 2.68 42.89 – 2.67	Depositor EDS
% Data completeness (in resolution range)	99.6 (38.48-2.68) 99.0 (42.89-2.67)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.89 (at 2.69Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.192 , 0.220 0.189 , 0.218	Depositor DCC
R_{free} test set	724 reflections (5.23%)	wwPDB-VP
Wilson B-factor (Å^2)	74.9	Xtriage
Anisotropy	0.353	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 54.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	0.025 for -h,-k,l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	2052	wwPDB-VP
Average B, all atoms (Å^2)	83.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.42% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: OMG, SO4, CA, OMC, T39, C5L, PO4, A2M

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	0.25	0/1193	0.49	0/1611
2	B	0.18	0/493	0.74	0/771
3	C	0.54	0/218	0.91	0/334
All	All	0.28	0/1904	0.63	0/2716

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

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	1168	0	1141	21	0
2	B	439	0	218	0	0
3	C	422	0	259	5	0
4	A	5	0	0	0	0
4	B	5	0	0	0	0
5	A	1	0	0	0	0
5	B	2	0	0	0	0
6	B	5	0	0	0	0
7	A	2	0	0	0	0
7	B	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	C	1	0	0	0	0
All	All	2052	0	1618	26	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 8.

All (26) 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:150:SER:HB3	1:A:153:ARG:HB2	1.81	0.61
3:C:23:OMC:H2'	3:C:24:A2M:H8	1.85	0.58
1:A:143:TYR:HH	1:A:267:PHE:HE1	1.51	0.58
1:A:208:TYR:HD2	1:A:261:VAL:HG22	1.72	0.55
1:A:198:LYS:NZ	1:A:250:LEU:O	2.40	0.55
1:A:138:ASP:OD2	1:A:138:ASP:N	2.38	0.55
1:A:194:ILE:HD12	1:A:194:ILE:H	1.73	0.53
1:A:215:ILE:O	1:A:219:THR:OG1	2.19	0.52
1:A:188:HIS:HA	1:A:191:CYS:HB2	1.91	0.52
1:A:274:ASP:O	1:A:278:ARG:HG2	2.12	0.49
1:A:163:TYR:HD2	1:A:272:GLU:HB3	1.79	0.48
1:A:176:LEU:HD22	1:A:185:ALA:HA	1.96	0.48
1:A:156:PRO:HG2	1:A:180:GLN:HB2	1.96	0.48
1:A:163:TYR:CD2	1:A:272:GLU:HB3	2.48	0.47
1:A:172:VAL:HB	1:A:174:ILE:HD11	1.97	0.47
3:C:23:OMC:H1'	3:C:23:OMC:HM23	1.73	0.46
1:A:176:LEU:HD12	1:A:177:PRO:HD2	1.97	0.46
1:A:247:LEU:O	1:A:251:THR:HG23	2.16	0.46
3:C:3:T39:H1'	3:C:3:T39:HCA2	1.47	0.45
3:C:14:DC:H2"	3:C:15:DT:H5"	2.00	0.44
3:C:23:OMC:H2'	3:C:24:A2M:C8	2.47	0.44
1:A:214:THR:O	1:A:218:ILE:HG22	2.18	0.44
1:A:190:ALA:O	1:A:193:ALA:N	2.49	0.44
1:A:170:LEU:HD13	1:A:196:GLN:HB3	1.99	0.43
1:A:195:GLU:O	1:A:199:THR:HG23	2.19	0.43
1:A:232:THR:HG23	1:A:236:LYS:O	2.21	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	147/294 (50%)	139 (95%)	8 (5%)	0	100 100

There are no Ramachandran outliers to report.

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	121/239 (51%)	116 (96%)	5 (4%)	30 56

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	153	ARG
1	A	163	TYR
1	A	265	SER
1	A	267	PHE
1	A	272	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	19/20 (95%)	5 (26%)	0

All (5) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	7	G
2	B	10	G
2	B	13	G
2	B	16	U
2	B	20	G

There are no RNA pucker outliers to report.

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

10 non-standard protein/DNA/RNA residues 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	A2M	C	5	2,3	18,25,26	2.01	5 (27%)	18,36,39	4.14	6 (33%)
3	OMC	C	16	2,3	19,22,23	1.50	4 (21%)	26,31,34	1.08	1 (3%)
3	OMC	C	23	2,3	19,22,23	1.41	4 (21%)	26,31,34	1.12	2 (7%)
3	A2M	C	24	2,3	18,25,26	2.03	5 (27%)	18,36,39	4.54	4 (22%)
3	C5L	C	2	2,3	22,26,27	0.98	3 (13%)	30,36,39	1.02	3 (10%)
3	OMC	C	22	2,3	19,22,23	1.40	3 (15%)	26,31,34	1.14	2 (7%)
3	OMC	C	1	2,3	19,19,23	1.44	3 (15%)	26,27,34	1.10	1 (3%)
3	T39	C	3	2,3	23,26,27	1.61	4 (17%)	32,36,39	2.18	4 (12%)
3	C5L	C	4	2,3	22,26,27	1.02	3 (13%)	30,36,39	0.98	2 (6%)
3	OMG	C	21	2,3	18,26,27	2.93	7 (38%)	19,38,41	1.42	3 (15%)

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	A2M	C	5	2,3	-	0/5/27/28	0/3/3/3
3	OMC	C	16	2,3	-	2/9/27/28	0/2/2/2
3	OMC	C	23	2,3	-	3/9/27/28	0/2/2/2
3	A2M	C	24	2,3	-	0/5/27/28	0/3/3/3
3	C5L	C	2	2,3	-	4/12/30/31	0/2/2/2
3	OMC	C	22	2,3	-	0/9/27/28	0/2/2/2
3	OMC	C	1	2,3	-	0/8/24/28	0/2/2/2
3	T39	C	3	2,3	-	3/12/30/31	0/2/2/2
3	C5L	C	4	2,3	-	2/12/30/31	0/2/2/2
3	OMG	C	21	2,3	-	2/5/27/28	0/3/3/3

All (41) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	21	OMG	O6-C6	8.39	1.40	1.23
3	C	5	A2M	O5'-C5'	-5.62	1.31	1.44
3	C	24	A2M	O5'-C5'	-5.59	1.31	1.44
3	C	21	OMG	C2-N2	4.69	1.45	1.34
3	C	21	OMG	O4'-C1'	4.50	1.47	1.41
3	C	3	T39	C2-N1	4.04	1.44	1.38
3	C	21	OMG	C8-N7	3.74	1.41	1.35
3	C	3	T39	C2-N3	3.60	1.44	1.38
3	C	3	T39	C4-C5	3.51	1.50	1.44
3	C	3	T39	O4-C4	-3.17	1.17	1.23
3	C	24	A2M	C5'-C4'	3.16	1.61	1.51
3	C	5	A2M	C5'-C4'	3.07	1.61	1.51
3	C	1	OMC	C2-N1	3.03	1.46	1.40
3	C	21	OMG	C5-C4	2.83	1.50	1.43
3	C	16	OMC	C2-N1	2.82	1.46	1.40
3	C	24	A2M	C2-N3	2.82	1.36	1.32
3	C	16	OMC	C2-N3	2.79	1.42	1.36
3	C	5	A2M	C2-N3	2.76	1.36	1.32
3	C	23	OMC	C2-N3	2.71	1.41	1.36
3	C	1	OMC	C2-N3	2.71	1.41	1.36
3	C	22	OMC	C2-N1	2.70	1.45	1.40
3	C	22	OMC	C2-N3	2.68	1.41	1.36
3	C	4	C5L	C2-N1	2.63	1.45	1.40
3	C	23	OMC	C2-N1	2.62	1.45	1.40
3	C	5	A2M	C2-N1	2.61	1.38	1.33
3	C	24	A2M	C2-N1	2.56	1.38	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	16	OMC	C4-N4	2.49	1.39	1.33
3	C	22	OMC	C4-N4	2.46	1.39	1.33
3	C	21	OMG	C3'-C2'	-2.44	1.47	1.52
3	C	1	OMC	C4-N4	2.43	1.39	1.33
3	C	23	OMC	C4-N4	2.36	1.39	1.33
3	C	2	C5L	C6-C5	2.26	1.38	1.34
3	C	2	C5L	C2-N1	2.26	1.44	1.40
3	C	2	C5L	C2-N3	2.17	1.40	1.36
3	C	4	C5L	C6-C5	2.16	1.38	1.34
3	C	5	A2M	C4-N3	2.15	1.38	1.35
3	C	4	C5L	C2-N3	2.09	1.40	1.36
3	C	21	OMG	C4-N3	2.09	1.42	1.37
3	C	16	OMC	O5'-C5'	-2.06	1.39	1.44
3	C	24	A2M	C4-N3	2.04	1.38	1.35
3	C	23	OMC	C5'-C4'	2.03	1.57	1.51

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	C	24	A2M	O2'-C2'-C1'	-18.21	72.98	109.09
3	C	5	A2M	O2'-C2'-C1'	-15.75	77.87	109.09
3	C	3	T39	C7-C5-C4	-7.99	109.98	118.77
3	C	3	T39	C5-C4-N3	-5.45	110.66	115.31
3	C	3	T39	C6-C5-C4	5.23	122.41	118.03
3	C	5	A2M	O4'-C1'-C2'	-4.20	99.30	106.59
3	C	3	T39	C7-C5-C6	3.57	127.62	122.85
3	C	24	A2M	O4'-C1'-C2'	-3.50	100.52	106.59
3	C	5	A2M	O4'-C4'-C5'	3.46	120.76	109.37
3	C	24	A2M	O4'-C4'-C5'	3.09	119.52	109.37
3	C	21	OMG	C2-N1-C6	-3.04	119.50	125.10
3	C	5	A2M	C3'-C2'-C1'	-3.03	97.18	102.89
3	C	21	OMG	C5-C6-N1	2.96	119.17	113.95
3	C	5	A2M	C5-C6-N6	2.92	124.79	120.35
3	C	24	A2M	C5-C6-N6	2.92	124.79	120.35
3	C	2	C5L	C7-C5-C6	-2.62	119.35	122.85
3	C	16	OMC	C6-N1-C2	-2.55	116.07	120.49
3	C	21	OMG	C8-N7-C5	2.54	107.83	102.99
3	C	1	OMC	C6-N1-C2	-2.54	116.09	120.49
3	C	4	C5L	C7-C5-C6	-2.46	119.56	122.85
3	C	22	OMC	C6-N1-C2	-2.40	116.33	120.49
3	C	22	OMC	O4'-C1'-N1	2.33	113.68	108.36
3	C	23	OMC	C6-N1-C2	-2.32	116.47	120.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	C	23	OMC	O4'-C1'-N1	2.31	113.65	108.36
3	C	4	C5L	C5-C4-N3	2.13	123.97	121.67
3	C	2	C5L	C5-C6-N1	-2.05	121.23	123.34
3	C	5	A2M	C2'-C3'-C4'	-2.03	97.58	101.99
3	C	2	C5L	C5-C4-N3	2.00	123.84	121.67

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	2	C5L	C1'-C2'-O2'-CA'
3	C	3	T39	C1'-C2'-O2'-CA'
3	C	16	OMC	C3'-C4'-C5'-O5'
3	C	21	OMG	C3'-C4'-C5'-O5'
3	C	23	OMC	C1'-C2'-O2'-CM2
3	C	16	OMC	O4'-C4'-C5'-O5'
3	C	23	OMC	C3'-C4'-C5'-O5'
3	C	23	OMC	O4'-C4'-C5'-O5'
3	C	2	C5L	O2'-CA'-CB'-OC'
3	C	21	OMG	O4'-C4'-C5'-O5'
3	C	4	C5L	CB'-CA'-O2'-C2'
3	C	2	C5L	CA'-CB'-OC'-CD'
3	C	3	T39	CB'-CA'-O2'-C2'
3	C	2	C5L	C3'-C2'-O2'-CA'
3	C	4	C5L	C1'-C2'-O2'-CA'
3	C	3	T39	O2'-CA'-CB'-OC'

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	23	OMC	3	0
3	C	24	A2M	2	0
3	C	3	T39	1	0

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 6 ligands modelled in this entry, 3 are monoatomic - leaving 3 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	PO4	A	301	-	4,4,4	0.92	0	6,6,6	0.42	0
6	SO4	B	101	-	4,4,4	0.14	0	6,6,6	0.05	0
4	PO4	B	102	-	4,4,4	0.92	0	6,6,6	0.43	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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^{th} 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	149/294 (50%)	0.40	9 (6%) 21 20	54, 76, 116, 133	0
2	B	20/20 (100%)	-0.65	0 100 100	61, 81, 126, 127	0
3	C	10/20 (50%)	-0.22	0 100 100	56, 62, 76, 84	0
All	All	179/334 (53%)	0.25	9 (5%) 28 26	54, 75, 120, 133	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	264	HIS	3.4
1	A	279	GLU	3.2
1	A	154	ARG	2.8
1	A	278	ARG	2.6
1	A	268	ILE	2.3
1	A	267	PHE	2.2
1	A	265	SER	2.2
1	A	237	GLU	2.1
1	A	275	ARG	2.0

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

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^{th} 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	OMG	C	21	24/25	0.91	0.15	61,92,105,110	0
3	A2M	C	5	23/24	0.92	0.13	76,97,112,116	0
3	C5L	C	4	25/26	0.94	0.15	85,102,117,127	0
3	C5L	C	2	25/26	0.94	0.16	93,119,126,127	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	T39	C	3	25/26	0.94	0.18	86,108,132,140	0
3	OMC	C	23	21/22	0.94	0.10	89,106,119,123	0
3	OMC	C	22	21/22	0.95	0.13	73,99,108,118	0
3	OMC	C	16	21/22	0.95	0.14	61,83,95,109	0
3	A2M	C	24	23/24	0.95	0.13	93,114,124,133	0
3	OMC	C	1	18/22	0.96	0.12	89,111,123,124	0

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
5	CA	B	103	1/1	0.43	0.13	107,107,107,107	0
6	SO4	B	101	5/5	0.61	0.39	130,142,156,163	0
5	CA	B	104	1/1	0.85	0.51	107,107,107,107	0
5	CA	A	302	1/1	0.87	0.14	99,99,99,99	0
4	PO4	A	301	5/5	0.88	0.28	78,97,117,145	0
4	PO4	B	102	5/5	0.94	0.10	94,97,110,138	0

6.5 Other polymers (i)

There are no such residues in this entry.