



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 18, 2026 – 05:43 PM UTC

PDB ID : 4NTP / pdb_00004ntp
Title : Crystal structure of macrocycles containing A 17-23 (LV(PHI)FAED) and A 30-36 (AII(SAR)L(ORN)V)
Authors : Nowick, J.S.; Spencer, R.K.; Li, H.
Deposited on : 2013-12-02
Resolution : 1.99 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

i

X-RAY DIFFRACTION

A.

the following graphic. The table shows the number of entries on which the scores are based.



Percentile relative to all X-ray structures

Percentile relative to X-ray structures of similar resolution

Metric

Whole archive
(#Entries)

Similar resolution
(#Entries, resolution range(Å))

 R_{free}

Clashscore

Ramachandran outliers

Sidechain outliers

RSRZ outliers

180053

190562

187476

187428

180081

1506 (1.98-1.98)

1534 (1.98-1.98)

1518 (1.98-1.98)

1518 (1.98-1.98)

1506 (1.98-1.98)

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.

Quality of chain

75%

25

75%

25

810

010

250

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Mol	Chain	Length	Quality of chain
1	F	16	 69% 31%
1	G	16	 81% 19%
1	H	16	 81% 19%
1	I	16	 81% 12% 6%
1	J	16	 81% 12% 6%
1	K	16	 75% 19% 6%
1	L	16	 75% 25%
1	M	16	 69% 25% 6%
1	N	16	 75% 25%
1	O	16	 75% 25%
1	P	16	 81% 19%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4484 atoms, of which 2190 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 Cyclic hexadecapeptide (ORN)LV(PHI)FAED(ORN)AII(SAR)L(ORN)V.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	16	Total	C	H	I	N	O	0	0	0
			262	85	137	1	19	20			
1	B	16	Total	C	H	I	N	O	0	0	0
			262	85	137	1	19	20			
1	C	16	Total	C	H	I	N	O	0	0	0
			262	85	137	1	19	20			
1	D	16	Total	C	H	I	N	O	0	0	0
			261	85	136	1	19	20			
1	E	16	Total	C	H	I	N	O	0	0	0
			262	85	137	1	19	20			
1	F	16	Total	C	H	I	N	O	0	0	0
			262	85	137	1	19	20			
1	G	16	Total	C	H	I	N	O	0	0	0
			262	85	137	1	19	20			
1	H	16	Total	C	H	I	N	O	0	0	0
			262	85	137	1	19	20			
1	I	16	Total	C	H	I	N	O	0	0	0
			262	85	137	1	19	20			
1	J	16	Total	C	H	I	N	O	0	0	0
			262	85	137	1	19	20			
1	K	16	Total	C	H	I	N	O	0	0	0
			262	85	137	1	19	20			
1	L	16	Total	C	H	I	N	O	0	0	0
			261	85	136	1	19	20			
1	M	16	Total	C	H	I	N	O	0	0	0
			262	85	137	1	19	20			
1	N	16	Total	C	H	I	N	O	0	0	0
			262	85	137	1	19	20			
1	O	16	Total	C	H	I	N	O	0	0	0
			262	85	137	1	19	20			
1	P	16	Total	C	H	I	N	O	0	0	0
			262	85	137	1	19	20			

- Molecule 2 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total 1	Na 1	0	0
2	E	1	Total 1	Na 1	0	0

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total 1	Cl 1	0	0
3	E	1	Total 1	Cl 1	0	0
3	G	1	Total 1	Cl 1	0	0
3	H	1	Total 1	Cl 1	0	0
3	P	1	Total 1	Cl 1	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	25	Total 25	O 25	0	0
4	B	16	Total 16	O 16	0	0
4	C	15	Total 15	O 15	0	0
4	D	22	Total 22	O 22	0	0
4	E	20	Total 20	O 20	0	0
4	F	18	Total 18	O 18	0	0
4	G	26	Total 26	O 26	0	0
4	H	20	Total 20	O 20	0	0
4	I	17	Total 17	O 17	0	0
4	J	14	Total 14	O 14	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	K	13	Total 13	O 13	0	0
4	L	13	Total 13	O 13	0	0
4	M	17	Total 17	O 17	0	0
4	N	15	Total 15	O 15	0	0
4	O	22	Total 22	O 22	0	0
4	P	14	Total 14	O 14	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: Cyclic hexadecapeptide (ORN)LV(PHI)FAED(ORN)AII(SAR)L(ORN)V

Chain A: 



- Molecule 1: Cyclic hexadecapeptide (ORN)LV(PHI)FAED(ORN)AII(SAR)L(ORN)V

Chain B: 



- Molecule 1: Cyclic hexadecapeptide (ORN)LV(PHI)FAED(ORN)AII(SAR)L(ORN)V

Chain C: 



- Molecule 1: Cyclic hexadecapeptide (ORN)LV(PHI)FAED(ORN)AII(SAR)L(ORN)V

Chain D: 



- Molecule 1: Cyclic hexadecapeptide (ORN)LV(PHI)FAED(ORN)AII(SAR)L(ORN)V

Chain E: 

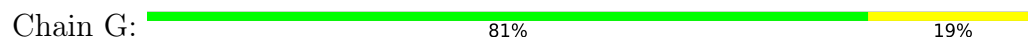


- Molecule 1: Cyclic hexadecapeptide (ORN)LV(PHI)FAED(ORN)AII(SAR)L(ORN)V

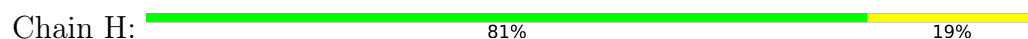
Chain F: 



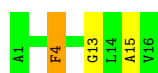
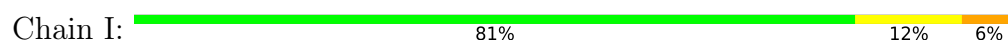
- Molecule 1: Cyclic hexadecapeptide (ORN)LV(PHI)FAED(ORN)AII(SAR)L(ORN)V



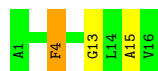
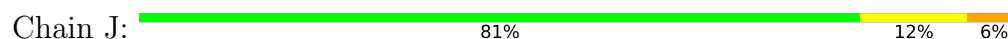
- Molecule 1: Cyclic hexadecapeptide (ORN)LV(PHI)FAED(ORN)AII(SAR)L(ORN)V



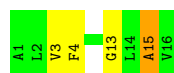
- Molecule 1: Cyclic hexadecapeptide (ORN)LV(PHI)FAED(ORN)AII(SAR)L(ORN)V



- Molecule 1: Cyclic hexadecapeptide (ORN)LV(PHI)FAED(ORN)AII(SAR)L(ORN)V



- Molecule 1: Cyclic hexadecapeptide (ORN)LV(PHI)FAED(ORN)AII(SAR)L(ORN)V



- Molecule 1: Cyclic hexadecapeptide (ORN)LV(PHI)FAED(ORN)AII(SAR)L(ORN)V

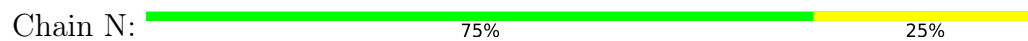


- Molecule 1: Cyclic hexadecapeptide (ORN)LV(PHI)FAED(ORN)AII(SAR)L(ORN)V

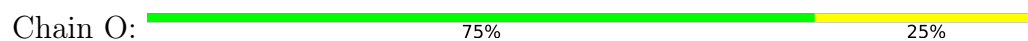




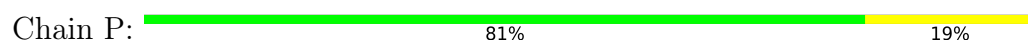
- Molecule 1: Cyclic hexadecapeptide (ORN)LV(PHI)FAED(ORN)AII(SAR)L(ORN)V



- Molecule 1: Cyclic hexadecapeptide (ORN)LV(PHI)FAED(ORN)AII(SAR)L(ORN)V



- Molecule 1: Cyclic hexadecapeptide (ORN)LV(PHI)FAED(ORN)AII(SAR)L(ORN)V



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	68.18Å 68.18Å 170.43Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	34.09 – 1.99 34.09 – 1.99	Depositor EDS
% Data completeness (in resolution range)	98.2 (34.09-1.99) 95.4 (34.09-1.99)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.82 (at 1.98Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.207 , 0.246 0.225 , 0.236	Depositor DCC
R_{free} test set	1010 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	21.9	Xtriage
Anisotropy	0.089	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 34.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.025 for $-1/3^*h+1/3^*k+1/3^*l, -k, 8/3^*h+4/3^*k+1/3^*l$ 0.040 for $-2/3^*h-1/3^*k-1/3^*l, -1/3^*h-2/3^*k+1/3^*l, -4/3^*h+4/3^*k+1/3^*l$ 0.019 for $-h, 1/3^*h-1/3^*k-1/3^*l, -4/3^*h-8/3^*k+1/3^*l$ 0.021 for $-h, 2/3^*h+1/3^*k+1/3^*l, 4/3^*h+8/3^*k-1/3^*l$ 0.025 for $1/3^*h+2/3^*k-1/3^*l, -k, -8/3^*h-4/3^*k-1/3^*l$ 0.040 for $-1/3^*h-2/3^*k+1/3^*l, -2/3^*h-1/3^*k-1/3^*l, 4/3^*h-4/3^*k-1/3^*l$ 0.489 for $-h-k, k, -l$	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4484	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 43.40 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.7731e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: SAR, PHI, ORN, NA, 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	0.30	0/80	0.48	0/102
1	B	0.29	0/80	0.44	0/102
1	C	0.31	0/80	0.41	0/102
1	D	0.22	0/80	0.42	0/102
1	E	0.27	0/80	0.51	0/102
1	F	0.31	0/80	0.44	0/102
1	G	0.22	0/80	0.42	0/102
1	H	0.27	0/80	0.47	0/102
1	I	0.36	0/80	0.52	0/102
1	J	0.29	0/80	0.47	0/102
1	K	0.27	0/80	0.45	0/102
1	L	0.29	0/80	0.46	0/102
1	M	0.33	0/80	0.51	0/102
1	N	0.29	0/80	0.49	0/102
1	O	0.35	0/80	0.54	0/102
1	P	0.28	0/80	0.47	0/102
All	All	0.29	0/1280	0.47	0/1632

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	125	137	130	1	0
1	B	125	137	130	0	1
1	C	125	137	130	0	0
1	D	125	136	130	0	0
1	E	125	137	130	0	0
1	F	125	137	130	2	0
1	G	125	137	130	0	0
1	H	125	137	130	0	0
1	I	125	137	130	1	0
1	J	125	137	130	1	0
1	K	125	137	130	1	0
1	L	125	136	130	0	0
1	M	125	137	130	1	1
1	N	125	137	130	0	0
1	O	125	137	130	1	0
1	P	125	137	130	0	0
2	A	1	0	0	0	0
2	E	1	0	0	0	0
3	B	1	0	0	1	0
3	E	1	0	0	0	0
3	G	1	0	0	0	0
3	H	1	0	0	1	0
3	P	1	0	0	0	0
4	A	25	0	0	1	0
4	B	16	0	0	1	0
4	C	15	0	0	0	0
4	D	22	0	0	0	0
4	E	20	0	0	0	0
4	F	18	0	0	2	0
4	G	26	0	0	0	0
4	H	20	0	0	1	0
4	I	17	0	0	1	0
4	J	14	0	0	0	0
4	K	13	0	0	0	0
4	L	13	0	0	0	0
4	M	17	0	0	0	0
4	N	15	0	0	0	1
4	O	22	0	0	0	1
4	P	14	0	0	0	0
All	All	2294	2190	2080	9	3

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 2.

All (9) 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:F:8:ASP:OD1	4:F:116:HOH:O	1.91	0.88
3:H:101:CL:CL	4:H:213:HOH:O	2.32	0.83
1:F:7:GLU:OE2	4:F:118:HOH:O	1.95	0.83
1:I:4:PHI:I	4:I:117:HOH:O	2.73	0.76
3:B:101:CL:CL	4:B:214:HOH:O	2.43	0.74
1:K:3:VAL:HG22	1:K:15:ORN:HB3	1.91	0.52
1:J:4:PHI:I	1:O:16:VAL:HG11	2.84	0.47
1:M:2:LEU:CD1	1:M:4:PHI:I	3.38	0.41
1:A:7:GLU:OE2	4:A:216:HOH:O	2.22	0.41

All (3) 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)
4:N:106:HOH:O	4:N:106:HOH:O[2_645]	1.93	0.27
4:O:105:HOH:O	4:O:105:HOH:O[3_755]	2.05	0.15
1:B:8:ASP:OD2	1:M:1:ORN:H2[3_765]	1.58	0.02

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	9/16 (56%)	9 (100%)	0	0	100	100
1	B	9/16 (56%)	9 (100%)	0	0	100	100
1	C	9/16 (56%)	9 (100%)	0	0	100	100
1	D	9/16 (56%)	9 (100%)	0	0	100	100
1	E	9/16 (56%)	9 (100%)	0	0	100	100
1	F	9/16 (56%)	9 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	G	9/16 (56%)	9 (100%)	0	0	100	100
1	H	9/16 (56%)	9 (100%)	0	0	100	100
1	I	9/16 (56%)	9 (100%)	0	0	100	100
1	J	9/16 (56%)	9 (100%)	0	0	100	100
1	K	9/16 (56%)	9 (100%)	0	0	100	100
1	L	9/16 (56%)	9 (100%)	0	0	100	100
1	M	9/16 (56%)	9 (100%)	0	0	100	100
1	N	9/16 (56%)	9 (100%)	0	0	100	100
1	O	9/16 (56%)	9 (100%)	0	0	100	100
1	P	9/16 (56%)	9 (100%)	0	0	100	100
All	All	144/256 (56%)	144 (100%)	0	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	9/9 (100%)	9 (100%)	0	100	100
1	B	9/9 (100%)	9 (100%)	0	100	100
1	C	9/9 (100%)	9 (100%)	0	100	100
1	D	9/9 (100%)	9 (100%)	0	100	100
1	E	9/9 (100%)	9 (100%)	0	100	100
1	F	9/9 (100%)	9 (100%)	0	100	100
1	G	9/9 (100%)	9 (100%)	0	100	100
1	H	9/9 (100%)	9 (100%)	0	100	100
1	I	9/9 (100%)	9 (100%)	0	100	100
1	J	9/9 (100%)	9 (100%)	0	100	100
1	K	9/9 (100%)	9 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	L	9/9 (100%)	8 (89%)	1 (11%)	6	1
1	M	9/9 (100%)	9 (100%)	0	100	100
1	N	9/9 (100%)	8 (89%)	1 (11%)	6	1
1	O	9/9 (100%)	9 (100%)	0	100	100
1	P	9/9 (100%)	9 (100%)	0	100	100
All	All	144/144 (100%)	142 (99%)	2 (1%)	59	54

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	2	LEU
1	N	2	LEU

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

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

80 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	$\# Z > 2$
1	PHI	C	4	1	11,12,13	1.76	3 (27%)	10,15,17	0.81	0
1	ORN	N	1	1	6,7,8	0.51	0	2,7,9	0.07	0
1	ORN	A	1	1	6,7,8	0.55	0	2,7,9	0.20	0
1	ORN	C	1	1	6,7,8	0.53	0	2,7,9	0.14	0
1	ORN	L	1	1	6,7,8	0.52	0	2,7,9	0.20	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	PHI	E	4	1	11,12,13	1.73	3 (27%)	10,15,17	0.89	0
1	ORN	N	15	1	6,7,8	2.30	1 (16%)	2,7,9	0.56	0
1	ORN	C	9	1	6,7,8	0.61	0	2,7,9	0.20	0
1	SAR	M	13	1	3,4,5	2.01	1 (33%)	1,3,5	1.31	0
1	ORN	E	1	1	6,7,8	0.54	0	2,7,9	0.12	0
1	PHI	M	4	1	11,12,13	1.75	3 (27%)	10,15,17	0.98	0
1	ORN	M	15	1	6,7,8	2.34	1 (16%)	2,7,9	0.78	0
1	ORN	A	9	1	6,7,8	0.50	0	2,7,9	0.01	0
1	SAR	F	13	1	3,4,5	2.03	1 (33%)	1,3,5	1.42	0
1	PHI	D	4	1	11,12,13	1.81	3 (27%)	10,15,17	0.97	0
1	ORN	P	15	1	6,7,8	2.32	1 (16%)	2,7,9	0.47	0
1	ORN	D	1	1	6,7,8	0.59	0	2,7,9	0.15	0
1	ORN	B	1	1	6,7,8	0.49	0	2,7,9	0.13	0
1	ORN	F	15	1	6,7,8	2.31	1 (16%)	2,7,9	1.00	0
1	ORN	F	1	1	6,7,8	0.46	0	2,7,9	0.25	0
1	ORN	I	15	1	6,7,8	2.25	1 (16%)	2,7,9	0.91	0
1	ORN	G	9	1	6,7,8	0.52	0	2,7,9	0.15	0
1	PHI	J	4	1	11,12,13	1.74	3 (27%)	10,15,17	1.02	0
1	SAR	I	13	1	3,4,5	2.07	1 (33%)	1,3,5	1.42	0
1	ORN	D	9	1	6,7,8	0.46	0	2,7,9	0.13	0
1	SAR	K	13	1	3,4,5	2.02	1 (33%)	1,3,5	1.30	0
1	ORN	O	9	1	6,7,8	0.63	0	2,7,9	0.11	0
1	PHI	K	4	1	11,12,13	1.74	3 (27%)	10,15,17	0.84	0
1	ORN	D	15	1	6,7,8	2.30	1 (16%)	2,7,9	0.41	0
1	ORN	K	15	1	6,7,8	2.42	1 (16%)	2,7,9	1.03	0
1	ORN	J	15	1	6,7,8	2.22	1 (16%)	2,7,9	0.56	0
1	ORN	P	1	1	6,7,8	0.63	0	2,7,9	0.07	0
1	SAR	D	13	1	3,4,5	2.04	1 (33%)	1,3,5	1.35	0
1	ORN	O	1	1	6,7,8	0.46	0	2,7,9	0.10	0
1	ORN	L	9	1	6,7,8	0.72	0	2,7,9	0.25	0
1	PHI	B	4	1	11,12,13	1.79	3 (27%)	10,15,17	0.80	0
1	ORN	M	9	1	6,7,8	0.72	0	2,7,9	0.16	0
1	PHI	I	4	1	11,12,13	1.78	3 (27%)	10,15,17	1.33	1 (10%)
1	SAR	J	13	1	3,4,5	2.05	1 (33%)	1,3,5	1.44	0
1	ORN	M	1	1	6,7,8	0.47	0	2,7,9	0.14	0
1	ORN	G	1	1	6,7,8	0.54	0	2,7,9	0.17	0
1	ORN	C	15	1	6,7,8	2.37	1 (16%)	2,7,9	0.64	0
1	SAR	N	13	1	3,4,5	1.98	1 (33%)	1,3,5	1.38	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	SAR	A	13	1	3,4,5	2.02	1 (33%)	1,3,5	1.39	0
1	ORN	P	9	1	6,7,8	0.71	0	2,7,9	0.33	0
1	ORN	N	9	1	6,7,8	0.67	0	2,7,9	0.21	0
1	SAR	E	13	1	3,4,5	2.09	1 (33%)	1,3,5	1.52	0
1	SAR	O	13	1	3,4,5	1.93	1 (33%)	1,3,5	1.24	0
1	SAR	L	13	1	3,4,5	2.05	1 (33%)	1,3,5	1.27	0
1	SAR	C	13	1	3,4,5	1.97	1 (33%)	1,3,5	1.30	0
1	SAR	H	13	1	3,4,5	2.03	1 (33%)	1,3,5	1.49	0
1	ORN	F	9	1	6,7,8	0.58	0	2,7,9	0.10	0
1	ORN	E	15	1	6,7,8	2.24	1 (16%)	2,7,9	0.25	0
1	ORN	H	9	1	6,7,8	0.67	0	2,7,9	0.08	0
1	PHI	H	4	1	11,12,13	1.70	3 (27%)	10,15,17	0.89	0
1	ORN	H	15	1	6,7,8	2.28	1 (16%)	2,7,9	0.33	0
1	PHI	P	4	1	11,12,13	1.77	3 (27%)	10,15,17	1.02	0
1	SAR	B	13	1	3,4,5	2.00	1 (33%)	1,3,5	1.19	0
1	ORN	K	1	1	6,7,8	0.65	0	2,7,9	0.15	0
1	PHI	A	4	1	11,12,13	1.78	3 (27%)	10,15,17	1.00	0
1	ORN	J	9	1	6,7,8	0.56	0	2,7,9	0.12	0
1	PHI	L	4	1	11,12,13	1.78	3 (27%)	10,15,17	0.81	0
1	ORN	I	9	1	6,7,8	0.90	0	2,7,9	0.37	0
1	ORN	A	15	1	6,7,8	2.27	1 (16%)	2,7,9	0.69	0
1	ORN	L	15	1	6,7,8	2.26	1 (16%)	2,7,9	0.60	0
1	ORN	K	9	1	6,7,8	0.63	0	2,7,9	0.34	0
1	ORN	B	15	1	6,7,8	2.25	1 (16%)	2,7,9	1.01	0
1	PHI	O	4	1	11,12,13	1.74	3 (27%)	10,15,17	1.20	0
1	ORN	O	15	1	6,7,8	2.30	1 (16%)	2,7,9	0.92	0
1	PHI	F	4	1	11,12,13	1.77	3 (27%)	10,15,17	0.83	0
1	ORN	B	9	1	6,7,8	0.88	0	2,7,9	0.21	0
1	PHI	G	4	1	11,12,13	1.81	3 (27%)	10,15,17	1.06	0
1	ORN	G	15	1	6,7,8	2.29	1 (16%)	2,7,9	0.81	0
1	PHI	N	4	1	11,12,13	1.74	3 (27%)	10,15,17	1.00	0
1	ORN	J	1	1	6,7,8	0.37	0	2,7,9	0.13	0
1	ORN	H	1	1	6,7,8	0.42	0	2,7,9	0.06	0
1	SAR	G	13	1	3,4,5	2.03	1 (33%)	1,3,5	1.41	0
1	ORN	E	9	1	6,7,8	0.56	0	2,7,9	0.17	0
1	SAR	P	13	1	3,4,5	1.90	1 (33%)	1,3,5	1.47	0
1	ORN	I	1	1	6,7,8	0.54	0	2,7,9	0.19	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
1	PHI	C	4	1	-	0/5/6/8	0/1/1/1
1	ORN	N	1	1	-	0/5/6/8	-
1	ORN	A	1	1	-	0/5/6/8	-
1	ORN	C	1	1	-	1/5/6/8	-
1	ORN	L	1	1	-	2/5/6/8	-
1	PHI	E	4	1	-	0/5/6/8	0/1/1/1
1	ORN	N	15	1	-	1/5/6/8	-
1	ORN	C	9	1	-	0/5/6/8	-
1	SAR	M	13	1	-	1/1/2/3	-
1	ORN	E	1	1	-	0/5/6/8	-
1	PHI	M	4	1	-	0/5/6/8	0/1/1/1
1	ORN	M	15	1	-	1/5/6/8	-
1	ORN	A	9	1	-	0/5/6/8	-
1	SAR	F	13	1	-	1/1/2/3	-
1	PHI	D	4	1	-	0/5/6/8	0/1/1/1
1	ORN	P	15	1	-	2/5/6/8	-
1	ORN	D	1	1	-	0/5/6/8	-
1	ORN	B	1	1	-	0/5/6/8	-
1	ORN	F	15	1	-	1/5/6/8	-
1	ORN	F	1	1	-	0/5/6/8	-
1	ORN	I	15	1	-	1/5/6/8	-
1	ORN	G	9	1	-	0/5/6/8	-
1	PHI	J	4	1	-	0/5/6/8	0/1/1/1
1	SAR	I	13	1	-	1/1/2/3	-
1	ORN	D	9	1	-	0/5/6/8	-
1	SAR	K	13	1	-	1/1/2/3	-
1	ORN	O	9	1	-	0/5/6/8	-
1	PHI	K	4	1	-	0/5/6/8	0/1/1/1
1	ORN	D	15	1	-	2/5/6/8	-
1	ORN	K	15	1	-	2/5/6/8	-
1	ORN	J	15	1	-	2/5/6/8	-
1	ORN	P	1	1	-	0/5/6/8	-
1	SAR	D	13	1	-	1/1/2/3	-
1	ORN	O	1	1	-	0/5/6/8	-
1	ORN	L	9	1	-	0/5/6/8	-
1	PHI	B	4	1	-	0/5/6/8	0/1/1/1
1	ORN	M	9	1	-	1/5/6/8	-
1	PHI	I	4	1	-	0/5/6/8	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	SAR	J	13	1	-	1/1/2/3	-
1	ORN	M	1	1	-	0/5/6/8	-
1	ORN	G	1	1	-	0/5/6/8	-
1	ORN	C	15	1	-	1/5/6/8	-
1	SAR	N	13	1	-	1/1/2/3	-
1	SAR	A	13	1	-	1/1/2/3	-
1	ORN	P	9	1	-	0/5/6/8	-
1	ORN	N	9	1	-	0/5/6/8	-
1	SAR	E	13	1	-	1/1/2/3	-
1	SAR	O	13	1	-	1/1/2/3	-
1	SAR	L	13	1	-	1/1/2/3	-
1	SAR	C	13	1	-	1/1/2/3	-
1	SAR	H	13	1	-	1/1/2/3	-
1	ORN	F	9	1	-	0/5/6/8	-
1	ORN	E	15	1	-	2/5/6/8	-
1	ORN	H	9	1	-	1/5/6/8	-
1	PHI	H	4	1	-	0/5/6/8	0/1/1/1
1	ORN	H	15	1	-	2/5/6/8	-
1	PHI	P	4	1	-	0/5/6/8	0/1/1/1
1	SAR	B	13	1	-	1/1/2/3	-
1	ORN	K	1	1	-	0/5/6/8	-
1	PHI	A	4	1	-	0/5/6/8	0/1/1/1
1	ORN	J	9	1	-	0/5/6/8	-
1	PHI	L	4	1	-	0/5/6/8	0/1/1/1
1	ORN	I	9	1	-	1/5/6/8	-
1	ORN	A	15	1	-	1/5/6/8	-
1	ORN	L	15	1	-	2/5/6/8	-
1	ORN	K	9	1	-	0/5/6/8	-
1	ORN	B	15	1	-	1/5/6/8	-
1	PHI	O	4	1	-	0/5/6/8	0/1/1/1
1	ORN	O	15	1	-	1/5/6/8	-
1	PHI	F	4	1	-	0/5/6/8	0/1/1/1
1	ORN	B	9	1	-	0/5/6/8	-
1	PHI	G	4	1	-	0/5/6/8	0/1/1/1
1	ORN	G	15	1	-	2/5/6/8	-
1	PHI	N	4	1	-	0/5/6/8	0/1/1/1
1	ORN	J	1	1	-	2/5/6/8	-
1	ORN	H	1	1	-	0/5/6/8	-
1	SAR	G	13	1	-	1/1/2/3	-
1	ORN	E	9	1	-	1/5/6/8	-
1	SAR	P	13	1	-	1/1/2/3	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	ORN	I	1	1	-	0/5/6/8	-

All (80) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	K	15	ORN	CA-N	-5.79	1.31	1.48
1	C	15	ORN	CA-N	-5.73	1.32	1.48
1	M	15	ORN	CA-N	-5.62	1.32	1.48
1	F	15	ORN	CA-N	-5.59	1.32	1.48
1	P	15	ORN	CA-N	-5.58	1.32	1.48
1	D	15	ORN	CA-N	-5.56	1.32	1.48
1	G	15	ORN	CA-N	-5.54	1.32	1.48
1	O	15	ORN	CA-N	-5.53	1.32	1.48
1	N	15	ORN	CA-N	-5.52	1.32	1.48
1	H	15	ORN	CA-N	-5.52	1.32	1.48
1	E	15	ORN	CA-N	-5.45	1.32	1.48
1	L	15	ORN	CA-N	-5.43	1.32	1.48
1	B	15	ORN	CA-N	-5.43	1.32	1.48
1	A	15	ORN	CA-N	-5.42	1.32	1.48
1	I	15	ORN	CA-N	-5.42	1.32	1.48
1	J	15	ORN	CA-N	-5.33	1.33	1.48
1	B	4	PHI	CD1-CG	3.66	1.46	1.38
1	C	4	PHI	CD1-CG	3.60	1.46	1.38
1	E	13	SAR	CA-C	-3.59	1.36	1.49
1	I	4	PHI	CD1-CG	3.59	1.46	1.38
1	I	13	SAR	CA-C	-3.56	1.36	1.49
1	D	4	PHI	CD1-CG	3.56	1.45	1.38
1	A	4	PHI	CD1-CG	3.55	1.45	1.38
1	M	4	PHI	CD1-CG	3.54	1.45	1.38
1	J	13	SAR	CA-C	-3.53	1.36	1.49
1	D	13	SAR	CA-C	-3.52	1.36	1.49
1	L	13	SAR	CA-C	-3.52	1.36	1.49
1	G	13	SAR	CA-C	-3.51	1.36	1.49
1	L	4	PHI	CD1-CG	3.51	1.45	1.38
1	F	13	SAR	CA-C	-3.50	1.36	1.49
1	H	13	SAR	CA-C	-3.49	1.36	1.49
1	G	4	PHI	CD1-CG	3.48	1.45	1.38
1	K	13	SAR	CA-C	-3.48	1.37	1.49
1	A	13	SAR	CA-C	-3.47	1.37	1.49
1	N	4	PHI	CD1-CG	3.46	1.45	1.38
1	M	13	SAR	CA-C	-3.46	1.37	1.49
1	P	4	PHI	CD1-CG	3.45	1.45	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	13	SAR	CA-C	-3.44	1.37	1.49
1	E	4	PHI	CD1-CG	3.43	1.45	1.38
1	N	13	SAR	CA-C	-3.41	1.37	1.49
1	J	4	PHI	CD1-CG	3.41	1.45	1.38
1	H	4	PHI	CD1-CG	3.40	1.45	1.38
1	C	13	SAR	CA-C	-3.39	1.37	1.49
1	K	4	PHI	CD1-CG	3.39	1.45	1.38
1	F	4	PHI	CD1-CG	3.39	1.45	1.38
1	O	4	PHI	CD1-CG	3.37	1.45	1.38
1	O	13	SAR	CA-C	-3.33	1.37	1.49
1	P	13	SAR	CA-C	-3.27	1.37	1.49
1	D	4	PHI	CE2-CZ	3.22	1.45	1.38
1	G	4	PHI	CE2-CZ	3.17	1.45	1.38
1	P	4	PHI	CE2-CZ	3.16	1.45	1.38
1	B	4	PHI	CE2-CZ	3.10	1.45	1.38
1	F	4	PHI	CE2-CZ	3.09	1.45	1.38
1	A	4	PHI	CE2-CZ	3.08	1.45	1.38
1	K	4	PHI	CE2-CZ	3.07	1.45	1.38
1	L	4	PHI	CE2-CZ	3.07	1.45	1.38
1	E	4	PHI	CE2-CZ	3.07	1.45	1.38
1	O	4	PHI	CE2-CZ	3.05	1.45	1.38
1	I	4	PHI	CE2-CZ	3.05	1.45	1.38
1	J	4	PHI	CE2-CZ	3.04	1.45	1.38
1	C	4	PHI	CE2-CZ	3.02	1.45	1.38
1	N	4	PHI	CE2-CZ	3.00	1.45	1.38
1	H	4	PHI	CE2-CZ	2.99	1.45	1.38
1	M	4	PHI	CE2-CZ	2.97	1.45	1.38
1	I	4	PHI	O-C	2.57	1.29	1.20
1	G	4	PHI	O-C	2.54	1.29	1.20
1	O	4	PHI	O-C	2.46	1.29	1.20
1	D	4	PHI	O-C	2.45	1.29	1.20
1	F	4	PHI	O-C	2.44	1.29	1.20
1	N	4	PHI	O-C	2.43	1.29	1.20
1	M	4	PHI	O-C	2.41	1.29	1.20
1	A	4	PHI	O-C	2.40	1.29	1.20
1	K	4	PHI	O-C	2.39	1.29	1.20
1	P	4	PHI	O-C	2.39	1.29	1.20
1	C	4	PHI	O-C	2.38	1.29	1.20
1	B	4	PHI	O-C	2.38	1.29	1.20
1	L	4	PHI	O-C	2.35	1.28	1.20
1	H	4	PHI	O-C	2.34	1.28	1.20
1	J	4	PHI	O-C	2.30	1.28	1.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	4	PHI	O-C	2.23	1.28	1.20

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	I	4	PHI	CD2-CG-CD1	2.07	121.31	118.23

There are no chirality outliers.

All (49) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	13	SAR	C-CA-N-CN
1	A	15	ORN	C-CA-CB-CG
1	B	13	SAR	C-CA-N-CN
1	B	15	ORN	C-CA-CB-CG
1	C	13	SAR	C-CA-N-CN
1	C	15	ORN	C-CA-CB-CG
1	D	13	SAR	C-CA-N-CN
1	D	15	ORN	C-CA-CB-CG
1	E	13	SAR	C-CA-N-CN
1	E	15	ORN	C-CA-CB-CG
1	E	15	ORN	O-C-CA-CB
1	F	13	SAR	C-CA-N-CN
1	F	15	ORN	C-CA-CB-CG
1	G	13	SAR	C-CA-N-CN
1	G	15	ORN	C-CA-CB-CG
1	G	15	ORN	O-C-CA-CB
1	H	13	SAR	C-CA-N-CN
1	H	15	ORN	C-CA-CB-CG
1	H	15	ORN	O-C-CA-CB
1	I	13	SAR	C-CA-N-CN
1	I	15	ORN	C-CA-CB-CG
1	J	1	ORN	O-C-CA-CB
1	J	13	SAR	C-CA-N-CN
1	J	15	ORN	C-CA-CB-CG
1	K	13	SAR	C-CA-N-CN
1	K	15	ORN	C-CA-CB-CG
1	L	1	ORN	O-C-CA-CB
1	L	13	SAR	C-CA-N-CN
1	L	15	ORN	C-CA-CB-CG
1	M	13	SAR	C-CA-N-CN
1	M	15	ORN	C-CA-CB-CG

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Mol	Chain	Res	Type	Atoms
1	N	13	SAR	C-CA-N-CN
1	N	15	ORN	C-CA-CB-CG
1	O	13	SAR	C-CA-N-CN
1	O	15	ORN	C-CA-CB-CG
1	P	13	SAR	C-CA-N-CN
1	P	15	ORN	C-CA-CB-CG
1	P	15	ORN	O-C-CA-CB
1	E	9	ORN	NE-CD-CG-CB
1	H	9	ORN	NE-CD-CG-CB
1	M	9	ORN	CA-CB-CG-CD
1	I	9	ORN	CA-CB-CG-CD
1	J	15	ORN	NE-CD-CG-CB
1	D	15	ORN	N-CA-CB-CG
1	K	15	ORN	N-CA-CB-CG
1	L	15	ORN	N-CA-CB-CG
1	C	1	ORN	C-CA-CB-CG
1	J	1	ORN	C-CA-CB-CG
1	L	1	ORN	C-CA-CB-CG

There are no ring outliers.

5 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	M	4	PHI	1	0
1	J	4	PHI	1	0
1	K	15	ORN	1	0
1	I	4	PHI	1	0
1	M	1	ORN	0	1

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 7 are monoatomic - leaving 0 for Mogul analysis.

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

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	11/16 (68%)	-1.35	0 100 100	18, 30, 32, 36	0
1	B	11/16 (68%)	-1.35	0 100 100	21, 27, 33, 36	0
1	C	11/16 (68%)	-1.26	0 100 100	19, 35, 39, 41	0
1	D	11/16 (68%)	-1.39	0 100 100	18, 28, 38, 38	0
1	E	11/16 (68%)	-1.29	0 100 100	20, 27, 44, 50	0
1	F	11/16 (68%)	-1.30	0 100 100	18, 32, 38, 49	0
1	G	11/16 (68%)	-1.29	0 100 100	17, 27, 36, 38	0
1	H	11/16 (68%)	-1.37	0 100 100	20, 24, 39, 55	0
1	I	11/16 (68%)	-1.24	0 100 100	21, 32, 40, 44	0
1	J	11/16 (68%)	-1.37	0 100 100	17, 29, 34, 38	0
1	K	11/16 (68%)	-1.24	0 100 100	20, 30, 39, 41	0
1	L	11/16 (68%)	-1.39	0 100 100	22, 29, 33, 40	0
1	M	11/16 (68%)	-1.28	0 100 100	17, 26, 34, 39	0
1	N	11/16 (68%)	-1.31	0 100 100	16, 27, 34, 40	0
1	O	11/16 (68%)	-1.32	0 100 100	18, 27, 39, 41	0
1	P	11/16 (68%)	-1.34	0 100 100	20, 26, 34, 37	0
All	All	176/256 (68%)	-1.32	0 100 100	16, 30, 40, 55	0

There are no RSRZ outliers to report.

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($\text{\AA}^2$)	Q<0.9
1	PHI	M	4	12/13	0.93	0.09	12,28,54,68	0
1	PHI	G	4	12/13	0.94	0.13	20,52,83,89	0
1	PHI	I	4	12/13	0.94	0.09	18,26,49,69	0
1	PHI	J	4	12/13	0.94	0.10	14,24,49,68	0
1	PHI	L	4	12/13	0.94	0.11	19,46,76,86	0
1	PHI	F	4	12/13	0.94	0.12	16,35,59,66	0
1	PHI	P	4	12/13	0.94	0.10	14,32,59,67	0
1	PHI	B	4	12/13	0.95	0.09	20,36,63,84	0
1	PHI	D	4	12/13	0.95	0.10	17,41,70,84	0
1	PHI	K	4	12/13	0.95	0.08	20,37,58,71	0
1	PHI	A	4	12/13	0.96	0.09	15,24,49,71	0
1	PHI	H	4	12/13	0.96	0.08	14,27,48,68	0
1	PHI	E	4	12/13	0.96	0.07	18,32,51,76	0
1	PHI	N	4	12/13	0.96	0.09	15,30,50,64	0
1	PHI	O	4	12/13	0.96	0.07	13,26,50,63	0
1	PHI	C	4	12/13	0.96	0.07	18,32,50,68	0
1	ORN	H	9	8/9	0.97	0.08	37,53,90,90	0
1	ORN	J	9	8/9	0.98	0.05	29,39,62,62	0
1	ORN	A	9	8/9	0.98	0.06	28,48,73,73	0
1	ORN	C	9	8/9	0.98	0.08	37,53,85,85	0
1	ORN	E	9	8/9	0.98	0.07	31,49,74,74	0
1	ORN	M	9	8/9	0.98	0.06	27,46,99,99	0
1	ORN	C	15	8/9	0.98	0.05	20,48,79,79	0
1	ORN	N	9	8/9	0.98	0.05	31,48,61,61	0
1	ORN	I	9	8/9	0.98	0.06	42,55,84,84	0
1	ORN	O	9	8/9	0.98	0.05	31,41,70,70	0
1	ORN	F	9	8/9	0.98	0.07	46,57,80,80	0
1	ORN	P	9	8/9	0.98	0.05	30,41,71,71	0
1	ORN	F	15	8/9	0.99	0.05	18,28,79,79	0
1	ORN	G	1	8/9	0.99	0.04	19,28,38,38	0
1	ORN	A	15	8/9	0.99	0.03	14,26,61,61	0
1	ORN	G	9	8/9	0.99	0.05	17,39,91,91	0
1	SAR	G	13	5/6	0.99	0.02	15,18,29,29	0
1	ORN	G	15	8/9	0.99	0.04	19,31,59,59	0
1	ORN	H	1	8/9	0.99	0.02	16,22,37,37	0
1	SAR	C	13	5/6	0.99	0.02	22,31,34,34	0
1	ORN	B	1	8/9	0.99	0.04	20,37,58,58	0
1	SAR	H	13	5/6	0.99	0.03	15,23,30,30	0
1	ORN	H	15	8/9	0.99	0.04	15,27,65,65	0
1	ORN	I	1	8/9	0.99	0.03	19,27,36,36	0
1	ORN	D	1	8/9	0.99	0.02	16,30,38,38	0
1	ORN	A	1	8/9	0.99	0.04	13,28,34,35	0
1	SAR	I	13	5/6	0.99	0.03	15,19,29,36	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
1	ORN	I	15	8/9	0.99	0.06	17,30,73,73	0
1	ORN	J	1	8/9	0.99	0.03	15,26,35,35	0
1	ORN	D	9	8/9	0.99	0.04	20,39,91,91	0
1	SAR	D	13	5/6	0.99	0.02	16,20,28,30	0
1	ORN	J	15	8/9	0.99	0.03	14,26,61,61	0
1	ORN	K	1	8/9	0.99	0.04	21,38,46,46	0
1	ORN	D	15	8/9	0.99	0.04	18,33,66,66	0
1	ORN	K	9	8/9	0.99	0.05	29,60,77,77	0
1	SAR	K	13	5/6	0.99	0.03	18,32,34,34	0
1	ORN	K	15	8/9	0.99	0.04	23,52,70,70	0
1	ORN	L	1	8/9	0.99	0.04	23,30,49,49	0
1	ORN	E	1	8/9	0.99	0.03	14,22,43,43	0
1	ORN	L	9	8/9	0.99	0.03	19,31,41,41	0
1	SAR	L	13	5/6	0.99	0.03	20,30,37,37	0
1	ORN	L	15	8/9	0.99	0.05	19,51,85,85	0
1	ORN	M	1	8/9	0.99	0.03	17,26,32,35	0
1	ORN	B	9	8/9	0.99	0.04	18,24,42,42	0
1	ORN	B	15	8/9	0.99	0.04	23,41,72,72	0
1	ORN	M	15	8/9	0.99	0.03	17,30,56,56	0
1	ORN	N	1	8/9	0.99	0.04	24,34,56,56	0
1	SAR	E	13	5/6	0.99	0.03	15,23,24,24	0
1	ORN	E	15	8/9	0.99	0.04	17,32,57,57	0
1	SAR	N	13	5/6	0.99	0.02	15,23,24,24	0
1	ORN	N	15	8/9	0.99	0.04	19,36,58,58	0
1	ORN	O	1	8/9	0.99	0.04	17,26,33,34	0
1	ORN	F	1	8/9	0.99	0.03	21,26,30,41	0
1	ORN	C	1	8/9	0.99	0.03	22,32,53,53	0
1	ORN	O	15	8/9	0.99	0.04	16,33,60,60	0
1	ORN	P	1	8/9	0.99	0.04	20,35,48,48	0
1	SAR	A	13	5/6	0.99	0.03	19,25,30,36	0
1	SAR	F	13	5/6	0.99	0.03	15,28,29,40	0
1	ORN	P	15	8/9	0.99	0.03	20,32,59,59	0
1	SAR	J	13	5/6	1.00	0.03	15,19,23,32	0
1	SAR	M	13	5/6	1.00	0.02	17,23,41,41	0
1	SAR	O	13	5/6	1.00	0.02	14,22,34,34	0
1	SAR	P	13	5/6	1.00	0.02	12,19,22,25	0
1	SAR	B	13	5/6	1.00	0.02	18,31,39,39	0

6.3 Carbohydrates ⓘ

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(Å ²)	Q<0.9
2	NA	E	102	1/1	0.98	0.07	42,42,42,42	0
2	NA	A	101	1/1	0.99	0.05	53,53,53,53	0
3	CL	B	101	1/1	0.99	0.04	73,73,73,73	0
3	CL	E	101	1/1	0.99	0.03	47,47,47,47	0
3	CL	G	101	1/1	0.99	0.04	43,43,43,43	0
3	CL	H	101	1/1	0.99	0.04	57,57,57,57	0
3	CL	P	101	1/1	0.99	0.05	54,54,54,54	0

6.5 Other polymers [i](#)

There are no such residues in this entry.