



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 26, 2026 – 02:40 AM EDT

PDB ID : 8SWB / pdb_00008swb
Title : RNase H complex with streopure ASO and RNA
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Deposited on : 2023-05-18
Resolution : 2.00 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	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

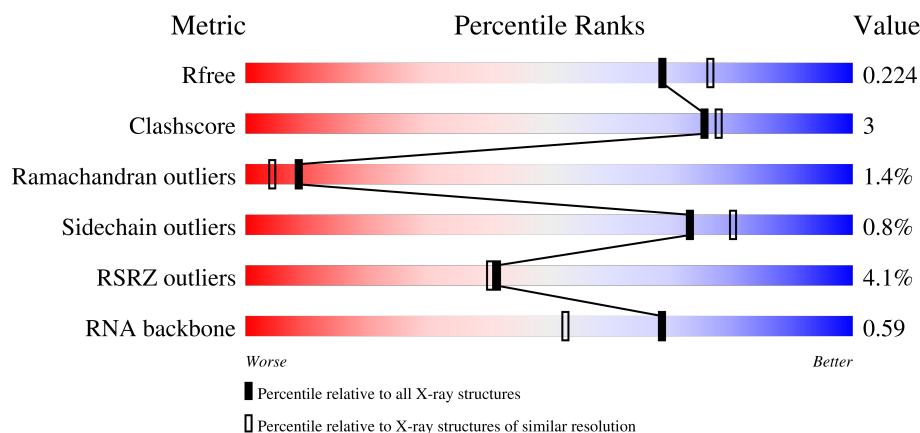
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)
RNA backbone	3983	1000 (2.36-1.64)

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	149	5% 93% 7%
2	B	20	85% 10% 5%
3	D	20	5% 75% 20%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 2226 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	8	0
			1237	767	239	225	6			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	210	ASN	ASP	conflict	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 DNA (5'-R(*(OMC)P*(N7X)P*(T39)P*(C5L)P*(A2M))-D(P*(SC)P*(PST)P*(SC)P*AP*(SC)P*(SC)P*(RC)P*(AS)P*(SC)P*(PST))-R(P*(6OO)P*(RFJ)P*(6OO)P*(6OO)P*(6NW))-3').

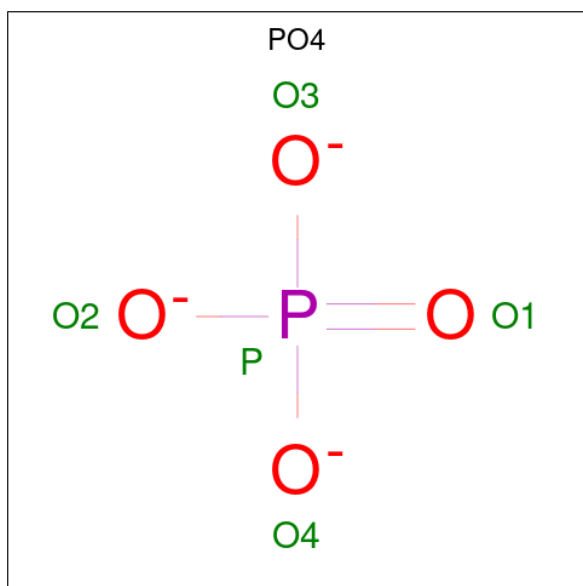
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	20	Total	C	N	O	P	S	0	0
			422	206	67	115	19	15		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



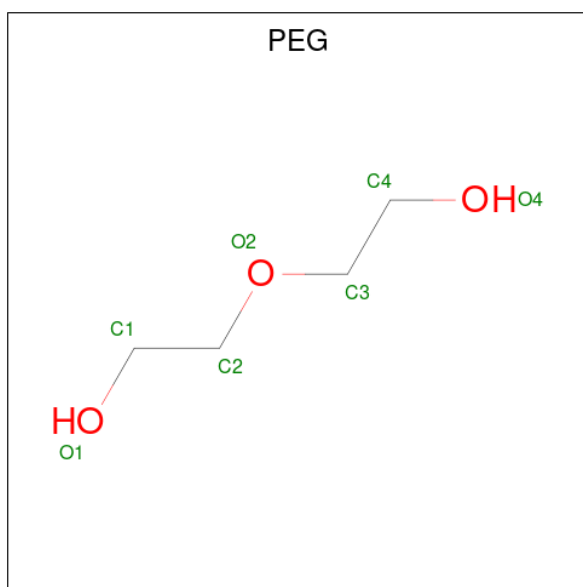
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



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

- Molecule 6 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



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

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	53	Total	O	0	0
			53	53		
7	B	29	Total	O	0	0
			29	29		
7	D	28	Total	O	0	0
			28	28		

3 Residue-property plots

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

- Molecule 1: Ribonuclease H1

Chain A: 

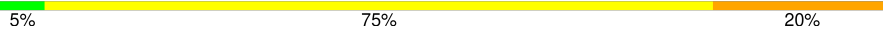


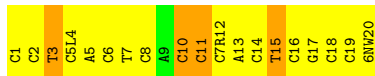
- Molecule 2: RNA (5'-R(*UP*GP*GP*CP*GP*AP*GP*UP*GP*GP*GP*UP*GP*AP*GP*UP*GP*AP*GP*G)-3')

Chain B: 



- Molecule 3: DNA (5'-R(*(OMC)P*(N7X)P*(T39)P*(C5L)P*(A2M))-D(P*(SC)P*(PST)P*(SC)P*AP*(SC)P*(SC)P*(RC)P*(AS)P*(SC)P*(PST))-R(P*(6OO)P*(RFJ)P*(6OO)P*(6OO)P*(6NW))-3')

Chain D: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	46.55Å 52.94Å 106.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.62 – 2.00 42.62 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.7 (42.62-2.00) 99.7 (42.62-2.00)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.35 (at 2.00Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.184 , 0.224 0.184 , 0.224	Depositor DCC
R_{free} test set	916 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	38.2	Xtriage
Anisotropy	0.396	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 50.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	2226	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.26% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: A2M, RFJ, OKQ, OMC, GOL, 6NW, C5L, AS, T39, SC, PEG, N7X, C7R, PST, PO4

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	0.23	0/1264	0.40	0/1706
2	B	0.26	0/493	0.47	0/771
3	D	0.43	0/23	0.70	0/33
All	All	0.24	0/1780	0.42	0/2510

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	1237	0	1201	6	1
2	B	439	0	218	1	0
3	D	422	0	163	3	1
4	A	6	0	8	0	0
5	A	5	0	0	0	0
6	D	7	0	10	1	0
7	A	53	0	0	0	0
7	B	29	0	0	0	0
7	D	28	0	0	0	0
All	All	2226	0	1600	10	1

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 3.

All (10) 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)
3:D:10:SC:HN42	6:D:201:PEG:H11	1.71	0.55
1:A:264:HIS:O	1:A:264:HIS:ND1	2.40	0.54
1:A:275:ARG:O	1:A:279:GLU:HG3	2.09	0.53
2:B:12:U:H2'	2:B:13:G:O4'	2.14	0.47
1:A:156:PRO:HG2	1:A:180:GLN:HB2	1.97	0.47
1:A:264:HIS:O	1:A:264:HIS:CG	2.68	0.46
3:D:3:T39:HCA1	3:D:3:T39:H1'	1.68	0.45
1:A:254:MET:HE3	1:A:254:MET:HB3	1.76	0.43
3:D:10:SC:H2'	3:D:11:SC:H6	2.00	0.43
1:A:188[B]:HIS:HA	1:A:191:CYS:HB2	2.01	0.43

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:195[A]:GLU:OE2	3:D:15:PST:SP[4_555]	1.47	0.73

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	155/149 (104%)	146 (94%)	7 (4%)	2 (1%)	9 5

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	263	GLY
1	A	264	HIS

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	128/120 (107%)	127 (99%)	1 (1%)	73	80

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	179	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	223	GLN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	19/20 (95%)	2 (10%)	0

All (2) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	13	G
2	B	19	G

There are no RNA pucker outliers to report.

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

19 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	SC	D	14	3,2	17,20,21	2.50	5 (29%)	23,28,31	1.74	5 (21%)
3	C5L	D	4	3,2	23,26,27	0.98	2 (8%)	30,36,39	3.76	4 (13%)
3	OKQ	D	16	3,2	19,22,23	2.30	6 (31%)	25,31,34	1.59	5 (20%)
3	T39	D	3	3,2	23,26,27	1.58	4 (17%)	31,36,39	2.10	4 (12%)
3	OMC	D	1	3,2	19,19,23	1.56	3 (15%)	26,27,34	1.13	2 (7%)
3	A2M	D	5	3,2	22,25,26	2.11	6 (27%)	30,36,39	2.61	12 (40%)
3	N7X	D	2	3,2	23,26,27	1.53	6 (26%)	30,36,39	1.76	5 (16%)
3	OKQ	D	18	3,2	19,22,23	2.39	7 (36%)	25,31,34	1.60	6 (24%)
3	6NW	D	20	3,2	22,25,26	1.67	6 (27%)	30,36,39	1.72	4 (13%)
3	RFJ	D	17	3,2	23,26,27	2.88	10 (43%)	32,38,41	2.56	8 (25%)
3	PST	D	7	3,2	18,21,22	2.29	8 (44%)	25,30,33	1.90	4 (16%)
3	C7R	D	12	3,2	17,20,21	0.68	1 (5%)	23,28,31	0.85	0
3	SC	D	11	3,2	17,20,21	2.50	4 (23%)	23,28,31	1.56	7 (30%)
3	SC	D	6	3,2	17,20,21	2.65	5 (29%)	23,28,31	1.40	5 (21%)
3	SC	D	10	3,2	17,20,21	2.19	6 (35%)	23,28,31	1.50	5 (21%)
3	PST	D	15	3,2	18,21,22	2.46	7 (38%)	25,30,33	1.94	4 (16%)
3	SC	D	8	3,2	17,20,21	2.59	5 (29%)	23,28,31	1.36	4 (17%)
3	OKQ	D	19	3,2	19,22,23	2.31	7 (36%)	25,31,34	1.55	3 (12%)
3	AS	D	13	3,2	20,23,24	1.94	6 (30%)	28,33,36	1.23	3 (10%)

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	SC	D	14	3,2	-	0/7/21/22	0/2/2/2
3	C5L	D	4	3,2	-	2/12/30/31	0/2/2/2
3	OKQ	D	16	3,2	-	1/9/27/28	0/2/2/2
3	T39	D	3	3,2	-	3/12/30/31	0/2/2/2
3	OMC	D	1	3,2	-	1/8/24/28	0/2/2/2
3	A2M	D	5	3,2	-	1/9/27/28	0/3/3/3
3	N7X	D	2	3,2	-	2/12/30/31	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	OKQ	D	18	3,2	-	0/9/27/28	0/2/2/2
3	6NW	D	20	3,2	-	0/9/27/28	0/3/3/3
3	RFJ	D	17	3,2	-	2/9/27/28	0/3/3/3
3	PST	D	7	3,2	-	0/7/21/22	0/2/2/2
3	C7R	D	12	3,2	-	1/7/21/22	0/2/2/2
3	SC	D	11	3,2	-	0/7/21/22	0/2/2/2
3	SC	D	6	3,2	-	0/7/21/22	0/2/2/2
3	SC	D	10	3,2	-	0/7/21/22	0/2/2/2
3	PST	D	15	3,2	-	0/7/21/22	0/2/2/2
3	SC	D	8	3,2	-	0/7/21/22	0/2/2/2
3	OKQ	D	19	3,2	-	0/9/27/28	0/2/2/2
3	AS	D	13	3,2	-	1/7/21/22	0/3/3/3

All (104) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	17	RFJ	O6-C6	8.68	1.40	1.23
3	D	6	SC	C2-N3	7.35	1.50	1.36
3	D	11	SC	C2-N3	7.15	1.50	1.36
3	D	14	SC	C2-N3	6.46	1.49	1.36
3	D	8	SC	C2-N3	6.40	1.49	1.36
3	D	8	SC	C4-N4	6.24	1.49	1.33
3	D	14	SC	C4-N4	6.05	1.48	1.33
3	D	15	PST	C1'-N1	5.77	1.62	1.48
3	D	11	SC	C4-N4	5.77	1.47	1.33
3	D	6	SC	C4-N4	5.74	1.47	1.33
3	D	10	SC	C4-N4	5.13	1.46	1.33
3	D	5	A2M	O5'-C5'	-5.12	1.29	1.44
3	D	18	OKQ	C4-N4	5.10	1.46	1.33
3	D	7	PST	C1'-N1	5.08	1.60	1.48
3	D	19	OKQ	C4-N4	5.02	1.46	1.33
3	D	19	OKQ	C1'-N1	4.76	1.61	1.47
3	D	17	RFJ	C2-N2	4.64	1.45	1.34
3	D	16	OKQ	C4-N4	4.64	1.45	1.33
3	D	16	OKQ	C2-N3	4.39	1.45	1.36
3	D	18	OKQ	O3'-C3'	4.39	1.53	1.43
3	D	16	OKQ	C1'-N1	4.39	1.60	1.47
3	D	13	AS	C1'-N9	4.39	1.54	1.46
3	D	10	SC	C2-N3	4.34	1.44	1.36
3	D	18	OKQ	C1'-N1	4.27	1.59	1.47
3	D	16	OKQ	O3'-C3'	4.23	1.53	1.43
3	D	19	OKQ	C2-N3	4.19	1.44	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	17	RFJ	C5-C4	4.12	1.50	1.38
3	D	15	PST	C6-N1	4.05	1.44	1.38
3	D	18	OKQ	C2-N3	4.04	1.44	1.36
3	D	7	PST	C2-N3	4.00	1.44	1.38
3	D	15	PST	C2-N3	3.95	1.44	1.38
3	D	17	RFJ	C5-N7	3.91	1.46	1.39
3	D	3	T39	C2-N3	3.90	1.44	1.38
3	D	20	6NW	C5-C4	3.90	1.46	1.39
3	D	13	AS	C5-C4	3.76	1.45	1.39
3	D	3	T39	C4-C5	3.70	1.50	1.44
3	D	6	SC	C5-C4	3.67	1.51	1.42
3	D	5	A2M	C2'-C1'	3.59	1.62	1.53
3	D	10	SC	C5-C4	3.56	1.51	1.42
3	D	17	RFJ	C4-N3	3.50	1.42	1.34
3	D	2	N7X	O5'-C5'	-3.48	1.34	1.44
3	D	7	PST	C4-N3	3.39	1.45	1.38
3	D	18	OKQ	C2-N1	3.38	1.47	1.40
3	D	1	OMC	C2-N3	3.36	1.43	1.36
3	D	5	A2M	C8-N7	3.33	1.38	1.31
3	D	1	OMC	C2-N1	3.32	1.47	1.40
3	D	19	OKQ	O3'-C3'	3.31	1.51	1.43
3	D	7	PST	C4-C5	3.29	1.50	1.44
3	D	17	RFJ	C8-N7	3.29	1.41	1.32
3	D	17	RFJ	C4-N9	3.28	1.46	1.38
3	D	8	SC	C5-C4	3.26	1.50	1.42
3	D	3	T39	C2-N1	3.25	1.43	1.38
3	D	15	PST	C4-N3	3.24	1.44	1.38
3	D	10	SC	C6-C5	3.23	1.42	1.35
3	D	3	T39	O4-C4	-3.09	1.17	1.23
3	D	14	SC	C5-C4	3.09	1.50	1.42
3	D	15	PST	C6-C5	3.03	1.39	1.34
3	D	11	SC	C5-C4	2.98	1.49	1.42
3	D	14	SC	C6-C5	2.98	1.42	1.35
3	D	16	OKQ	C2-N1	2.95	1.46	1.40
3	D	19	OKQ	C2-N1	2.88	1.46	1.40
3	D	4	C5L	C6-C5	2.87	1.39	1.34
3	D	5	A2M	C8-N9	2.81	1.42	1.37
3	D	20	6NW	C8-N7	2.79	1.37	1.31
3	D	2	N7X	C6-C5	2.75	1.39	1.34
3	D	7	PST	C6-N1	2.70	1.42	1.38
3	D	7	PST	C5M-C5	2.69	1.57	1.50
3	D	15	PST	C4-C5	2.68	1.49	1.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	20	6NW	C6-N6	2.67	1.41	1.34
3	D	15	PST	C5M-C5	2.65	1.57	1.50
3	D	12	C7R	C2-N3	2.65	1.41	1.36
3	D	13	AS	C8-N9	2.64	1.42	1.37
3	D	20	6NW	C2'-C1'	2.60	1.59	1.53
3	D	5	A2M	C2-N1	2.57	1.38	1.33
3	D	19	OKQ	C6-C5	2.52	1.40	1.35
3	D	2	N7X	C1'-N1	2.50	1.54	1.47
3	D	18	OKQ	C6-C5	2.48	1.40	1.35
3	D	13	AS	O4'-C1'	-2.41	1.37	1.42
3	D	11	SC	C6-C5	2.40	1.40	1.35
3	D	16	OKQ	C6-C5	2.37	1.40	1.35
3	D	1	OMC	C4-N4	2.36	1.39	1.33
3	D	13	AS	C8-N7	2.34	1.36	1.31
3	D	8	SC	C6-C5	2.31	1.40	1.35
3	D	2	N7X	C2'-C1'	2.25	1.58	1.53
3	D	20	6NW	C8-N9	2.24	1.41	1.37
3	D	4	C5L	C2-N3	2.24	1.40	1.36
3	D	14	SC	C6-N1	2.20	1.43	1.38
3	D	10	SC	C2-N1	2.20	1.44	1.40
3	D	6	SC	C6-C5	2.19	1.40	1.35
3	D	5	A2M	C5-C4	2.18	1.43	1.39
3	D	6	SC	C6-N1	2.17	1.43	1.38
3	D	13	AS	C5'-C4'	2.17	1.58	1.51
3	D	18	OKQ	C5-C4	2.16	1.47	1.42
3	D	8	SC	C6-N1	2.15	1.43	1.38
3	D	17	RFJ	O4'-C1'	2.15	1.47	1.42
3	D	17	RFJ	C8-N9	2.13	1.42	1.37
3	D	20	6NW	C2-N1	2.12	1.37	1.33
3	D	19	OKQ	C5-C4	2.12	1.47	1.42
3	D	7	PST	O4'-C1'	-2.10	1.37	1.42
3	D	2	N7X	C4-N4	2.09	1.39	1.34
3	D	7	PST	C6-C5	2.09	1.38	1.34
3	D	10	SC	C6-N1	2.09	1.43	1.38
3	D	2	N7X	C2-N3	2.07	1.40	1.36
3	D	17	RFJ	C3'-C2'	-2.05	1.48	1.53

All (90) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	4	C5L	CA'-O2'-C2'	17.69	153.30	114.02
3	D	17	RFJ	C5-C4-N3	-8.69	114.55	128.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	4	C5L	O2'-C2'-C3'	8.25	132.30	111.05
3	D	3	T39	C7-C5-C4	-7.36	110.92	118.78
3	D	5	A2M	C2'-C1'-N9	7.27	125.71	113.75
3	D	5	A2M	CM'-O2'-C2'	7.19	132.94	114.47
3	D	17	RFJ	N9-C4-N3	6.56	139.06	125.95
3	D	20	6NW	CM'-O2'-C2'	6.29	130.63	114.47
3	D	17	RFJ	C2-N3-C4	5.84	122.36	112.30
3	D	7	PST	C4-N3-C2	-5.80	119.74	127.34
3	D	3	T39	C5-C4-N3	-5.62	110.44	115.32
3	D	15	PST	C4-N3-C2	-5.33	120.35	127.34
3	D	2	N7X	C12-O12-C1	5.33	145.03	112.90
3	D	3	T39	C6-C5-C4	5.15	122.27	118.02
3	D	15	PST	C5-C4-N3	4.99	119.66	115.32
3	D	5	A2M	O2'-C2'-C1'	4.90	118.29	108.99
3	D	7	PST	C5-C4-N3	4.70	119.41	115.32
3	D	14	SC	O4'-C1'-N1	4.48	115.81	107.86
3	D	2	N7X	C6-N1-C2	-4.42	115.06	120.95
3	D	5	A2M	O4'-C1'-C2'	-4.23	99.31	106.59
3	D	10	SC	C4-N3-C2	3.88	126.38	120.26
3	D	14	SC	C4-N3-C2	3.78	126.22	120.26
3	D	20	6NW	C5-C4-N3	-3.71	121.61	126.72
3	D	19	OKQ	C1'-N1-C6	3.53	128.33	120.78
3	D	18	OKQ	C1'-N1-C6	3.41	128.06	120.78
3	D	7	PST	N3-C2-N1	3.38	119.29	114.89
3	D	15	PST	N3-C2-N1	3.35	119.25	114.89
3	D	8	SC	C4-N3-C2	3.29	125.44	120.26
3	D	11	SC	C4-N3-C2	3.28	125.42	120.26
3	D	6	SC	O4'-C1'-N1	3.14	113.43	107.86
3	D	10	SC	O4'-C1'-N1	3.13	113.42	107.86
3	D	16	OKQ	C2'-C1'-N1	3.05	120.03	114.24
3	D	19	OKQ	C2'-C1'-N1	2.99	119.92	114.24
3	D	3	T39	C7-C5-C6	2.92	126.81	122.85
3	D	14	SC	C1'-N1-C6	2.91	127.26	121.53
3	D	17	RFJ	C2-N1-C6	-2.85	119.94	125.11
3	D	5	A2M	O4'-C4'-C5'	2.84	118.45	109.33
3	D	16	OKQ	C1'-N1-C6	2.81	126.78	120.78
3	D	4	C5L	O2'-CA'-CB'	-2.80	100.19	109.67
3	D	6	SC	C4-N3-C2	2.79	124.66	120.26
3	D	18	OKQ	O3'-C3'-C4'	2.77	119.03	111.08
3	D	11	SC	C5-C4-N3	-2.75	116.68	121.32
3	D	1	OMC	C6-N1-C2	-2.67	115.95	120.46
3	D	15	PST	O4'-C4'-C3'	-2.66	99.59	105.65

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	16	OKQ	O2'-C2'-C1'	2.65	114.02	108.99
3	D	18	OKQ	C2'-C1'-N1	2.64	119.25	114.24
3	D	14	SC	C1'-N1-C2	-2.63	113.31	117.83
3	D	16	OKQ	O4'-C1'-N1	-2.61	102.44	108.36
3	D	11	SC	C1'-N1-C6	2.60	126.65	121.53
3	D	4	C5L	O2'-C2'-C1'	-2.60	104.28	108.97
3	D	2	N7X	C3'-C2'-C1'	-2.59	97.84	102.81
3	D	11	SC	C1'-N1-C2	-2.55	113.45	117.83
3	D	13	AS	C5-C4-N3	-2.53	123.23	126.72
3	D	10	SC	C1'-N1-C6	2.47	126.39	121.53
3	D	10	SC	C1'-N1-C2	-2.47	113.59	117.83
3	D	14	SC	C5-C4-N3	-2.46	117.17	121.32
3	D	17	RFJ	C4-C5-N7	-2.46	106.77	110.67
3	D	20	6NW	C4'-O4'-C1'	2.43	114.84	109.47
3	D	5	A2M	N9-C8-N7	-2.41	110.51	113.94
3	D	17	RFJ	C5-C6-N1	2.39	119.34	113.25
3	D	6	SC	C5-C4-N3	-2.39	117.30	121.32
3	D	8	SC	N4-C4-N3	2.38	122.17	117.91
3	D	5	A2M	C2'-C3'-C4'	-2.34	96.96	101.99
3	D	5	A2M	C5-C4-N3	-2.32	123.52	126.72
3	D	8	SC	C5-C4-N3	-2.31	117.43	121.32
3	D	2	N7X	C1'-N1-C6	2.30	124.93	121.15
3	D	13	AS	C2'-C1'-N9	2.29	120.01	114.63
3	D	6	SC	N4-C4-N3	2.27	121.97	117.91
3	D	17	RFJ	C2'-C1'-N9	-2.25	109.97	114.24
3	D	5	A2M	C4'-O4'-C1'	2.24	114.42	109.47
3	D	11	SC	O4'-C1'-N1	2.23	111.82	107.86
3	D	5	A2M	C3'-C2'-C1'	-2.22	98.56	102.81
3	D	18	OKQ	O4'-C4'-C3'	-2.21	100.76	105.15
3	D	18	OKQ	O2'-C2'-C1'	2.19	113.15	108.99
3	D	11	SC	N4-C4-N3	2.15	121.74	117.91
3	D	20	6NW	N3-C4-N9	2.14	130.81	127.17
3	D	13	AS	O4'-C1'-N9	-2.14	104.11	107.96
3	D	18	OKQ	C4-N3-C2	2.14	123.64	120.26
3	D	7	PST	O2-C2-N3	-2.14	117.54	121.49
3	D	19	OKQ	C4-N3-C2	2.11	123.58	120.26
3	D	2	N7X	N1-C2-N3	2.10	122.45	118.80
3	D	1	OMC	C4'-O4'-C1'	2.09	114.09	109.47
3	D	6	SC	C1'-N1-C6	2.08	125.63	121.53
3	D	5	A2M	C4-C5-N7	-2.08	108.20	110.58
3	D	5	A2M	C5-N7-C8	2.07	106.71	103.45
3	D	16	OKQ	O4'-C4'-C3'	-2.01	101.16	105.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	8	SC	N1-C2-N3	-2.01	115.31	118.80
3	D	10	SC	C5-C4-N3	-2.00	117.94	121.32
3	D	11	SC	C2'-C3'-C4'	2.00	106.86	102.80
3	D	17	RFJ	O6-C6-C5	-2.00	121.25	126.53

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	3	T39	C1'-C2'-O2'-CA'
3	D	1	OMC	O4'-C4'-C5'-O5'
3	D	17	RFJ	C3'-C4'-C5'-O5'
3	D	4	C5L	O2'-CA'-CB'-OC'
3	D	2	N7X	O2'-C-C1-O12
3	D	4	C5L	CA'-CB'-OC'-CD'
3	D	2	N7X	C-C1-O12-C12
3	D	5	A2M	C3'-C2'-O2'-CM'
3	D	12	C7R	C3'-C4'-C5'-O5'
3	D	16	OKQ	O4'-C4'-C5'-O5'
3	D	17	RFJ	O4'-C4'-C5'-O5'
3	D	3	T39	O2'-CA'-CB'-OC'
3	D	3	T39	CA'-CB'-OC'-CD'
3	D	13	AS	O4'-C1'-N9-C4

There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	3	T39	1	0
3	D	11	SC	1	0
3	D	10	SC	2	0
3	D	15	PST	0	1

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

3 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	$\# Z > 2$
6	PEG	D	201	-	6,6,6	0.09	0	5,5,5	0.16	0
5	PO4	A	302	-	4,4,4	0.91	0	6,6,6	0.44	0
4	GOL	A	301	-	5,5,5	1.13	0	5,5,5	0.90	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
6	PEG	D	201	-	-	3/4/4/4	-
4	GOL	A	301	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	301	GOL	O1-C1-C2-C3
6	D	201	PEG	O2-C3-C4-O4
4	A	301	GOL	O1-C1-C2-O2
6	D	201	PEG	O1-C1-C2-O2
6	D	201	PEG	C4-C3-O2-C2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	D	201	PEG	1	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	149/149 (100%)	0.28	7 (4%) 36 35	15, 37, 65, 118	8 (5%)
2	B	20/20 (100%)	-0.59	0 100 100	30, 45, 53, 54	0
3	D	1/20 (5%)	-0.81	0 100 100	32, 32, 32, 32	0
All	All	170/189 (89%)	0.17	7 (4%) 41 40	15, 38, 62, 118	8 (4%)

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	137	GLY	3.9
1	A	263	GLY	3.3
1	A	267	PHE	3.0
1	A	264	HIS	2.8
1	A	188[A]	HIS	2.7
1	A	262	PRO	2.6
1	A	285	GLU	2.4

6.2 Non-standard residues in protein, DNA, RNA chains [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	OMC	D	1	18/22	0.84	0.13	41,50,56,64	0
3	N7X	D	2	25/26	0.91	0.10	40,49,55,75	0
3	OKQ	D	16	21/22	0.91	0.09	37,42,53,62	0
3	SC	D	6	19/20	0.93	0.07	22,32,44,49	0
3	PST	D	15	20/21	0.93	0.09	36,44,54,56	0
3	C5L	D	4	25/26	0.93	0.09	36,47,59,77	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	RFJ	D	17	24/25	0.93	0.09	31,44,52,54	0
3	OKQ	D	19	21/22	0.93	0.08	29,37,43,50	0
3	A2M	D	5	23/24	0.94	0.07	28,34,49,50	0
3	T39	D	3	25/26	0.94	0.09	35,43,54,60	0
3	OKQ	D	18	21/22	0.95	0.07	27,39,42,45	0
3	6NW	D	20	23/24	0.95	0.09	33,47,59,62	0
3	SC	D	11	19/20	0.96	0.07	31,35,39,43	0
3	C7R	D	12	19/20	0.96	0.07	27,33,40,54	0
3	AS	D	13	21/22	0.96	0.07	28,36,40,44	0
3	SC	D	14	19/20	0.96	0.07	33,40,52,53	0
3	PST	D	7	20/21	0.96	0.07	27,31,37,39	0
3	SC	D	8	19/20	0.97	0.06	18,28,36,39	0
3	SC	D	10	19/20	0.97	0.06	26,29,36,39	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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($\text{\AA}^2$)	Q<0.9
4	GOL	A	301	6/6	0.71	0.17	55,62,73,75	0
6	PEG	D	201	7/7	0.87	0.13	48,55,57,62	0
5	PO4	A	302	5/5	0.91	0.08	53,56,58,62	0

6.5 Other polymers [i](#)

There are no such residues in this entry.