



Full wwPDB X-ray Structure Validation Report

Feb 26, 2014 – 03:16 PM GMT

PDB ID : 4KVO
Title : The NatA (Naa10p/Naa15p) amino-terminal acetyltransferase complex bound to AcCoA
Authors : Liszczak, G.P.; Marmorstein, R.Q.
Deposited on : 2013-05-22
Resolution : 3.15 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

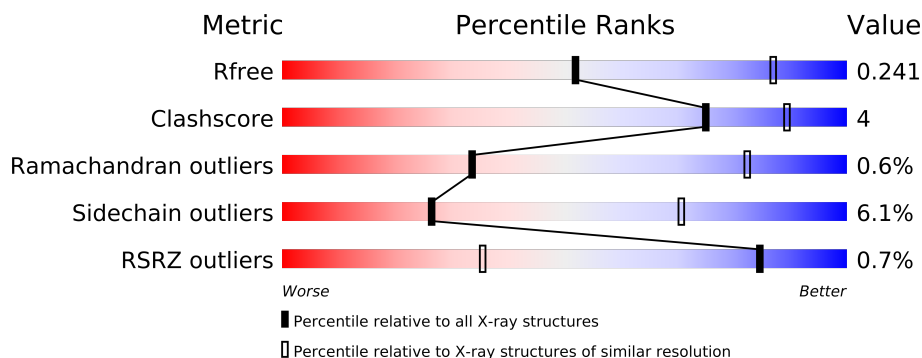
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 3.15 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}	66092	1360 (3.22-3.10)
Clashscore	79885	1681 (3.22-3.10)
Ramachandran outliers	78287	1639 (3.22-3.10)
Sidechain outliers	78261	1638 (3.22-3.10)
RSRZ outliers	66119	1361 (3.22-3.10)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	731	
1	B	731	
1	C	731	
1	D	731	
2	E	156	
2	F	156	
2	G	156	
2	H	156	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	NA	A	801	-	X

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Mol	Type	Chain	Res	Geometry	Electron density
3	NA	E	202	-	X
3	NA	E	203	-	X
4	CL	A	802	-	X
4	CL	A	807	-	X
4	CL	A	808	-	X
4	CL	A	809	-	X
4	CL	A	810	-	X
4	CL	B	803	-	X
4	CL	B	809	-	X
4	CL	B	810	-	X
4	CL	B	811	-	X
4	CL	C	801	-	X
4	CL	C	802	-	X
4	CL	C	803	-	X
4	CL	C	806	-	X
4	CL	C	809	-	X
4	CL	C	810	-	X
4	CL	D	803	-	X
4	CL	D	805	-	X
4	CL	D	808	-	X
4	CL	E	204	-	X
4	CL	E	206	-	X
4	CL	E	207	-	X
4	CL	F	203	-	X
4	CL	H	202	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 28430 atoms, of which 0 are hydrogen and 0 are deuterium.

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 N-terminal acetyltransferase A complex subunit nat1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	718	Total	C	N	O	S	0	1	0
			5764	3679	968	1099	18			
1	B	722	Total	C	N	O	S	0	1	0
			5777	3686	971	1102	18			
1	C	717	Total	C	N	O	S	0	1	0
			5753	3671	968	1096	18			
1	D	717	Total	C	N	O	S	0	1	0
			5761	3678	969	1096	18			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	730	ALA	-	EXPRESSION TAG	UNP O74985
A	731	ALA	-	EXPRESSION TAG	UNP O74985
B	730	ALA	-	EXPRESSION TAG	UNP O74985
B	731	ALA	-	EXPRESSION TAG	UNP O74985
C	730	ALA	-	EXPRESSION TAG	UNP O74985
C	731	ALA	-	EXPRESSION TAG	UNP O74985
D	730	ALA	-	EXPRESSION TAG	UNP O74985
D	731	ALA	-	EXPRESSION TAG	UNP O74985

- Molecule 2 is a protein called N-terminal acetyltransferase A complex catalytic subunit ard1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	153	Total	C	N	O	S	0	1	0
			1247	790	221	225	11			
2	F	153	Total	C	N	O	S	0	1	0
			1247	790	221	225	11			
2	G	153	Total	C	N	O	S	0	1	0
			1247	790	221	225	11			
2	H	153	Total	C	N	O	S	0	1	0
			1247	790	221	225	11			

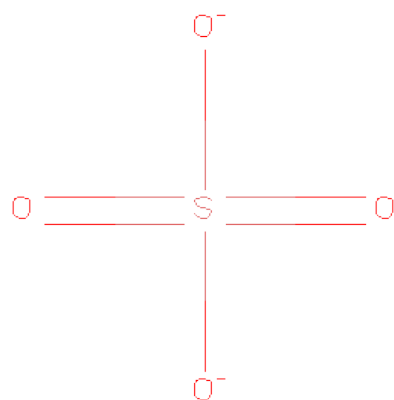
- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

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

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

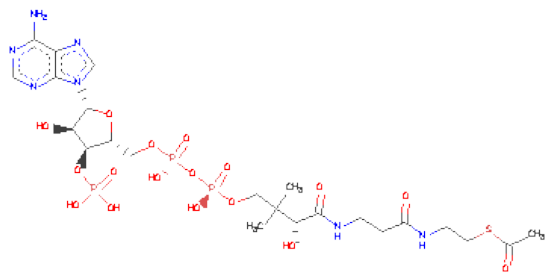
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	G	2	Total 2	Cl 2	0	0
4	D	10	Total 10	Cl 10	0	0
4	E	7	Total 7	Cl 7	0	0
4	H	3	Total 3	Cl 3	0	0
4	B	14	Total 14	Cl 14	0	0
4	C	15	Total 15	Cl 15	0	0
4	A	10	Total 10	Cl 10	0	0
4	F	2	Total 2	Cl 2	0	0

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

- Molecule 6 is ACETYL COENZYME *A (three-letter code: ACO) (formula: C₂₃H₃₈N₇O₁₇P₃S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
6	E	1	Total	C	N	O	P	S	0	0
			51	23	7	17	3	1		
6	F	1	Total	C	N	O	P	S	0	0
			51	23	7	17	3	1		

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Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
6	G	1	Total	C	N	O	P	S	0	0
			51	23	7	17	3	1		
6	H	1	Total	C	N	O	P	S	0	0
			51	23	7	17	3	1		

- Molecule 7 is water.

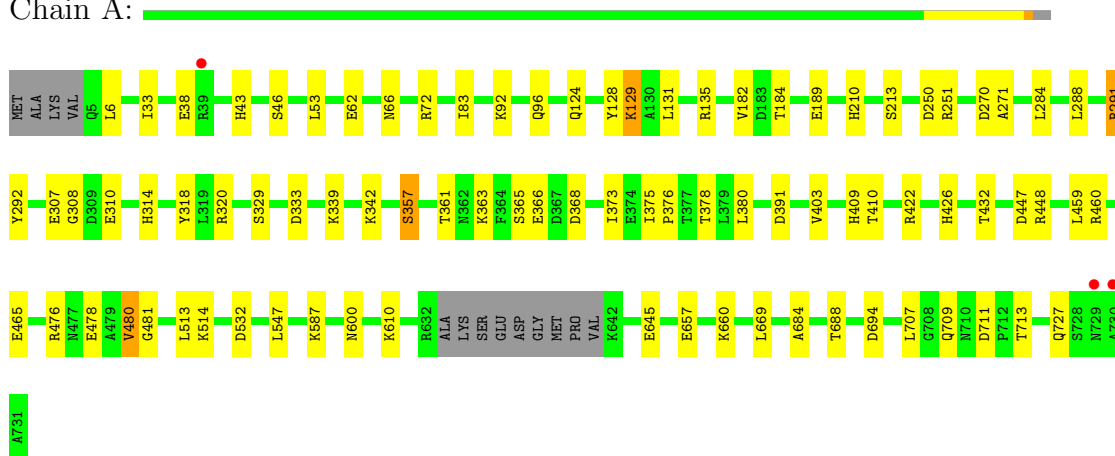
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	22	Total	O	0	0
			22	22		
7	B	22	Total	O	0	0
			22	22		
7	C	25	Total	O	0	0
			25	25		
7	D	11	Total	O	0	0
			11	11		
7	E	4	Total	O	0	0
			4	4		
7	F	6	Total	O	0	0
			6	6		
7	G	5	Total	O	0	0
			5	5		
7	H	10	Total	O	0	0
			10	10		

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

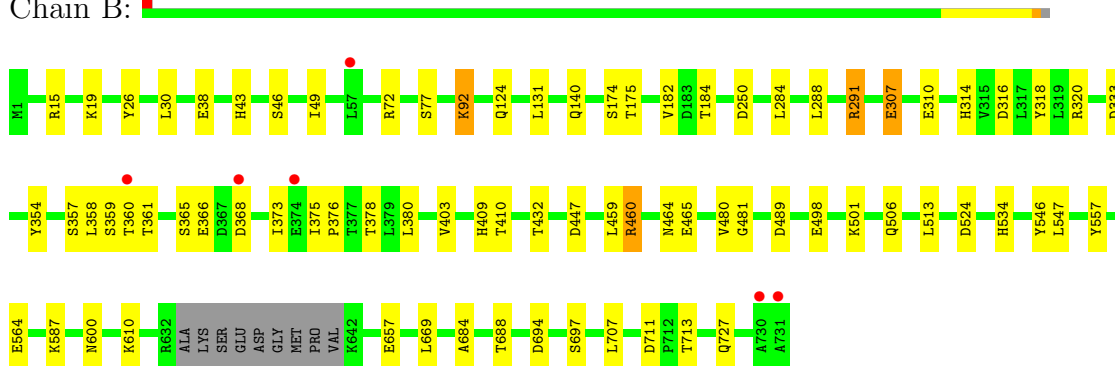
- Molecule 1: N-terminal acetyltransferase A complex subunit nat1

Chain A:



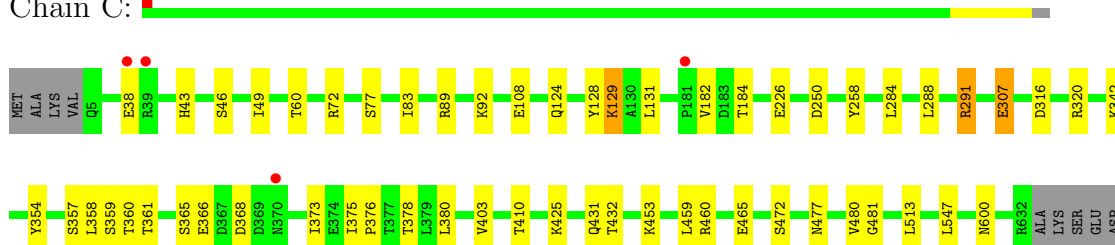
- Molecule 1: N-terminal acetyltransferase A complex subunit nat1

Chain B:



- Molecule 1: N-terminal acetyltransferase A complex subunit nat1

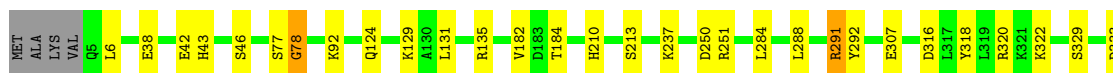
Chain C:





- Molecule 1: N-terminal acetyltransferase A complex subunit nat1

Chain D:



- Molecule 2: N-terminal acetyltransferase A complex catalytic subunit ard1

Chain E:



- Molecule 2: N-terminal acetyltransferase A complex catalytic subunit ard1

Chain F:



- Molecule 2: N-terminal acetyltransferase A complex catalytic subunit ard1

Chain G:



- Molecule 2: N-terminal acetyltransferase A complex catalytic subunit ard1

Chain H:



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	80.74Å 119.69Å 132.02Å 80.28° 76.85° 70.65°	Depositor
Resolution (Å)	49.05 – 3.15 49.05 – 3.15	Depositor EDS
% Data completeness (in resolution range)	98.2 (49.05-3.15) 91.9 (49.05-3.15)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.29 (at 3.12Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1069)	Depositor
R, R_{free}	0.217 , 0.246 0.212 , 0.241	Depositor DCC
R_{free} test set	3608 reflections (4.92%)	DCC
Wilson B-factor (Å ²)	75.0	Xtriage
Anisotropy	0.590	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 28.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 76796 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	28430	wwPDB-VP
Average B, all atoms (Å ²)	85.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.25% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: NA, ACO, SO4, 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	0.23	0/5878	0.41	0/7929
1	B	0.22	0/5891	0.41	0/7948
1	C	0.22	0/5867	0.41	0/7915
1	D	0.23	0/5876	0.42	0/7927
2	E	0.24	0/1277	0.52	4/1722 (0.2%)
2	F	0.22	0/1277	0.41	0/1722
2	G	0.22	0/1277	0.40	0/1722
2	H	0.25	0/1277	0.50	3/1722 (0.2%)
All	All	0.23	0/28620	0.42	7/38607 (0.0%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	23	PRO	CA-C-N	6.96	132.52	117.20
2	E	23	PRO	C-N-CA	6.10	136.95	121.70
2	H	23	PRO	CA-C-N	6.10	130.62	117.20
2	H	23	PRO	N-CA-C	5.79	127.16	112.10
2	E	23	PRO	N-CA-C	5.69	126.90	112.10
2	E	23	PRO	CA-C-O	-5.49	107.03	120.20
2	H	23	PRO	C-N-CA	5.30	134.95	121.70

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens

added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5764	0	0	30	0
1	B	5777	0	13	24	0
1	C	5753	0	0	17	0
1	D	5761	0	0	24	1
2	E	1247	0	0	8	0
2	F	1247	0	0	8	0
2	G	1247	0	0	5	0
2	H	1247	0	0	6	0
3	A	1	0	0	0	0
3	D	1	0	0	0	0
3	E	2	0	0	0	0
3	G	1	0	0	0	0
4	A	10	0	0	4	0
4	B	14	0	0	1	0
4	C	15	0	0	4	0
4	D	10	0	0	4	0
4	E	7	0	0	1	0
4	F	2	0	0	0	0
4	G	2	0	0	0	0
4	H	3	0	0	0	0
5	A	5	0	0	0	0
5	D	5	0	0	0	0
6	E	51	0	34	0	0
6	F	51	0	34	1	0
6	G	51	0	34	0	0
6	H	51	0	34	0	1
7	A	22	0	0	3	0
7	B	22	0	0	4	0
7	C	25	0	0	1	0
7	D	11	0	0	0	0
7	E	4	0	0	0	0
7	F	6	0	0	0	0
7	G	5	0	0	0	0
7	H	10	0	0	1	0
All	All	28430	0	149	113	1

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 4.

All (113) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:89:ARG:NH1	4:C:801:CL:CL	2.25	1.06
1:A:320:ARG:NH1	1:A:368:ASP:OD2	2.02	0.93
1:C:342:LYS:NZ	4:C:815:CL:CL	2.39	0.92
1:D:320:ARG:NH1	1:D:368:ASP:OD2	2.07	0.88
2:H:80:ARG:NH1	7:H:310:HOH:O	2.08	0.87
1:C:320:ARG:NH1	1:C:368:ASP:OD2	2.10	0.83
1:B:320:ARG:NH1	1:B:368:ASP:OD2	2.13	0.81
1:D:250:ASP:OD2	1:D:291:ARG:NH1	2.16	0.78
1:A:250:ASP:OD2	1:A:291:ARG:NH1	2.19	0.76
1:A:135:ARG:NH1	4:A:802:CL:CL	2.56	0.76
1:D:135:ARG:NH1	4:D:803:CL:CL	2.58	0.74
1:A:422:ARG:NH2	4:A:809:CL:CL	2.61	0.70
2:E:17:CYS:O	2:E:21:ASN:ND2	2.23	0.70
1:A:308:GLY:N	4:A:808:CL:CL	2.62	0.69
1:B:175:THR:OG1	7:B:902:HOH:O	2.11	0.68
1:B:447:ASP:OD1	7:B:907:HOH:O	2.14	0.65
1:A:62:GLU:O	1:A:66:ASN:ND2	2.30	0.65
2:G:17:CYS:O	2:G:21:ASN:ND2	2.30	0.64
1:A:43:HIS:ND1	1:A:46:SER:OG	2.32	0.63
1:A:342:LYS:NZ	4:A:808:CL:CL	2.69	0.63
1:C:425:LYS:NZ	4:C:809:CL:CL	2.69	0.63
1:D:329:SER:OG	2:H:81:SER:O	2.17	0.63
1:C:250:ASP:OD2	1:C:291:ARG:NH1	2.32	0.62
2:H:21:ASN:OD1	2:H:21:ASN:N	2.35	0.60
2:F:21:ASN:OD1	2:F:21:ASN:N	2.36	0.59
2:E:55:TYR:OH	2:E:75:SER:OG	2.21	0.59
1:D:251:ARG:NH1	2:H:97:ARG:NH1	2.51	0.59
1:B:489:ASP:O	2:F:27:GLN:NE2	2.36	0.58
1:B:250:ASP:OD2	1:B:291:ARG:NH1	2.37	0.58
1:C:43:HIS:ND1	1:C:46:SER:OG	2.37	0.57
1:A:251:ARG:NH1	2:E:97:ARG:NH1	2.52	0.57
2:F:55:TYR:OH	2:F:75:SER:OG	2.22	0.56
1:D:587:LYS:NZ	1:D:610:LYS:NZ	2.53	0.55
1:D:422:ARG:NH2	4:D:807:CL:CL	2.77	0.55
2:H:55:TYR:OH	2:H:75:SER:OG	2.25	0.54
1:D:657:GLU:OE2	1:D:657:GLU:N	2.40	0.54
1:D:77:SER:OG	1:D:78:GLY:N	2.40	0.54
1:B:318:TYR:OH	1:B:333:ASP:OD2	2.26	0.53
1:C:650:GLU:OE1	7:C:915:HOH:O	2.19	0.53
1:B:307:GLU:OE2	1:B:307:GLU:N	2.42	0.53
1:D:316:ASP:OD2	1:D:354:TYR:OH	2.27	0.53
1:B:524:ASP:OD1	1:B:557:TYR:OH	2.28	0.52
1:D:684:ALA:O	1:D:688:THR:OG1	2.28	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:329:SER:OG	2:E:81:SER:O	2.28	0.52
1:A:447:ASP:OD1	7:A:914:HOH:O	2.19	0.51
1:B:140:GLN:NE2	7:B:901:HOH:O	2.45	0.50
1:D:711:ASP:OD2	1:D:713:THR:OG1	2.30	0.50
1:D:357:SER:OG	1:D:363:LYS:O	2.30	0.50
1:A:128:TYR:N	1:A:129[A]:LYS:NZ	2.60	0.50
1:B:92:LYS:NZ	4:B:802:CL:CL	2.82	0.49
1:D:318:TYR:OH	1:D:333:ASP:OD2	2.31	0.49
2:F:21:ASN:ND2	2:F:78:VAL:O	2.45	0.49
1:A:514:LYS:NZ	1:A:645:GLU:OE1	2.45	0.49
1:A:711:ASP:OD2	1:A:713:THR:OG1	2.31	0.48
1:D:514:LYS:NZ	1:D:645:GLU:OE1	2.47	0.48
1:A:391:ASP:OD1	1:A:426:HIS:ND1	2.46	0.48
1:B:684:ALA:O	1:B:688:THR:OG1	2.32	0.48
1:B:534:HIS:ND1	1:B:546:TYR:OH	2.47	0.47
1:B:310:GLU:O	1:B:314:HIS:ND1	2.48	0.47
1:B:19:LYS:NZ	1:D:598:GLU:OE1	2.48	0.46
1:A:684:ALA:O	1:A:688:THR:OG1	2.33	0.46
1:A:587:LYS:NZ	1:A:610:LYS:NZ	2.64	0.46
1:D:365:SER:OG	1:D:366:GLU:N	2.48	0.46
1:A:448:ARG:NH1	2:E:22:LEU:O	2.49	0.46
2:G:132:GLN:N	2:G:147:ALA:O	2.49	0.45
1:D:694:ASP:OD1	1:D:694:ASP:N	2.49	0.45
2:E:4:ARG:NH2	4:E:206:CL:CL	2.86	0.45
1:C:128:TYR:N	1:C:129:LYS:NZ	2.65	0.45
1:A:532:ASP:OD2	2:E:29:LYS:NZ	2.50	0.45
1:A:694:ASP:N	1:A:694:ASP:OD1	2.50	0.45
1:C:226:GLU:OE2	1:C:258:TYR:OH	2.33	0.45
1:D:359:SER:OG	1:D:360:THR:N	2.51	0.44
1:A:476:ARG:NH2	1:A:478:GLU:OE1	2.51	0.44
1:A:357:SER:OG	1:A:363:LYS:O	2.35	0.44
1:B:460:ARG:NH2	1:B:498:GLU:OE1	2.51	0.44
1:C:453:LYS:NZ	4:C:808:CL:CL	2.88	0.44
7:B:907:HOH:O	2:F:83:ARG:NH2	2.51	0.44
1:A:310:GLU:O	1:A:314:HIS:ND1	2.51	0.44
1:B:464:ASN:OD1	1:B:506:GLN:NE2	2.51	0.44
1:C:472:SER:O	1:C:477:ASN:ND2	2.51	0.43
1:B:694:ASP:N	1:B:694:ASP:OD1	2.51	0.43
1:A:339:LYS:NZ	7:A:921:HOH:O	2.51	0.43
1:A:365:SER:OG	1:A:366:GLU:N	2.51	0.43
1:C:359:SER:OG	1:C:360:THR:N	2.52	0.43
1:B:359:SER:OG	1:B:360:THR:N	2.52	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:711:ASP:OD2	1:C:713:THR:OG1	2.36	0.43
1:B:711:ASP:OD2	1:B:713:THR:OG1	2.37	0.43
2:H:1[B]:MET:N	2:H:1[B]:MET:SD	2.92	0.42
1:A:189:GLU:OE1	7:A:902:HOH:O	2.21	0.42
1:C:694:ASP:OD1	1:C:694:ASP:N	2.52	0.42
1:A:270:ASP:OD1	1:A:271:ALA:N	2.52	0.42
1:B:365:SER:OG	1:B:366:GLU:N	2.51	0.42
2:F:4:ARG:NH2	2:F:10:ASP:OD2	2.52	0.42
1:C:365:SER:OG	1:C:366:GLU:N	2.51	0.42
1:D:322:LYS:NZ	1:D:333:ASP:OD2	2.52	0.42
2:G:115:SER:OG	2:G:143:GLU:OE1	2.38	0.42
2:F:86:GLY:N	6:F:201:ACO:O1A	2.53	0.42
1:B:43:HIS:ND1	1:B:46:SER:OG	2.53	0.42
2:F:115:SER:OG	2:F:143:GLU:OE1	2.38	0.42
1:D:43:HIS:ND1	1:D:46:SER:OG	2.53	0.42
1:C:316:ASP:OD2	1:C:354:TYR:OH	2.37	0.42
2:G:66:ASP:N	2:G:66:ASP:OD2	2.53	0.42
1:C:307:GLU:OE2	1:C:307:GLU:O	2.39	0.41
1:D:237:LYS:NZ	4:D:811:CL:CL	2.90	0.41
1:A:318:TYR:OH	1:A:333:ASP:OD2	2.39	0.41
1:A:210:HIS:O	1:A:213:SER:OG	2.38	0.41
1:A:251:ARG:CZ	2:E:97:ARG:NH1	2.84	0.41
1:B:501:LYS:NZ	1:B:564:GLU:OE1	2.54	0.40
1:D:425:LYS:NZ	4:D:808:CL:CL	2.91	0.40
2:G:4:ARG:NH2	2:G:10:ASP:OD2	2.54	0.40
1:D:210:HIS:O	1:D:213:SER:OG	2.39	0.40
1:B:316:ASP:OD2	1:B:354:TYR:OH	2.39	0.40
1:B:587:LYS:NZ	1:B:610:LYS:NZ	2.69	0.40

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	Distance(Å)	Clash(Å)
1:D:42:GLU:OE1	6:H:201:ACO:O2B[1_655]	2.16	0.04

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	715/731 (98%)	686 (96%)	23 (3%)	6 (1%)	27	77
1	B	719/731 (98%)	688 (96%)	26 (4%)	5 (1%)	30	79
1	C	714/731 (98%)	683 (96%)	26 (4%)	5 (1%)	30	79
1	D	714/731 (98%)	684 (96%)	25 (4%)	5 (1%)	30	79
2	E	151/156 (97%)	148 (98%)	2 (1%)	1 (1%)	30	79
2	F	151/156 (97%)	149 (99%)	2 (1%)	0	100	100
2	G	151/156 (97%)	149 (99%)	2 (1%)	0	100	100
2	H	151/156 (97%)	148 (98%)	3 (2%)	0	100	100
All	All	3466/3548 (98%)	3335 (96%)	109 (3%)	22 (1%)	33	82

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	373	ILE
1	A	480	VAL
1	B	373	ILE
1	C	373	ILE
1	D	373	ILE
2	E	24	GLU
1	A	182	VAL
1	A	481	GLY
1	B	182	VAL
1	B	481	GLY
1	C	77	SER
1	C	182	VAL
1	C	481	GLY
1	D	182	VAL
1	B	77	SER
1	D	6	LEU
1	A	6	LEU
1	A	376	PRO
1	D	78	GLY
1	D	376	PRO
1	B	376	PRO
1	C	376	PRO

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	610/638 (96%)	571 (94%)	39 (6%)	25	68
1	B	610/638 (96%)	573 (94%)	37 (6%)	26	71
1	C	609/638 (96%)	572 (94%)	37 (6%)	26	71
1	D	610/638 (96%)	581 (95%)	29 (5%)	35	79
2	E	133/135 (98%)	124 (93%)	9 (7%)	22	65
2	F	133/135 (98%)	123 (92%)	10 (8%)	19	61
2	G	133/135 (98%)	120 (90%)	13 (10%)	12	43
2	H	133/135 (98%)	123 (92%)	10 (8%)	19	61
All	All	2971/3092 (96%)	2787 (94%)	184 (6%)	26	70

All (184) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	ILE
1	A	38	GLU
1	A	53	LEU
1	A	72	ARG
1	A	83	ILE
1	A	92	LYS
1	A	96	GLN
1	A	124	GLN
1	A	129[A]	LYS
1	A	129[B]	LYS
1	A	131	LEU
1	A	184	THR
1	A	284	LEU
1	A	288	LEU
1	A	291	ARG
1	A	292	TYR
1	A	307	GLU
1	A	357	SER
1	A	361	THR
1	A	375	ILE
1	A	378	THR

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Mol	Chain	Res	Type
1	A	380	LEU
1	A	403	VAL
1	A	409	HIS
1	A	410	THR
1	A	432	THR
1	A	459	LEU
1	A	460	ARG
1	A	465	GLU
1	A	480	VAL
1	A	513	LEU
1	A	547	LEU
1	A	600	ASN
1	A	657	GLU
1	A	660	LYS
1	A	669	LEU
1	A	707	LEU
1	A	709	GLN
1	A	727	GLN
1	B	15	ARG
1	B	26	TYR
1	B	30	LEU
1	B	38	GLU
1	B	49	ILE
1	B	72	ARG
1	B	92	LYS
1	B	124	GLN
1	B	131	LEU
1	B	174	SER
1	B	184	THR
1	B	284	LEU
1	B	288	LEU
1	B	291	ARG
1	B	307	GLU
1	B	357	SER
1	B	358	LEU
1	B	361	THR
1	B	375	ILE
1	B	378	THR
1	B	380	LEU
1	B	403	VAL
1	B	409	HIS
1	B	410	THR

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Mol	Chain	Res	Type
1	B	432	THR
1	B	459	LEU
1	B	460	ARG
1	B	465	GLU
1	B	480	VAL
1	B	513	LEU
1	B	547	LEU
1	B	600	ASN
1	B	657	GLU
1	B	669	LEU
1	B	697	SER
1	B	707	LEU
1	B	727	GLN
1	C	38	GLU
1	C	49	ILE
1	C	60	THR
1	C	72	ARG
1	C	83	ILE
1	C	92	LYS
1	C	108	GLU
1	C	124	GLN
1	C	129	LYS
1	C	131	LEU
1	C	184	THR
1	C	284	LEU
1	C	288	LEU
1	C	291	ARG
1	C	307	GLU
1	C	357	SER
1	C	358	LEU
1	C	361	THR
1	C	375	ILE
1	C	378	THR
1	C	380	LEU
1	C	403	VAL
1	C	410	THR
1	C	431	GLN
1	C	432	THR
1	C	459	LEU
1	C	460	ARG
1	C	465	GLU
1	C	480	VAL

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Mol	Chain	Res	Type
1	C	513	LEU
1	C	547	LEU
1	C	600	ASN
1	C	657	GLU
1	C	660	LYS
1	C	669	LEU
1	C	707	LEU
1	C	727	GLN
1	D	38	GLU
1	D	92	LYS
1	D	124	GLN
1	D	129	LYS
1	D	131	LEU
1	D	184	THR
1	D	284	LEU
1	D	288	LEU
1	D	291	ARG
1	D	292	TYR
1	D	307	GLU
1	D	357	SER
1	D	358	LEU
1	D	361	THR
1	D	375	ILE
1	D	378	THR
1	D	380	LEU
1	D	403	VAL
1	D	410	THR
1	D	459	LEU
1	D	460	ARG
1	D	465	GLU
1	D	513	LEU
1	D	547	LEU
1	D	600	ASN
1	D	669	LEU
1	D	707	LEU
1	D	709	GLN
1	D	727	GLN
2	E	3	ILE
2	E	21	ASN
2	E	24	GLU
2	E	25	ASN
2	E	51	ARG

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Mol	Chain	Res	Type
2	E	57	LEU
2	E	81	SER
2	E	106	LYS
2	E	120	ILE
2	F	1[A]	MET
2	F	1[B]	MET
2	F	3	ILE
2	F	21	ASN
2	F	23	PRO
2	F	51	ARG
2	F	57	LEU
2	F	81	SER
2	F	120	ILE
2	F	130	ASP
2	G	1[A]	MET
2	G	1[B]	MET
2	G	3	ILE
2	G	7	ARG
2	G	21	ASN
2	G	25	ASN
2	G	51	ARG
2	G	57	LEU
2	G	66	ASP
2	G	68	ILE
2	G	81	SER
2	G	120	ILE
2	G	130	ASP
2	H	1[A]	MET
2	H	1[B]	MET
2	H	3	ILE
2	H	7	ARG
2	H	21	ASN
2	H	27	GLN
2	H	51	ARG
2	H	57	LEU
2	H	81	SER
2	H	120	ILE

Some 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 chains in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 74 ligands modelled in this entry, 68 are monoatomic - leaving 6 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	SO4	A	812	-	4,4,4	0.20	0	6,6,6	0.07	0
5	SO4	D	812	-	4,4,4	0.19	0	6,6,6	0.07	0
6	ACO	E	201	-	53,53,53	1.86	8 (15%)	79,79,79	1.88	19 (24%)
6	ACO	F	201	-	53,53,53	1.87	8 (15%)	79,79,79	1.91	20 (25%)
6	ACO	G	201	-	53,53,53	1.86	8 (15%)	79,79,79	1.86	19 (24%)
6	ACO	H	201	-	53,53,53	1.86	8 (15%)	79,79,79	1.89	19 (24%)

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
5	SO4	A	812	-	-	0/0/0/0	0/0/0/0
5	SO4	D	812	-	-	0/0/0/0	0/0/0/0
6	ACO	E	201	-	-	0/51/67/67	0/1/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	ACO	F	201	-	-	0/51/67/67	0/1/3/3
6	ACO	G	201	-	-	0/51/67/67	0/1/3/3
6	ACO	H	201	-	-	1/51/67/67	0/1/3/3

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	H	201	ACO	C9P-N8P	7.01	1.48	1.33
6	G	201	ACO	C9P-N8P	7.00	1.48	1.33
6	F	201	ACO	C9P-N8P	6.99	1.48	1.33
6	E	201	ACO	C9P-N8P	6.92	1.48	1.33
6	F	201	ACO	C5P-N4P	5.27	1.45	1.33
6	H	201	ACO	C5P-N4P	5.26	1.45	1.33
6	E	201	ACO	C5P-N4P	5.25	1.45	1.33
6	G	201	ACO	C5P-N4P	5.23	1.45	1.33
6	E	201	ACO	O-C	4.64	1.38	1.21
6	F	201	ACO	O-C	4.61	1.38	1.21
6	G	201	ACO	O-C	4.59	1.38	1.21
6	H	201	ACO	O-C	4.58	1.38	1.21
6	F	201	ACO	C2B-C3B	-3.82	1.44	1.53
6	G	201	ACO	C2B-C3B	-3.73	1.44	1.53
6	E	201	ACO	C2B-C3B	-3.66	1.44	1.53
6	H	201	ACO	C2B-C3B	-3.52	1.44	1.53
6	F	201	ACO	C2B-C1B	-3.27	1.48	1.53
6	H	201	ACO	CBP-CAP	2.97	1.58	1.55
6	G	201	ACO	CBP-CAP	2.95	1.58	1.55
6	G	201	ACO	C2B-C1B	-2.93	1.49	1.53
6	F	201	ACO	CBP-CAP	2.92	1.58	1.55
6	H	201	ACO	C2B-C1B	-2.87	1.49	1.53
6	E	201	ACO	C2B-C1B	-2.86	1.49	1.53
6	E	201	ACO	CBP-CAP	2.82	1.58	1.55
6	E	201	ACO	OAP-CAP	-2.47	1.37	1.42
6	F	201	ACO	OAP-CAP	-2.47	1.37	1.42
6	G	201	ACO	OAP-CAP	-2.39	1.37	1.42
6	H	201	ACO	OAP-CAP	-2.38	1.37	1.42
6	E	201	ACO	C3B-C4B	-2.34	1.45	1.52
6	G	201	ACO	C3B-C4B	-2.30	1.46	1.52
6	F	201	ACO	C3B-C4B	-2.27	1.46	1.52
6	H	201	ACO	C3B-C4B	-2.27	1.46	1.52

All (77) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	F	201	ACO	N3A-C2A-N1A	-8.13	121.91	128.71
6	G	201	ACO	N3A-C2A-N1A	-8.09	121.94	128.71
6	E	201	ACO	N3A-C2A-N1A	-7.95	122.06	128.71
6	H	201	ACO	N3A-C2A-N1A	-7.94	122.07	128.71
6	E	201	ACO	CBP-CAP-C9P	4.86	117.45	112.73
6	F	201	ACO	CBP-CAP-C9P	4.83	117.41	112.73
6	H	201	ACO	CBP-CAP-C9P	4.62	117.21	112.73
6	E	201	ACO	N3A-C4A-N9A	4.60	133.75	125.43
6	H	201	ACO	N3A-C4A-N9A	4.60	133.74	125.43
6	G	201	ACO	N3A-C4A-N9A	4.57	133.68	125.43
6	F	201	ACO	N3A-C4A-N9A	4.54	133.63	125.43
6	H	201	ACO	O6A-CCP-CBP	4.17	117.55	110.57
6	G	201	ACO	CBP-CAP-C9P	4.07	116.68	112.73
6	G	201	ACO	O6A-CCP-CBP	4.07	117.38	110.57
6	F	201	ACO	P2A-O3A-P1A	-4.06	119.79	131.68
6	H	201	ACO	O-C-S1P	-4.04	107.50	122.06
6	F	201	ACO	O6A-CCP-CBP	4.03	117.31	110.57
6	G	201	ACO	O-C-S1P	-3.98	107.73	122.06
6	E	201	ACO	O6A-CCP-CBP	3.92	117.13	110.57
6	F	201	ACO	O-C-S1P	-3.84	108.22	122.06
6	E	201	ACO	O-C-S1P	-3.83	108.26	122.06
6	E	201	ACO	O-C-CH3	-3.71	110.48	123.20
6	G	201	ACO	O-C-CH3	-3.68	110.56	123.20
6	G	201	ACO	P2A-O3A-P1A	-3.61	121.08	131.68
6	F	201	ACO	O-C-CH3	-3.61	110.81	123.20
6	H	201	ACO	O-C-CH3	-3.61	110.83	123.20
6	H	201	ACO	P2A-O3A-P1A	-3.59	121.16	131.68
6	G	201	ACO	C7P-C6P-C5P	3.53	118.27	112.25
6	F	201	ACO	O4B-C1B-N9A	3.47	111.67	108.44
6	E	201	ACO	C7P-C6P-C5P	3.41	118.06	112.25
6	E	201	ACO	P2A-O3A-P1A	-3.28	122.06	131.68
6	H	201	ACO	C3B-C2B-C1B	3.04	107.19	99.98
6	E	201	ACO	C3B-C2B-C1B	3.01	107.13	99.98
6	F	201	ACO	C7P-C6P-C5P	2.94	117.26	112.25
6	H	201	ACO	C7P-C6P-C5P	2.93	117.24	112.25
6	E	201	ACO	C5A-C4A-N3A	-2.83	119.53	125.70
6	H	201	ACO	C5A-C4A-N3A	-2.80	119.61	125.70
6	G	201	ACO	C5A-C4A-N3A	-2.78	119.66	125.70
6	F	201	ACO	C5A-C4A-N3A	-2.77	119.67	125.70
6	G	201	ACO	O4B-C1B-N9A	2.72	110.97	108.44
6	H	201	ACO	C4B-O4B-C1B	-2.60	106.93	109.75
6	G	201	ACO	O5B-C5B-C4B	2.55	118.30	108.94
6	F	201	ACO	C2P-C3P-N4P	2.55	118.27	112.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	E	201	ACO	O4B-C1B-N9A	2.48	110.75	108.44
6	E	201	ACO	C2P-C3P-N4P	2.48	118.12	112.50
6	F	201	ACO	C4B-O4B-C1B	-2.46	107.08	109.75
6	H	201	ACO	C2P-C3P-N4P	2.44	118.01	112.50
6	F	201	ACO	O5B-C5B-C4B	2.42	117.83	108.94
6	G	201	ACO	C2P-C3P-N4P	2.36	117.85	112.50
6	G	201	ACO	C3B-C2B-C1B	2.35	105.56	99.98
6	E	201	ACO	O5B-C5B-C4B	2.34	117.54	108.94
6	F	201	ACO	C2P-S1P-C	2.33	110.86	101.38
6	G	201	ACO	C3P-C2P-S1P	2.33	118.21	111.04
6	E	201	ACO	C2P-S1P-C	2.27	110.62	101.38
6	E	201	ACO	C3P-C2P-S1P	2.27	118.03	111.04
6	F	201	ACO	C4A-C5A-N7A	-2.27	107.58	109.52
6	F	201	ACO	C3P-C2P-S1P	2.26	118.01	111.04
6	F	201	ACO	C2A-N3A-C4A	2.24	120.40	114.01
6	H	201	ACO	C3P-C2P-S1P	2.24	117.93	111.04
6	E	201	ACO	C2A-N3A-C4A	2.23	120.35	114.01
6	G	201	ACO	C2A-N3A-C4A	2.22	120.34	114.01
6	E	201	ACO	C4A-C5A-N7A	-2.21	107.63	109.52
6	G	201	ACO	C4A-C5A-N7A	-2.21	107.63	109.52
6	H	201	ACO	C2A-N3A-C4A	2.20	120.28	114.01
6	H	201	ACO	C2P-S1P-C	2.20	110.32	101.38
6	H	201	ACO	C4A-C5A-N7A	-2.19	107.64	109.52
6	G	201	ACO	C2P-S1P-C	2.18	110.26	101.38
6	F	201	ACO	C6P-C7P-N8P	2.17	116.54	111.87
6	H	201	ACO	O5B-C5B-C4B	2.15	116.81	108.94
6	E	201	ACO	C4B-O4B-C1B	-2.14	107.42	109.75
6	F	201	ACO	C8A-N9A-C4A	2.11	108.51	106.90
6	H	201	ACO	C6P-C5P-N4P	2.10	120.32	116.50
6	E	201	ACO	C6P-C5P-N4P	2.10	120.31	116.50
6	G	201	ACO	C6P-C7P-N8P	2.10	116.39	111.87
6	H	201	ACO	C8A-N9A-C4A	2.04	108.45	106.90
6	F	201	ACO	C6P-C5P-N4P	2.02	120.17	116.50
6	G	201	ACO	C8A-N9A-C4A	2.01	108.44	106.90

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	H	201	ACO	CH3-C-S1P-C2P

There are no ring outliers.

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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	718/731 (98%)	-0.17	3 (0%)	90 49	38, 69, 149, 217	0
1	B	722/731 (98%)	-0.06	6 (0%)	83 32	37, 90, 170, 249	0
1	C	717/731 (98%)	-0.08	7 (0%)	79 26	30, 84, 160, 213	0
1	D	717/731 (98%)	-0.13	2 (0%)	91 56	40, 73, 158, 203	0
2	E	153/156 (98%)	-0.20	0	100 100	38, 57, 101, 149	0
2	F	153/156 (98%)	-0.08	2 (1%)	74 21	36, 80, 146, 169	0
2	G	153/156 (98%)	0.02	1 (0%)	84 35	35, 86, 156, 185	0
2	H	153/156 (98%)	-0.11	0	100 100	36, 60, 135, 172	0
All	All	3486/3548 (98%)	-0.11	21 (0%)	84 39	30, 76, 159, 249	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	730	ALA	6.1
1	B	731	ALA	6.1
1	B	730	ALA	5.2
1	D	729	ASN	5.2
1	D	374	GLU	3.3
1	B	360	THR	3.1
1	B	374	GLU	2.8
1	C	370	ASN	2.7
1	A	729	ASN	2.7
1	B	57	LEU	2.7
1	B	368	ASP	2.6
1	C	181	PRO	2.4
1	C	707	LEU	2.3
2	G	135	GLU	2.3
1	A	730	ALA	2.3
1	C	644	ASP	2.3

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Mol	Chain	Res	Type	RSRZ
2	F	146	TYR	2.3
1	C	39	ARG	2.3
2	F	66	ASP	2.1
1	A	39	ARG	2.1
1	C	38	GLU	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	CL	C	803	1/1	1.30	41.10	111,111,111,111	0
4	CL	E	207	1/1	0.53	40.73	100,100,100,100	0
4	CL	A	802	1/1	0.68	23.53	108,108,108,108	0
4	CL	C	801	1/1	1.01	17.98	117,117,117,117	0
4	CL	C	809	1/1	0.43	16.41	89,89,89,89	0
4	CL	C	802	1/1	0.29	13.46	77,77,77,77	0
4	CL	B	810	1/1	0.53	11.89	98,98,98,98	0
4	CL	F	203	1/1	0.43	9.57	80,80,80,80	0
4	CL	D	803	1/1	0.45	9.15	94,94,94,94	0
4	CL	B	803	1/1	0.30	8.38	83,83,83,83	0
4	CL	D	805	1/1	0.30	7.98	87,87,87,87	0
4	CL	A	810	1/1	0.34	7.73	72,72,72,72	0
4	CL	A	809	1/1	0.30	6.41	55,55,55,55	0
3	NA	E	203	1/1	0.44	5.53	39,39,39,39	0
3	NA	A	801	1/1	0.29	5.01	38,38,38,38	0
4	CL	D	808	1/1	0.34	4.25	97,97,97,97	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	CL	B	811	1/1	0.30	4.10	81,81,81,81	0
3	NA	E	202	1/1	0.33	4.02	72,72,72,72	0
4	CL	B	809	1/1	0.32	3.88	78,78,78,78	0
4	CL	E	206	1/1	0.31	3.88	91,91,91,91	0
4	CL	E	204	1/1	0.27	3.70	47,47,47,47	0
4	CL	H	202	1/1	0.27	3.19	44,44,44,44	0
4	CL	C	810	1/1	0.37	2.63	83,83,83,83	0
4	CL	C	806	1/1	0.24	2.56	77,77,77,77	0
4	CL	A	808	1/1	0.28	2.13	85,85,85,85	0
4	CL	A	807	1/1	0.27	2.12	104,104,104,104	0
4	CL	G	203	1/1	0.43	1.47	77,77,77,77	0
4	CL	C	804	1/1	0.22	1.28	82,82,82,82	0
6	ACO	F	201	51/51	0.32	1.16	89,154,252,337	0
6	ACO	G	201	51/51	0.32	1.13	117,151,232,233	0
4	CL	C	808	1/1	0.25	1.08	57,57,57,57	0
3	NA	D	801	1/1	0.23	1.07	60,60,60,60	0
4	CL	C	815	1/1	0.28	0.98	133,133,133,133	0
6	ACO	H	201	51/51	0.32	0.97	77,128,209,268	0
4	CL	G	204	1/1	0.24	0.93	72,72,72,72	0
4	CL	E	208	1/1	0.25	0.78	65,65,65,65	0
4	CL	B	807	1/1	0.24	0.69	64,64,64,64	0
4	CL	D	806	1/1	0.21	0.38	99,99,99,99	0
4	CL	D	809	1/1	0.33	0.23	88,88,88,88	0
4	CL	B	808	1/1	0.22	0.20	59,59,59,59	0
4	CL	C	814	1/1	0.20	0.07	77,77,77,77	0
4	CL	F	202	1/1	0.28	-0.09	43,43,43,43	0
6	ACO	E	201	51/51	0.22	-0.09	30,100,183,184	0
4	CL	E	209	1/1	0.18	-0.13	91,91,91,91	0
4	CL	B	806	1/1	0.24	-0.17	91,91,91,91	0
4	CL	A	805	1/1	0.26	-0.20	122,122,122,122	0
4	CL	B	801	1/1	0.18	-0.21	65,65,65,65	0
4	CL	B	802	1/1	0.18	-0.28	87,87,87,87	0
4	CL	E	210	1/1	0.21	-0.37	73,73,73,73	0
4	CL	H	204	1/1	0.21	-0.40	59,59,59,59	0
4	CL	A	806	1/1	0.17	-0.53	77,77,77,77	0
4	CL	E	205	1/1	0.15	-0.57	76,76,76,76	0
4	CL	D	810	1/1	0.20	-0.60	61,61,61,61	0
4	CL	H	203	1/1	0.20	-0.81	121,121,121,121	0
4	CL	D	804	1/1	0.13	-1.02	100,100,100,100	0
4	CL	D	807	1/1	0.14	-1.57	47,47,47,47	0
4	CL	C	813	1/1	0.12	-1.67	86,86,86,86	0
4	CL	A	804	1/1	0.15	-1.69	62,62,62,62	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	SO4	D	812	5/5	0.12	-2.01	66,66,68,69	0
4	CL	A	803	1/1	0.15	-2.05	59,59,59,59	0
4	CL	C	805	1/1	0.14	-2.24	62,62,62,62	0
4	CL	B	814	1/1	0.09	-2.40	76,76,76,76	0
4	CL	D	811	1/1	0.14	-2.56	78,78,78,78	0
4	CL	C	811	1/1	0.16	-2.85	43,43,43,43	0
5	SO4	A	812	5/5	0.12	-2.88	61,62,63,65	0
4	CL	B	804	1/1	0.14	-3.12	92,92,92,92	0
4	CL	D	802	1/1	0.09	-3.37	70,70,70,70	0
3	NA	G	202	1/1	0.09	-3.93	65,65,65,65	0
4	CL	B	805	1/1	0.16	-4.12	73,73,73,73	0
4	CL	C	807	1/1	0.14	-4.32	61,61,61,61	0
4	CL	A	811	1/1	0.09	-5.68	74,74,74,74	0
4	CL	B	812	1/1	0.15	-6.43	103,103,103,103	0
4	CL	B	813	1/1	0.14	-6.74	80,80,80,80	0
4	CL	C	812	1/1	0.21	-	78,78,78,78	0

6.5 Other polymers ⓘ

There are no such residues in this entry.