



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 08:56 PM GMT

PDB ID : 2Q4Y
Title : Ensemble refinement of the protein crystal structure of At1g77540-coenzyme A complex
Authors : Levin, E.J.; Kondrashov, D.A.; Wesenberg, G.E.; Phillips Jr., G.N.; Center for Eukaryotic Structural Genomics (CESG)
Deposited on : 2007-05-31
Resolution : 2.06 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>

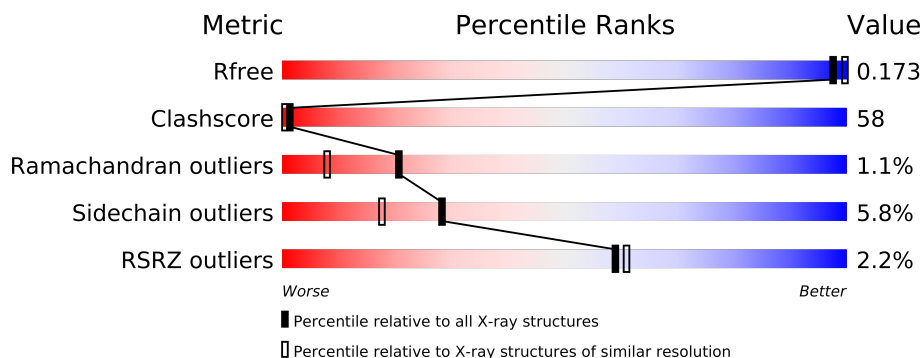
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.06 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	1224 (2.08-2.04)
Clashscore	79885	1390 (2.08-2.04)
Ramachandran outliers	78287	1381 (2.08-2.04)
Sidechain outliers	78261	1381 (2.08-2.04)
RSRZ outliers	66119	1225 (2.08-2.04)

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	1-A	103	
1	2-A	103	
1	3-A	103	
1	4-A	103	

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

Mol	Type	Chain	Res	Geometry	Electron density
2	COA	1-A	104	-	X
2	COA	2-A	104	-	X
2	COA	3-A	104	-	X
2	COA	4-A	104	-	X

2 Entry composition i

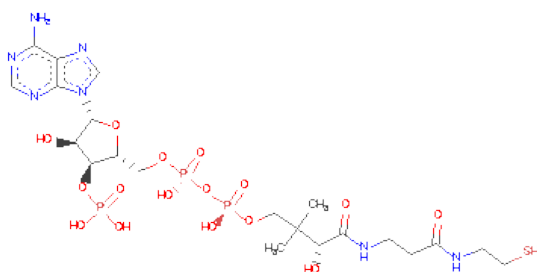
There are 3 unique types of molecules in this entry. The entry contains 3404 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Uncharacterized protein At1g77540.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	1-A	91	Total	C	N	O	S	0	0	0
			735	472	130	129	4			
1	2-A	91	Total	C	N	O	S	0	0	0
			735	472	130	129	4			
1	3-A	91	Total	C	N	O	S	0	0	0
			735	472	130	129	4			
1	4-A	91	Total	C	N	O	S	0	0	0
			735	472	130	129	4			

- Molecule 2 is COENZYME A (three-letter code: COA) (formula: $C_{21}H_{36}N_7O_{16}P_3S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	1-A	1	Total	C	N	O	P	S	0	0
			48	21	7	16	3	1		
2	2-A	1	Total	C	N	O	P	S	0	0
			48	21	7	16	3	1		

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Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	3-A	1	Total	C	N	O	P	S	0	0
			48	21	7	16	3	1		
2	4-A	1	Total	C	N	O	P	S	0	0
			48	21	7	16	3	1		

- Molecule 3 is water.

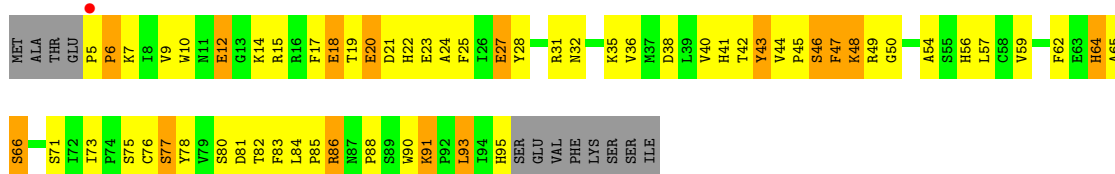
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	1-A	68	Total	O	0	0
			68	68		
3	2-A	68	Total	O	0	0
			68	68		
3	3-A	68	Total	O	0	0
			68	68		
3	4-A	68	Total	O	0	0
			68	68		

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.

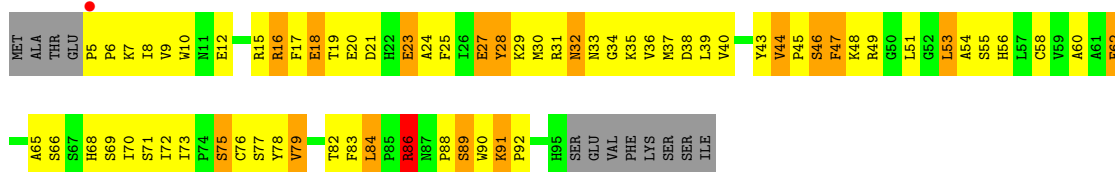
- Molecule 1: Uncharacterized protein At1g77540

Chain 1-A: 



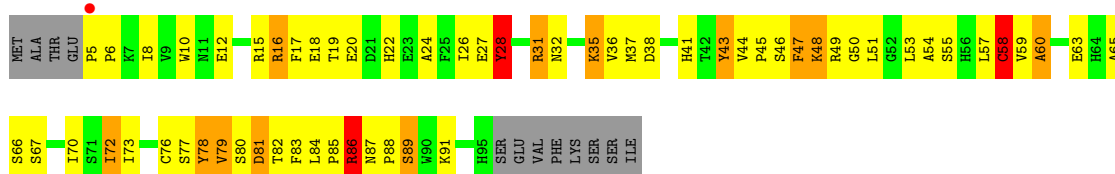
- Molecule 1: Uncharacterized protein At1g77540

Chain 2-A: 



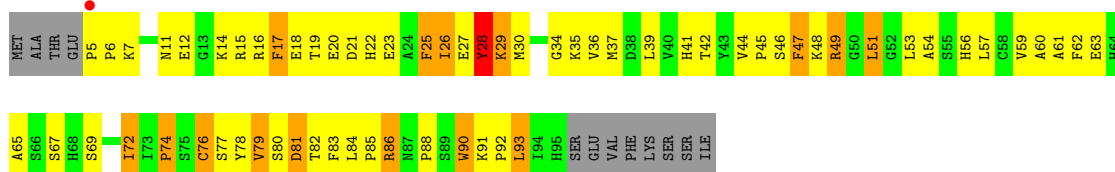
- Molecule 1: Uncharacterized protein At1g77540

Chain 3-A: 



- Molecule 1: Uncharacterized protein At1g77540

Chain 4-A: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	27.91Å 63.94Å 29.52Å 90.00° 90.86° 90.00°	Depositor
Resolution (Å)	31.97 – 2.06 31.97 – 2.05	Depositor EDS
% Data completeness (in resolution range)	99.2 (31.97-2.06) 99.0 (31.97-2.05)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.73 (at 2.05Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.171 , 0.225 0.179 , 0.173	Depositor DCC
R_{free} test set	636 reflections (10.92%)	DCC
Wilson B-factor (Å ²)	29.0	Xtriage
Anisotropy	0.036	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 51.2	EDS
Estimated twinning fraction	0.066 for h,-k,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 6458 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3404	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.36% 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: COA

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	1-A	1.89	12/759 (1.6%)	1.44	4/1028 (0.4%)
1	2-A	1.95	18/759 (2.4%)	1.43	3/1028 (0.3%)
1	3-A	1.89	15/759 (2.0%)	1.38	7/1028 (0.7%)
1	4-A	1.91	19/759 (2.5%)	1.53	11/1028 (1.1%)
All	All	1.91	64/3036 (2.1%)	1.45	25/4112 (0.6%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	1-A	0	1
1	3-A	0	2
1	4-A	0	1
All	All	0	4

All (64) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	3-A	58	CYS	CB-SG	-10.24	1.64	1.82
1	2-A	28	TYR	CD1-CE1	10.22	1.54	1.39
1	4-A	76	CYS	CB-SG	-9.52	1.66	1.82
1	2-A	77	SER	CB-OG	-8.87	1.30	1.42
1	2-A	79	VAL	CB-CG2	8.76	1.71	1.52
1	1-A	12	GLU	CB-CG	-8.56	1.35	1.52
1	2-A	83	PHE	CD2-CE2	8.46	1.56	1.39
1	4-A	79	VAL	CB-CG2	8.28	1.70	1.52
1	2-A	16	ARG	CG-CD	8.20	1.72	1.51
1	3-A	78	TYR	CD1-CE1	8.11	1.51	1.39
1	4-A	25	PHE	CB-CG	7.64	1.64	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	1-A	43	TYR	CB-CG	7.63	1.63	1.51
1	1-A	43	TYR	CD2-CE2	7.57	1.50	1.39
1	3-A	35	LYS	CD-CE	7.35	1.69	1.51
1	3-A	55	SER	CB-OG	-7.31	1.32	1.42
1	4-A	29	LYS	CE-NZ	7.22	1.67	1.49
1	2-A	17	PHE	CE1-CZ	7.21	1.51	1.37
1	1-A	23	GLU	CD-OE1	-7.04	1.18	1.25
1	1-A	27	GLU	CB-CG	6.91	1.65	1.52
1	2-A	23	GLU	CG-CD	6.80	1.62	1.51
1	2-A	27	GLU	CG-CD	6.60	1.61	1.51
1	4-A	36	VAL	CB-CG1	6.45	1.66	1.52
1	1-A	9	VAL	CB-CG1	6.31	1.66	1.52
1	1-A	62	PHE	CB-CG	6.26	1.61	1.51
1	1-A	27	GLU	CG-CD	6.21	1.61	1.51
1	3-A	79	VAL	CB-CG2	6.18	1.65	1.52
1	2-A	32	ASN	CB-CG	6.15	1.65	1.51
1	4-A	60	ALA	CA-CB	6.09	1.65	1.52
1	1-A	18	GLU	CB-CG	6.07	1.63	1.52
1	4-A	74	PRO	CB-CG	5.89	1.79	1.50
1	3-A	60	ALA	CA-CB	5.89	1.64	1.52
1	3-A	28	TYR	CD1-CE1	5.87	1.48	1.39
1	2-A	62	PHE	CB-CG	5.84	1.61	1.51
1	3-A	16	ARG	CG-CD	5.78	1.66	1.51
1	3-A	48	LYS	CE-NZ	5.76	1.63	1.49
1	2-A	58	CYS	CB-SG	-5.73	1.72	1.81
1	4-A	49	ARG	CZ-NH2	5.73	1.40	1.33
1	2-A	43	TYR	CD1-CE1	-5.73	1.30	1.39
1	3-A	43	TYR	CD1-CE1	-5.71	1.30	1.39
1	4-A	56	HIS	C-O	-5.70	1.12	1.23
1	4-A	44	VAL	CB-CG1	5.67	1.64	1.52
1	4-A	62	PHE	CB-CG	5.62	1.60	1.51
1	3-A	28	TYR	CD2-CE2	5.60	1.47	1.39
1	2-A	28	TYR	CD2-CE2	5.46	1.47	1.39
1	4-A	81	ASP	CB-CG	5.45	1.63	1.51
1	3-A	59	VAL	CA-CB	5.44	1.66	1.54
1	2-A	66	SER	CB-OG	-5.37	1.35	1.42
1	4-A	25	PHE	CD1-CE1	5.36	1.50	1.39
1	4-A	28	TYR	CE2-CZ	5.34	1.45	1.38
1	4-A	35	LYS	CD-CE	5.28	1.64	1.51
1	3-A	18	GLU	CB-CG	5.28	1.62	1.52
1	4-A	18	GLU	CB-CG	5.27	1.62	1.52
1	4-A	27	GLU	CD-OE1	-5.24	1.19	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	3-A	63	GLU	CD-OE2	5.22	1.31	1.25
1	3-A	43	TYR	CD2-CE2	5.17	1.47	1.39
1	2-A	44	VAL	CB-CG1	5.16	1.63	1.52
1	4-A	17	PHE	CD2-CE2	5.11	1.49	1.39
1	1-A	83	PHE	CB-CG	-5.11	1.42	1.51
1	1-A	91	LYS	CD-CE	5.08	1.64	1.51
1	2-A	60	ALA	CA-CB	5.07	1.63	1.52
1	4-A	90	TRP	CE3-CZ3	5.07	1.47	1.38
1	1-A	66	SER	CB-OG	-5.07	1.35	1.42
1	2-A	75	SER	CB-OG	-5.06	1.35	1.42
1	2-A	91	LYS	CD-CE	5.04	1.63	1.51

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	4-A	86	ARG	NE-CZ-NH1	-16.34	112.13	120.30
1	2-A	86	ARG	NE-CZ-NH2	-9.51	115.54	120.30
1	4-A	86	ARG	NE-CZ-NH2	8.76	124.68	120.30
1	1-A	86	ARG	NE-CZ-NH1	-8.17	116.22	120.30
1	4-A	49	ARG	NE-CZ-NH1	-7.75	116.43	120.30
1	1-A	17	PHE	N-CA-C	-7.15	91.69	111.00
1	3-A	31	ARG	NE-CZ-NH2	-6.58	117.01	120.30
1	4-A	49	ARG	NE-CZ-NH2	6.45	123.52	120.30
1	4-A	35	LYS	CD-CE-NZ	6.33	126.26	111.70
1	4-A	29	LYS	CD-CE-NZ	6.20	125.97	111.70
1	3-A	86	ARG	CB-CG-CD	-6.14	95.64	111.60
1	2-A	53	LEU	CB-CG-CD1	-5.89	100.99	111.00
1	4-A	16	ARG	NE-CZ-NH1	-5.86	117.37	120.30
1	4-A	17	PHE	CZ-CE2-CD2	-5.81	113.13	120.10
1	3-A	16	ARG	NE-CZ-NH2	5.70	123.15	120.30
1	4-A	26	ILE	N-CA-C	-5.68	95.66	111.00
1	1-A	27	GLU	CA-CB-CG	5.48	125.45	113.40
1	3-A	17	PHE	N-CA-C	-5.33	96.62	111.00
1	2-A	84	LEU	CA-CB-CG	-5.26	103.20	115.30
1	1-A	48	LYS	CD-CE-NZ	5.24	123.75	111.70
1	4-A	16	ARG	NE-CZ-NH2	5.18	122.89	120.30
1	4-A	51	LEU	CB-CG-CD1	-5.17	102.22	111.00
1	3-A	86	ARG	NE-CZ-NH1	5.16	122.88	120.30
1	3-A	86	ARG	CG-CD-NE	-5.16	100.96	111.80
1	3-A	81	ASP	CB-CG-OD1	5.03	122.82	118.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	1-A	64	HIS	Sidechain
1	3-A	28	TYR	Sidechain
1	3-A	78	TYR	Sidechain
1	4-A	28	TYR	Sidechain

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	1-A	735	0	720	76	1
1	2-A	735	0	720	86	1
1	3-A	735	0	720	95	0
1	4-A	735	0	720	69	0
2	1-A	48	0	32	24	0
2	2-A	48	0	31	13	0
2	3-A	48	0	31	21	0
2	4-A	48	0	31	22	0
3	1-A	68	0	0	11	1
3	2-A	68	0	0	11	1
3	3-A	68	0	0	11	0
3	4-A	68	0	0	22	0
All	All	3404	0	3005	353	2

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 58.

All (353) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:73:ILE:CD1	1:A:73:ILE:CG1	1.78	1.58
1:A:74:PRO:CB	1:A:74:PRO:CG	1.79	1.46
1:A:85:PRO:CG	1:A:85:PRO:CB	1.74	1.44
1:A:46:SER:HA	1:A:49:ARG:CZ	1.64	1.27
2:A:104:COA:H121	2:A:104:COA:HN8	1.06	1.15
1:A:29:LYS:HE3	3:A:162:HOH:O	1.46	1.15
1:A:42:THR:H	2:A:104:COA:H142	1.01	1.13

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:23:GLU:OE2	1:A:45:PRO:HG3	1.48	1.11
2:A:104:COA:H121	2:A:104:COA:N8P	1.61	1.11
1:A:82:THR:HG23	1:A:86:ARG:HH21	1.01	1.08
1:A:24:ALA:CB	1:A:53:LEU:HD21	1.85	1.05
1:A:86:ARG:NH1	2:A:104:COA:O8A	1.93	1.02
1:A:45:PRO:HG2	1:A:48:LYS:HD3	1.41	1.01
1:A:45:PRO:HG2	1:A:48:LYS:HB2	1.41	1.01
1:A:47:PHE:O	1:A:48:LYS:HG3	1.60	0.99
1:A:45:PRO:CG	1:A:48:LYS:HD3	1.95	0.95
1:A:11:ASN:OD1	1:A:14:LYS:N	1.99	0.95
1:A:86:ARG:CZ	2:A:104:COA:O8A	2.14	0.95
1:A:8:ILE:HD13	1:A:57:LEU:HA	1.48	0.94
1:A:82:THR:HG23	1:A:86:ARG:NH2	1.82	0.94
1:A:46:SER:CA	1:A:49:ARG:CZ	2.47	0.92
1:A:42:THR:N	2:A:104:COA:H142	1.85	0.91
1:A:82:THR:CG2	1:A:86:ARG:HH21	1.84	0.90
2:A:104:COA:O2B	2:A:104:COA:P3B	2.30	0.90
2:A:104:COA:N8P	2:A:104:COA:H133	1.88	0.88
1:A:82:THR:O	1:A:86:ARG:HG2	1.74	0.87
2:A:104:COA:H133	2:A:104:COA:HN8	1.36	0.87
1:A:44:VAL:HB	2:A:104:COA:H122	1.55	0.86
1:A:78:TYR:HA	2:A:104:COA:H2A	1.58	0.86
1:A:84:LEU:HB3	1:A:91:LYS:HE2	1.58	0.84
1:A:46:SER:HA	1:A:49:ARG:NE	1.92	0.84
2:A:104:COA:O9A	3:A:136:HOH:O	1.95	0.83
2:A:104:COA:P3B	3:A:144:HOH:O	2.36	0.83
2:A:104:COA:CDP	2:A:104:COA:N8P	2.42	0.83
1:A:21:ASP:OD2	3:A:123:HOH:O	1.96	0.82
1:A:86:ARG:HB2	1:A:86:ARG:CZ	2.09	0.80
2:A:104:COA:CDP	2:A:104:COA:HN8	1.95	0.80
1:A:31:ARG:HG2	1:A:31:ARG:HH11	1.46	0.80
1:A:54:ALA:CB	2:A:104:COA:H143	2.11	0.79
1:A:27:GLU:HG3	1:A:40:VAL:HG23	1.61	0.79
2:A:104:COA:HN8	2:A:104:COA:CCP	1.90	0.79
1:A:8:ILE:HD13	1:A:57:LEU:HD23	1.65	0.78
1:A:27:GLU:HG2	1:A:41:HIS:HB3	1.64	0.78
1:A:49:ARG:HD3	2:A:104:COA:H71	1.66	0.78
1:A:42:THR:HG22	2:A:104:COA:OAP	1.83	0.78
1:A:7:LYS:O	1:A:19:THR:HA	1.83	0.78
1:A:86:ARG:HD2	2:A:104:COA:O8A	1.85	0.77
1:A:24:ALA:HB1	1:A:53:LEU:HD21	1.67	0.77
1:A:77:SER:HB3	2:A:104:COA:H132	1.67	0.76

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:7:LYS:HD3	1:A:20:GLU:OE2	1.86	0.76
1:A:7:LYS:HE2	1:A:9:VAL:CG1	2.16	0.75
1:A:45:PRO:HG2	1:A:48:LYS:CD	2.15	0.75
1:A:7:LYS:N	3:A:149:HOH:O	2.19	0.75
1:A:76:CYS:SG	2:A:104:COA:H143	2.27	0.75
1:A:44:VAL:HG13	1:A:53:LEU:HD23	1.68	0.74
1:A:69:SER:OG	3:A:157:HOH:O	2.03	0.74
1:A:54:ALA:HB1	2:A:104:COA:H143	1.70	0.74
1:A:8:ILE:CD1	1:A:57:LEU:HA	2.18	0.73
1:A:37:MET:SD	1:A:72:ILE:HD11	2.27	0.73
1:A:21:ASP:HB3	3:A:156:HOH:O	1.88	0.73
1:A:5:PRO:HB2	1:A:6:PRO:O	1.86	0.73
1:A:66:SER:HB2	1:A:93:LEU:HD21	1.69	0.73
1:A:21:ASP:HB3	3:A:156:HOH:O	1.88	0.73
1:A:23:GLU:OE2	1:A:48:LYS:NZ	2.21	0.73
1:A:32:ASN:HD22	1:A:35:LYS:HB2	1.52	0.72
1:A:31:ARG:NH1	3:A:113:HOH:O	2.23	0.72
1:A:12:GLU:HG2	3:A:127:HOH:O	1.88	0.72
2:A:104:COA:H61	2:A:104:COA:O5A	1.90	0.72
1:A:86:ARG:CZ	2:A:104:COA:P3B	2.78	0.72
1:A:17:PHE:CE2	1:A:61:ALA:HA	2.25	0.72
1:A:90:TRP:CE3	1:A:93:LEU:HD23	2.24	0.72
1:A:9:VAL:N	1:A:18:GLU:O	2.23	0.72
1:A:37:MET:O	1:A:72:ILE:HA	1.89	0.71
1:A:29:LYS:HE3	3:A:162:HOH:O	1.91	0.71
1:A:27:GLU:CG	1:A:41:HIS:HB3	2.21	0.69
1:A:24:ALA:CB	1:A:53:LEU:CD2	2.67	0.69
1:A:15:ARG:HD2	3:A:106:HOH:O	1.93	0.69
1:A:82:THR:HG21	2:A:104:COA:H4B	1.76	0.68
2:A:104:COA:HN8	2:A:104:COA:H133	1.59	0.68
1:A:24:ALA:HB3	1:A:53:LEU:HD21	1.75	0.68
1:A:20:GLU:OE1	3:A:165:HOH:O	2.12	0.67
1:A:86:ARG:NH2	2:A:104:COA:O3B	2.28	0.67
1:A:38:ASP:HB2	1:A:73:ILE:HB	1.76	0.67
1:A:45:PRO:HG2	1:A:48:LYS:HD3	1.77	0.67
1:A:42:THR:O	2:A:104:COA:OAP	2.13	0.67
1:A:23:GLU:OE2	1:A:48:LYS:HD3	1.96	0.66
1:A:47:PHE:N	1:A:47:PHE:CD1	2.64	0.66
1:A:46:SER:N	1:A:49:ARG:NH2	2.44	0.66
2:A:104:COA:HN8	2:A:104:COA:H131	1.60	0.66
1:A:30:MET:CE	1:A:34:GLY:HA2	2.26	0.66
1:A:46:SER:CA	1:A:49:ARG:NH2	2.59	0.66

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:86:ARG:HG3	1:A:87:ASN:ND2	2.11	0.65
1:A:19:THR:HG23	1:A:24:ALA:HB3	1.79	0.65
1:A:31:ARG:HB2	1:A:36:VAL:CG1	2.28	0.64
1:A:32:ASN:ND2	3:A:147:HOH:O	2.23	0.64
1:A:26:ILE:HG23	1:A:26:ILE:O	1.97	0.64
1:A:15:ARG:CD	3:A:106:HOH:O	2.46	0.64
2:A:104:COA:H133	2:A:104:COA:N8P	2.13	0.63
1:A:56:HIS:HE1	3:A:158:HOH:O	1.81	0.63
1:A:56:HIS:CE1	3:A:158:HOH:O	2.51	0.63
1:A:8:ILE:HD12	1:A:60:ALA:HB2	1.80	0.63
1:A:5:PRO:N	1:A:6:PRO:HA	2.14	0.62
1:A:31:ARG:HD2	1:A:73:ILE:HD12	1.80	0.62
1:A:29:LYS:HG3	3:A:161:HOH:O	1.99	0.62
1:A:24:ALA:HA	1:A:43:TYR:O	2.00	0.62
1:A:10:TRP:CE2	1:A:12:GLU:HG2	2.34	0.62
1:A:47:PHE:H	1:A:47:PHE:HD1	1.47	0.61
1:A:44:VAL:HB	2:A:104:COA:CCP	2.28	0.61
1:A:45:PRO:HG3	1:A:48:LYS:HD3	1.82	0.61
2:A:104:COA:O1A	2:A:104:COA:O5A	2.15	0.61
1:A:31:ARG:CG	1:A:31:ARG:HH11	2.14	0.61
2:A:104:COA:N8P	2:A:104:COA:CCP	2.46	0.61
1:A:80:SER:HA	1:A:84:LEU:HD12	1.82	0.61
1:A:23:GLU:OE2	1:A:45:PRO:CG	2.39	0.61
1:A:46:SER:HA	1:A:49:ARG:HD2	1.83	0.60
1:A:46:SER:O	1:A:48:LYS:N	2.34	0.60
1:A:19:THR:HG23	1:A:24:ALA:HB3	1.84	0.60
1:A:44:VAL:HB	2:A:104:COA:H122	1.83	0.59
1:A:44:VAL:CG2	2:A:104:COA:H122	2.32	0.59
1:A:84:LEU:HB2	1:A:85:PRO:HD3	1.83	0.59
1:A:90:TRP:CZ3	1:A:93:LEU:HD23	2.37	0.59
1:A:29:LYS:CE	3:A:162:HOH:O	2.23	0.59
1:A:25:PHE:CD1	1:A:27:GLU:HG3	2.37	0.59
1:A:45:PRO:HB2	1:A:47:PHE:CE1	2.36	0.59
1:A:59:VAL:O	1:A:63:GLU:HG2	2.03	0.59
1:A:81:ASP:O	1:A:85:PRO:HG2	2.02	0.58
1:A:84:LEU:HB2	1:A:85:PRO:HD3	1.85	0.58
1:A:45:PRO:HG2	1:A:48:LYS:HB2	1.86	0.58
1:A:7:LYS:HE2	1:A:9:VAL:HG12	1.85	0.58
1:A:77:SER:HB3	2:A:104:COA:CDP	2.32	0.58
1:A:54:ALA:CB	2:A:104:COA:CEP	2.82	0.58
1:A:41:HIS:HE1	3:A:163:HOH:O	1.86	0.58
1:A:54:ALA:HB3	1:A:78:TYR:CE1	2.39	0.58

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:29:LYS:NZ	3:A:162:HOH:O	2.37	0.58
1:A:14:LYS:O	1:A:15:ARG:HB2	2.03	0.58
1:A:54:ALA:HB2	2:A:104:COA:H143	1.85	0.57
1:A:5:PRO:HG2	3:A:149:HOH:O	2.04	0.57
1:A:42:THR:HG21	2:A:104:COA:H142	1.86	0.57
2:A:104:COA:O2B	3:A:144:HOH:O	2.17	0.57
1:A:29:LYS:CE	3:A:162:HOH:O	2.47	0.57
1:A:89:SER:HB3	3:A:109:HOH:O	2.04	0.57
1:A:45:PRO:HB2	3:A:124:HOH:O	2.03	0.57
1:A:39:LEU:HD13	1:A:79:VAL:HG21	1.85	0.57
1:A:86:ARG:NE	2:A:104:COA:O9A	2.38	0.56
1:A:49:ARG:O	1:A:51:LEU:HG	2.05	0.56
1:A:54:ALA:HB2	2:A:104:COA:H143	1.86	0.56
1:A:15:ARG:NH1	1:A:29:LYS:HG2	2.21	0.56
1:A:25:PHE:CE1	1:A:41:HIS:CD2	2.93	0.56
1:A:19:THR:CG2	1:A:24:ALA:HB3	2.35	0.56
1:A:31:ARG:HB2	1:A:36:VAL:HB	1.87	0.56
1:A:54:ALA:HB2	2:A:104:COA:O6A	2.06	0.56
1:A:71:SER:HB2	1:A:95:HIS:HA	1.87	0.56
1:A:35:LYS:O	1:A:70:ILE:HG23	2.06	0.56
1:A:50:GLY:N	2:A:104:COA:O1A	2.38	0.56
2:A:104:COA:H132	3:A:167:HOH:O	2.06	0.56
1:A:44:VAL:CB	2:A:104:COA:H122	2.32	0.56
1:A:49:ARG:O	1:A:51:LEU:HG	2.06	0.56
1:A:5:PRO:CD	1:A:6:PRO:HA	2.36	0.56
1:A:47:PHE:N	1:A:47:PHE:CD1	2.70	0.55
1:A:45:PRO:CB	3:A:124:HOH:O	2.54	0.55
1:A:21:ASP:OD2	1:A:23:GLU:HB2	2.07	0.55
1:A:10:TRP:HZ2	1:A:28:TYR:HH	1.53	0.55
1:A:32:ASN:ND2	1:A:35:LYS:HD2	2.22	0.54
1:A:64:HIS:CD2	1:A:64:HIS:C	2.80	0.54
1:A:82:THR:O	1:A:86:ARG:HG3	2.07	0.54
1:A:41:HIS:CE1	3:A:163:HOH:O	2.60	0.54
1:A:65:ALA:CB	1:A:72:ILE:HD11	2.37	0.54
1:A:38:ASP:O	1:A:40:VAL:HG13	2.08	0.54
1:A:30:MET:HE2	1:A:34:GLY:HA2	1.90	0.54
1:A:34:GLY:O	3:A:170:HOH:O	2.19	0.54
1:A:25:PHE:CE2	1:A:43:TYR:HB3	2.42	0.54
2:A:104:COA:H8A	3:A:144:HOH:O	2.07	0.54
1:A:31:ARG:C	1:A:36:VAL:HB	2.28	0.54
1:A:75:SER:O	1:A:76:CYS:C	2.46	0.54
1:A:73:ILE:HG22	1:A:75:SER:OG	2.08	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:88:PRO:HA	1:A:91:LYS:HG3	1.89	0.54
1:A:8:ILE:HA	1:A:18:GLU:O	2.08	0.53
1:A:82:THR:CG2	2:A:104:COA:H1B	2.39	0.53
1:A:30:MET:HE3	1:A:34:GLY:HA2	1.90	0.53
2:A:104:COA:H133	2:A:104:COA:N8P	2.13	0.53
1:A:26:ILE:CG2	1:A:26:ILE:O	2.56	0.53
1:A:46:SER:HA	1:A:49:ARG:NH1	2.24	0.53
1:A:42:THR:O	2:A:104:COA:H122	2.09	0.52
1:A:54:ALA:HB2	2:A:104:COA:CEP	2.39	0.52
1:A:6:PRO:HB2	1:A:19:THR:HG22	1.90	0.52
1:A:76:CYS:SG	2:A:104:COA:CEP	2.97	0.52
1:A:6:PRO:HB2	3:A:141:HOH:O	2.08	0.52
1:A:28:TYR:HA	1:A:38:ASP:O	2.09	0.52
1:A:19:THR:HB	3:A:123:HOH:O	2.09	0.52
1:A:15:ARG:HD2	3:A:106:HOH:O	2.09	0.52
1:A:28:TYR:C	1:A:28:TYR:CD1	2.83	0.52
1:A:38:ASP:HA	1:A:73:ILE:O	2.10	0.52
1:A:42:THR:CG2	2:A:104:COA:H142	2.40	0.52
1:A:18:GLU:OE1	1:A:22:HIS:HA	2.09	0.52
1:A:85:PRO:O	1:A:88:PRO:HD3	2.10	0.52
1:A:5:PRO:HD2	1:A:7:LYS:H	1.74	0.52
1:A:31:ARG:O	1:A:36:VAL:HB	2.10	0.52
1:A:5:PRO:HD2	1:A:6:PRO:HA	1.92	0.52
2:A:104:COA:CDP	3:A:167:HOH:O	2.57	0.51
1:A:84:LEU:CB	1:A:91:LYS:HE2	2.35	0.51
1:A:38:ASP:HA	1:A:73:ILE:HB	1.92	0.51
1:A:47:PHE:N	1:A:47:PHE:CD1	2.78	0.51
1:A:49:ARG:NH1	2:A:104:COA:H22	2.26	0.51
1:A:25:PHE:CE1	1:A:27:GLU:HG3	2.45	0.51
1:A:45:PRO:HG2	1:A:48:LYS:HD3	1.93	0.51
1:A:45:PRO:HB2	1:A:47:PHE:CD1	2.46	0.51
1:A:49:ARG:CD	2:A:104:COA:H71	2.38	0.50
1:A:24:ALA:CB	1:A:53:LEU:HD21	2.41	0.50
1:A:23:GLU:CD	1:A:48:LYS:HZ3	2.12	0.50
1:A:65:ALA:HB2	1:A:72:ILE:HD11	1.94	0.50
1:A:7:LYS:HE2	1:A:9:VAL:HG13	1.92	0.50
1:A:46:SER:OG	1:A:49:ARG:NH2	2.45	0.50
1:A:37:MET:HB3	1:A:72:ILE:HG13	1.92	0.50
1:A:16:ARG:HH21	1:A:27:GLU:CD	2.14	0.50
1:A:25:PHE:HE1	1:A:41:HIS:CD2	2.29	0.50
1:A:45:PRO:HB2	3:A:124:HOH:O	2.11	0.50
1:A:7:LYS:HB3	3:A:149:HOH:O	2.12	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:31:ARG:NH1	1:A:31:ARG:CG	2.75	0.50
1:A:49:ARG:HD3	2:A:104:COA:C7P	2.40	0.49
1:A:64:HIS:CD2	1:A:65:ALA:N	2.80	0.49
1:A:37:MET:N	1:A:71:SER:O	2.39	0.49
2:A:104:COA:H51A	3:A:167:HOH:O	2.11	0.49
1:A:54:ALA:O	1:A:58:CYS:HB3	2.12	0.49
1:A:50:GLY:HA2	2:A:104:COA:O1A	2.13	0.49
1:A:90:TRP:HE3	1:A:93:LEU:HD23	1.72	0.49
1:A:49:ARG:NE	2:A:104:COA:H62	2.27	0.49
1:A:5:PRO:HB2	1:A:6:PRO:O	2.12	0.49
1:A:55:SER:HB2	1:A:78:TYR:OH	2.13	0.49
1:A:89:SER:O	1:A:92:PRO:HD2	2.12	0.49
1:A:54:ALA:HB3	1:A:78:TYR:CZ	2.48	0.49
1:A:44:VAL:HB	2:A:104:COA:CCP	2.43	0.49
1:A:59:VAL:O	1:A:59:VAL:HG12	2.13	0.49
1:A:45:PRO:CB	3:A:124:HOH:O	2.61	0.48
1:A:37:MET:SD	1:A:39:LEU:HD21	2.53	0.48
1:A:20:GLU:HB2	3:A:165:HOH:O	2.14	0.48
1:A:82:THR:O	1:A:86:ARG:HG3	2.13	0.48
1:A:45:PRO:CG	1:A:48:LYS:HD3	2.44	0.48
1:A:54:ALA:HB1	2:A:104:COA:CEP	2.39	0.48
1:A:24:ALA:HA	1:A:43:TYR:O	2.14	0.48
1:A:78:TYR:CE2	1:A:83:PHE:HB2	2.48	0.48
1:A:50:GLY:C	1:A:51:LEU:HD23	2.34	0.47
1:A:15:ARG:HD2	3:A:106:HOH:O	2.13	0.47
1:A:20:GLU:HG3	3:A:165:HOH:O	2.14	0.47
1:A:45:PRO:HB2	1:A:47:PHE:CE1	2.50	0.47
1:A:41:HIS:HE1	3:A:163:HOH:O	1.98	0.47
1:A:47:PHE:CG	1:A:47:PHE:O	2.67	0.47
1:A:44:VAL:HG22	1:A:53:LEU:HG	1.97	0.47
1:A:88:PRO:HA	1:A:91:LYS:HD2	1.96	0.47
2:A:104:COA:H131	3:A:112:HOH:O	2.13	0.47
1:A:66:SER:O	1:A:67:SER:C	2.54	0.47
1:A:46:SER:N	1:A:49:ARG:HH21	2.11	0.47
1:A:20:GLU:C	1:A:22:HIS:H	2.18	0.46
1:A:84:LEU:N	1:A:85:PRO:CD	2.78	0.46
1:A:8:ILE:HD11	1:A:56:HIS:HB3	1.97	0.46
1:A:78:TYR:O	1:A:79:VAL:C	2.49	0.46
1:A:25:PHE:CZ	1:A:41:HIS:CD2	3.03	0.46
1:A:25:PHE:C	1:A:57:LEU:HD13	2.36	0.46
1:A:16:ARG:HA	1:A:26:ILE:O	2.15	0.46
1:A:17:PHE:CZ	1:A:61:ALA:HA	2.49	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:80:SER:OG	1:A:81:ASP:N	2.48	0.46
1:A:19:THR:CG2	1:A:24:ALA:HB3	2.46	0.46
1:A:32:ASN:ND2	1:A:35:LYS:HB2	2.26	0.46
1:A:31:ARG:HB2	1:A:36:VAL:HG11	1.97	0.46
1:A:49:ARG:NH2	3:A:124:HOH:O	2.48	0.46
1:A:91:LYS:NZ	3:A:139:HOH:O	2.48	0.46
1:A:32:ASN:HB3	1:A:36:VAL:HG23	1.98	0.46
1:A:79:VAL:O	1:A:84:LEU:HG	2.16	0.46
1:A:31:ARG:HD2	1:A:73:ILE:HD12	1.97	0.45
1:A:65:ALA:HB2	1:A:72:ILE:CD1	2.46	0.45
1:A:77:SER:O	1:A:81:ASP:HB2	2.15	0.45
1:A:36:VAL:CG1	1:A:73:ILE:HG13	2.46	0.45
1:A:84:LEU:N	1:A:85:PRO:CD	2.78	0.45
1:A:45:PRO:CG	1:A:48:LYS:CD	2.82	0.45
1:A:25:PHE:HE1	1:A:27:GLU:OE2	2.00	0.45
1:A:53:LEU:HD12	1:A:53:LEU:O	2.16	0.45
1:A:84:LEU:HB2	1:A:85:PRO:HD3	1.99	0.45
2:A:104:COA:C5B	3:A:167:HOH:O	2.65	0.45
1:A:84:LEU:N	1:A:85:PRO:CD	2.80	0.45
1:A:10:TRP:CZ2	1:A:12:GLU:HG2	2.51	0.45
1:A:15:ARG:HA	1:A:28:TYR:CZ	2.52	0.45
1:A:86:ARG:CZ	2:A:104:COA:O9A	2.65	0.45
2:A:104:COA:H51A	3:A:120:HOH:O	2.17	0.44
1:A:31:ARG:HG2	1:A:31:ARG:NH1	2.23	0.44
1:A:62:PHE:CD2	1:A:90:TRP:HB3	2.53	0.44
1:A:37:MET:SD	1:A:72:ILE:CD1	3.03	0.44
1:A:80:SER:OG	3:A:114:HOH:O	2.21	0.44
1:A:44:VAL:HB	2:A:104:COA:H122	1.99	0.44
1:A:83:PHE:O	1:A:87:ASN:ND2	2.37	0.44
1:A:21:ASP:OD1	1:A:23:GLU:HB2	2.18	0.44
1:A:44:VAL:HG13	1:A:45:PRO:HD2	1.99	0.44
1:A:62:PHE:HB3	1:A:90:TRP:CE3	2.53	0.44
1:A:5:PRO:HB2	1:A:6:PRO:O	2.17	0.44
1:A:65:ALA:HB1	1:A:93:LEU:HD11	1.99	0.44
1:A:10:TRP:CE2	1:A:12:GLU:HG2	2.53	0.44
2:A:104:COA:N8P	2:A:104:COA:CDP	2.79	0.44
1:A:68:HIS:O	1:A:69:SER:HB2	2.17	0.44
1:A:10:TRP:CZ2	1:A:12:GLU:HG2	2.53	0.43
2:A:104:COA:HN8	2:A:104:COA:CDP	2.19	0.43
1:A:38:ASP:CB	1:A:73:ILE:HB	2.46	0.43
1:A:22:HIS:HB3	3:A:154:HOH:O	2.17	0.43
1:A:25:PHE:CZ	1:A:43:TYR:HB3	2.53	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:59:VAL:O	1:A:59:VAL:CG1	2.67	0.43
1:A:15:ARG:HB3	1:A:28:TYR:O	2.18	0.43
1:A:46:SER:O	1:A:47:PHE:C	2.55	0.43
1:A:31:ARG:HB2	1:A:36:VAL:CB	2.48	0.43
1:A:32:ASN:ND2	3:A:147:HOH:O	2.52	0.42
1:A:24:ALA:HB2	1:A:45:PRO:CD	2.49	0.42
1:A:37:MET:HB3	1:A:72:ILE:HG12	2.00	0.42
1:A:24:ALA:HB2	1:A:45:PRO:HD3	2.00	0.42
1:A:6:PRO:HB2	3:A:149:HOH:O	2.19	0.42
1:A:6:PRO:HB2	1:A:7:LYS:H	1.44	0.42
1:A:16:ARG:NH2	1:A:27:GLU:CD	2.73	0.42
1:A:49:ARG:O	1:A:51:LEU:HG	2.19	0.42
1:A:41:HIS:CE1	3:A:163:HOH:O	2.71	0.42
1:A:84:LEU:CB	1:A:85:PRO:HD3	2.48	0.42
1:A:37:MET:O	1:A:72:ILE:HA	2.20	0.42
1:A:28:TYR:HE1	1:A:30:MET:HG2	1.85	0.42
1:A:15:ARG:HA	1:A:28:TYR:OH	2.20	0.42
1:A:50:GLY:CA	2:A:104:COA:O1A	2.68	0.42
1:A:40:VAL:HA	1:A:75:SER:HB2	2.01	0.42
1:A:86:ARG:HB2	1:A:86:ARG:NH1	2.34	0.42
1:A:91:LYS:O	1:A:92:PRO:C	2.58	0.42
1:A:65:ALA:CB	1:A:72:ILE:HD12	2.50	0.42
1:A:77:SER:O	1:A:81:ASP:HB2	2.19	0.42
1:A:48:LYS:HE2	1:A:53:LEU:CD2	2.50	0.42
1:A:20:GLU:C	1:A:22:HIS:N	2.73	0.42
1:A:78:TYR:CA	2:A:104:COA:H2A	2.38	0.41
1:A:19:THR:HG21	1:A:53:LEU:HD21	2.02	0.41
1:A:91:LYS:NZ	3:A:139:HOH:O	2.50	0.41
1:A:39:LEU:CD1	1:A:79:VAL:HG21	2.50	0.41
1:A:36:VAL:HG11	1:A:73:ILE:HD11	2.02	0.41
1:A:54:ALA:HA	1:A:57:LEU:HD12	2.01	0.41
1:A:49:ARG:HG2	2:A:104:COA:H121	2.03	0.41
1:A:47:PHE:H	1:A:47:PHE:HD1	1.57	0.41
1:A:82:THR:HG21	2:A:104:COA:H1B	2.02	0.41
1:A:20:GLU:HG3	3:A:165:HOH:O	2.20	0.41
1:A:78:TYR:HA	2:A:104:COA:C2A	2.51	0.41
1:A:90:TRP:CZ3	1:A:93:LEU:CD2	3.04	0.41
1:A:10:TRP:CE2	1:A:12:GLU:HG2	2.56	0.41
1:A:24:ALA:HB1	1:A:53:LEU:CD2	2.40	0.41
1:A:86:ARG:HG2	2:A:104:COA:O8A	2.21	0.41
1:A:49:ARG:NE	2:A:104:COA:O9P	2.55	0.40
1:A:51:LEU:N	1:A:51:LEU:HD23	2.36	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:88:PRO:C	1:A:90:TRP:H	2.25	0.40
1:A:24:ALA:HB2	1:A:53:LEU:CD2	2.48	0.40
1:A:78:TYR:HA	2:A:104:COA:H2A	2.03	0.40
1:A:65:ALA:HA	1:A:70:ILE:HB	2.03	0.40
1:A:65:ALA:CB	1:A:93:LEU:HD11	2.52	0.40
1:A:80:SER:HA	1:A:84:LEU:HD12	2.03	0.40
1:A:44:VAL:CG2	2:A:104:COA:H122	2.51	0.40
1:A:16:ARG:NH2	1:A:27:GLU:OE2	2.55	0.40
1:A:59:VAL:O	1:A:63:GLU:CG	2.68	0.40
1:A:44:VAL:HG23	2:A:104:COA:H122	2.03	0.40

All (2) 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:A:81:ASP:OD2	3:A:129:HOH:O[1_554]	2.12	0.08
1:A:86:ARG:NH2	3:A:107:HOH:O[1_554]	2.12	0.08

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	1-A	89/103 (86%)	79 (89%)	6 (7%)	4 (4%)	4	0
1	2-A	89/103 (86%)	86 (97%)	3 (3%)	0	100	100
1	3-A	89/103 (86%)	81 (91%)	8 (9%)	0	100	100
1	4-A	89/103 (86%)	82 (92%)	7 (8%)	0	100	100
All	All	356/412 (86%)	328 (92%)	24 (7%)	4 (1%)	21	8

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	1-A	46	SER
1	1-A	47	PHE
1	1-A	6	PRO

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Mol	Chain	Res	Type
1	1-A	20	GLU

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	1-A	82/93 (88%)	80 (98%)	2 (2%)	61	57
1	2-A	82/93 (88%)	75 (92%)	7 (8%)	15	7
1	3-A	82/93 (88%)	76 (93%)	6 (7%)	20	10
1	4-A	82/93 (88%)	78 (95%)	4 (5%)	35	24
All	All	328/372 (88%)	309 (94%)	19 (6%)	28	18

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	1-A	77	SER
1	1-A	93	LEU
1	2-A	18	GLU
1	2-A	33	ASN
1	2-A	46	SER
1	2-A	47	PHE
1	2-A	76	CYS
1	2-A	86	ARG
1	2-A	89	SER
1	3-A	47	PHE
1	3-A	58	CYS
1	3-A	72	ILE
1	3-A	76	CYS
1	3-A	86	ARG
1	3-A	89	SER
1	4-A	47	PHE
1	4-A	67	SER
1	4-A	72	ILE
1	4-A	93	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such

sidechains are listed below:

Mol	Chain	Res	Type
1	1-A	32	ASN
1	1-A	33	ASN
1	1-A	64	HIS
1	2-A	32	ASN
1	2-A	41	HIS
1	2-A	64	HIS
1	3-A	64	HIS
1	4-A	33	ASN
1	4-A	64	HIS

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 ⓘ

4 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	$\# Z > 2$
2	COA	1-A	104	-	50,50,50	2.69	18 (36%)	75,75,75	3.35	33 (44%)
2	COA	2-A	104	-	50,50,50	3.56	24 (48%)	75,75,75	3.21	30 (40%)
2	COA	3-A	104	-	50,50,50	3.49	23 (46%)	75,75,75	3.43	32 (42%)
2	COA	4-A	104	-	50,50,50	4.37	26 (52%)	75,75,75	3.73	35 (46%)

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
2	COA	1-A	104	-	-	1/48/64/64	0/1/3/3
2	COA	2-A	104	-	-	2/48/64/64	0/1/3/3
2	COA	3-A	104	-	-	0/48/64/64	0/1/3/3
2	COA	4-A	104	-	-	0/48/64/64	0/1/3/3

All (91) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	4-A	104	COA	P2A-O3A	-14.52	1.33	1.59
2	2-A	104	COA	C2B-C1B	12.08	1.70	1.53
2	3-A	104	COA	CCP-CBP	12.02	1.71	1.52
2	2-A	104	COA	CBP-CAP	11.51	1.67	1.55
2	4-A	104	COA	C2B-C1B	10.96	1.69	1.53
2	3-A	104	COA	CBP-CAP	9.30	1.65	1.55
2	4-A	104	COA	CCP-CBP	9.18	1.67	1.52
2	1-A	104	COA	C2B-C1B	9.09	1.66	1.53
2	1-A	104	COA	CBP-CAP	-8.91	1.46	1.55
2	3-A	104	COA	C2B-C1B	8.83	1.66	1.53
2	4-A	104	COA	C9P-N8P	8.22	1.51	1.33
2	4-A	104	COA	C3P-N4P	-7.29	1.28	1.46
2	4-A	104	COA	C5P-N4P	-6.95	1.16	1.33
2	2-A	104	COA	C7P-N8P	6.41	1.61	1.46
2	4-A	104	COA	C8A-N9A	6.11	1.45	1.36
2	4-A	104	COA	O9P-C9P	6.09	1.35	1.23
2	3-A	104	COA	O5B-C5B	-5.86	1.20	1.44
2	4-A	104	COA	C4A-N9A	5.46	1.45	1.37
2	3-A	104	COA	C5P-N4P	5.31	1.45	1.33
2	2-A	104	COA	C2A-N1A	-5.21	1.23	1.33
2	3-A	104	COA	OAP-CAP	-5.19	1.32	1.42
2	3-A	104	COA	P2A-O3A	5.08	1.69	1.59
2	1-A	104	COA	CCP-CBP	4.93	1.60	1.52
2	3-A	104	COA	O4B-C1B	-4.86	1.33	1.41
2	4-A	104	COA	P2A-O5A	4.84	1.76	1.55
2	1-A	104	COA	P1A-O5B	4.76	1.80	1.59
2	4-A	104	COA	O6A-CCP	4.71	1.61	1.43
2	4-A	104	COA	P1A-O1A	4.65	1.69	1.51
2	4-A	104	COA	C2P-S1P	-4.64	1.63	1.80
2	4-A	104	COA	C6P-C5P	4.63	1.60	1.51
2	3-A	104	COA	C4A-N9A	-4.61	1.31	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	1-A	104	COA	O9P-C9P	-4.59	1.14	1.23
2	2-A	104	COA	C9P-N8P	4.57	1.43	1.33
2	4-A	104	COA	O4B-C4B	-4.56	1.34	1.45
2	2-A	104	COA	OAP-CAP	-4.45	1.34	1.42
2	1-A	104	COA	CAP-C9P	4.36	1.66	1.53
2	2-A	104	COA	O5B-C5B	-4.32	1.26	1.44
2	2-A	104	COA	C2B-C3B	4.28	1.63	1.53
2	2-A	104	COA	P2A-O3A	4.25	1.67	1.59
2	4-A	104	COA	C5A-C4A	4.22	1.50	1.40
2	2-A	104	COA	P1A-O5B	-4.05	1.40	1.59
2	4-A	104	COA	P3B-O3B	3.95	1.72	1.59
2	2-A	104	COA	P3B-O8A	3.91	1.69	1.54
2	2-A	104	COA	C8A-N9A	3.73	1.42	1.36
2	4-A	104	COA	C5B-C4B	3.69	1.63	1.51
2	4-A	104	COA	CBP-CAP	3.60	1.59	1.55
2	3-A	104	COA	P1A-O5B	-3.54	1.43	1.59
2	3-A	104	COA	C3P-N4P	3.52	1.54	1.46
2	1-A	104	COA	C7P-N8P	-3.50	1.37	1.46
2	4-A	104	COA	CEP-CBP	3.49	1.61	1.53
2	2-A	104	COA	CAP-C9P	-3.48	1.42	1.53
2	2-A	104	COA	O6A-CCP	3.47	1.56	1.43
2	1-A	104	COA	O4B-C1B	-3.41	1.36	1.41
2	3-A	104	COA	O9P-C9P	3.35	1.30	1.23
2	3-A	104	COA	P3B-O9A	3.24	1.66	1.54
2	2-A	104	COA	C6A-N6A	3.19	1.45	1.35
2	2-A	104	COA	O4B-C1B	3.13	1.46	1.41
2	2-A	104	COA	C3B-C4B	3.13	1.62	1.52
2	3-A	104	COA	C6A-N6A	3.00	1.44	1.35
2	2-A	104	COA	P2A-O5A	-2.97	1.41	1.55
2	1-A	104	COA	C4A-N9A	2.84	1.41	1.37
2	2-A	104	COA	P3B-O9A	2.83	1.65	1.54
2	2-A	104	COA	C1B-N9A	-2.73	1.40	1.48
2	3-A	104	COA	C9P-N8P	2.69	1.39	1.33
2	3-A	104	COA	P1A-O1A	-2.66	1.41	1.51
2	4-A	104	COA	P1A-O5B	2.66	1.71	1.59
2	1-A	104	COA	C2B-C3B	2.59	1.59	1.53
2	3-A	104	COA	C2P-C3P	2.58	1.61	1.51
2	2-A	104	COA	CCP-CBP	-2.58	1.48	1.52
2	4-A	104	COA	C1B-N9A	2.58	1.56	1.48
2	1-A	104	COA	P3B-O9A	2.57	1.64	1.54
2	3-A	104	COA	O6A-CCP	2.52	1.53	1.43
2	4-A	104	COA	P3B-O8A	2.46	1.63	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	4-A	104	COA	C3B-C4B	2.45	1.60	1.52
2	3-A	104	COA	P2A-O5A	-2.43	1.44	1.55
2	1-A	104	COA	C3P-N4P	-2.42	1.40	1.46
2	1-A	104	COA	O4B-C4B	2.39	1.50	1.45
2	1-A	104	COA	P3B-O3B	-2.38	1.52	1.59
2	3-A	104	COA	C2P-S1P	2.34	1.89	1.80
2	4-A	104	COA	C2P-C3P	-2.32	1.41	1.51
2	1-A	104	COA	C9P-N8P	-2.31	1.28	1.33
2	3-A	104	COA	C2A-N3A	2.28	1.36	1.32
2	2-A	104	COA	C4A-N9A	2.27	1.41	1.37
2	3-A	104	COA	P1A-O2A	-2.25	1.45	1.55
2	2-A	104	COA	P1A-O2A	-2.22	1.45	1.55
2	1-A	104	COA	CDP-CBP	-2.22	1.48	1.53
2	1-A	104	COA	O2B-C2B	2.21	1.48	1.43
2	1-A	104	COA	C5A-C4A	2.18	1.45	1.40
2	4-A	104	COA	OAP-CAP	-2.15	1.38	1.42
2	3-A	104	COA	C7P-N8P	-2.09	1.41	1.46
2	2-A	104	COA	C8A-N7A	-2.00	1.30	1.34

All (130) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	4-A	104	COA	CBP-CAP-C9P	-10.64	102.41	112.73
2	2-A	104	COA	O9A-P3B-O3B	-10.59	76.61	107.09
2	3-A	104	COA	CCP-CBP-CAP	10.45	123.97	108.70
2	3-A	104	COA	N3A-C2A-N1A	-10.20	120.18	128.71
2	4-A	104	COA	C3P-C2P-S1P	-10.01	87.25	113.50
2	4-A	104	COA	O6A-CCP-CBP	9.59	126.62	110.57
2	4-A	104	COA	O9A-P3B-O8A	-9.06	72.32	107.61
2	3-A	104	COA	P3B-O3B-C3B	-8.96	103.09	121.96
2	1-A	104	COA	P2A-O3A-P1A	-8.78	105.94	131.68
2	1-A	104	COA	CCP-CBP-CAP	8.62	121.30	108.70
2	4-A	104	COA	O9A-P3B-O3B	-8.56	82.44	107.09
2	1-A	104	COA	N3A-C2A-N1A	-8.54	121.57	128.71
2	1-A	104	COA	O9A-P3B-O7A	-8.15	83.80	110.44
2	4-A	104	COA	CEP-CBP-CCP	7.96	120.23	108.76
2	1-A	104	COA	O9A-P3B-O3B	-7.95	84.20	107.09
2	3-A	104	COA	O9A-P3B-O3B	-7.72	84.85	107.09
2	2-A	104	COA	C4B-O4B-C1B	-7.42	101.68	109.75
2	2-A	104	COA	P3B-O3B-C3B	-7.30	106.59	121.96
2	4-A	104	COA	N3A-C2A-N1A	-7.21	122.68	128.71
2	1-A	104	COA	P3B-O3B-C3B	-7.12	106.96	121.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	3-A	104	COA	O9A-P3B-O7A	-7.12	87.16	110.44
2	4-A	104	COA	CEP-CBP-CAP	-7.00	96.68	108.82
2	2-A	104	COA	O3B-P3B-O7A	6.72	125.57	106.79
2	1-A	104	COA	O3B-P3B-O7A	6.70	125.50	106.79
2	2-A	104	COA	N3A-C2A-N1A	-6.50	123.28	128.71
2	1-A	104	COA	O9A-P3B-O8A	-6.46	82.44	107.61
2	2-A	104	COA	N3A-C4A-N9A	6.42	137.02	125.43
2	4-A	104	COA	P2A-O3A-P1A	-6.35	113.06	131.68
2	4-A	104	COA	O9A-P3B-O7A	-6.29	89.89	110.44
2	2-A	104	COA	O8A-P3B-O3B	6.23	125.02	107.09
2	2-A	104	COA	O4B-C1B-N9A	-6.15	102.72	108.44
2	2-A	104	COA	O9A-P3B-O7A	-6.14	90.36	110.44
2	2-A	104	COA	CDP-CBP-CCP	6.13	117.59	108.76
2	4-A	104	COA	N3A-C4A-N9A	6.13	136.50	125.43
2	4-A	104	COA	O9P-C9P-CAP	-5.82	109.87	120.48
2	3-A	104	COA	CBP-CAP-C9P	-5.78	107.12	112.73
2	3-A	104	COA	O9A-P3B-O8A	-5.66	85.57	107.61
2	2-A	104	COA	CEP-CBP-CCP	-5.59	100.70	108.76
2	2-A	104	COA	O9A-P3B-O8A	-5.35	86.77	107.61
2	3-A	104	COA	O6A-CCP-CBP	-5.21	101.86	110.57
2	1-A	104	COA	O4B-C1B-N9A	5.15	113.23	108.44
2	4-A	104	COA	O4B-C4B-C3B	5.05	116.35	104.88
2	1-A	104	COA	C6P-C7P-N8P	-5.01	101.08	111.87
2	1-A	104	COA	C4A-C5A-N7A	-4.96	105.27	109.52
2	3-A	104	COA	O8A-P3B-O3B	4.85	121.05	107.09
2	1-A	104	COA	N3A-C4A-N9A	4.83	134.15	125.43
2	3-A	104	COA	C8A-N9A-C4A	4.82	110.58	106.90
2	3-A	104	COA	O3B-P3B-O7A	4.82	120.25	106.79
2	3-A	104	COA	O4B-C4B-C3B	4.80	115.78	104.88
2	3-A	104	COA	C7P-N8P-C9P	4.77	132.37	122.57
2	1-A	104	COA	CEP-CBP-CAP	-4.69	100.68	108.82
2	3-A	104	COA	P2A-O3A-P1A	-4.35	118.91	131.68
2	2-A	104	COA	C2A-N3A-C4A	4.32	126.31	114.01
2	1-A	104	COA	C3P-C2P-S1P	-4.26	102.32	113.50
2	4-A	104	COA	O3A-P1A-O5B	-4.24	84.46	103.41
2	1-A	104	COA	O3A-P1A-O5B	-4.14	84.89	103.41
2	4-A	104	COA	O3B-P3B-O7A	4.14	118.35	106.79
2	3-A	104	COA	CDP-CBP-CAP	-4.12	101.68	108.82
2	2-A	104	COA	O9P-C9P-CAP	-4.12	112.97	120.48
2	3-A	104	COA	CAP-C9P-N8P	-4.11	107.77	116.57
2	2-A	104	COA	O4B-C4B-C3B	3.95	113.85	104.88
2	1-A	104	COA	O9P-C9P-N8P	-3.94	115.05	123.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	3-A	104	COA	CEP-CBP-CCP	-3.93	103.10	108.76
2	1-A	104	COA	C5A-C4A-N3A	-3.91	117.18	125.70
2	4-A	104	COA	O8A-P3B-O3B	3.91	118.34	107.09
2	2-A	104	COA	C5A-C4A-N3A	-3.89	117.23	125.70
2	1-A	104	COA	O5B-C5B-C4B	-3.80	94.99	108.94
2	4-A	104	COA	C8A-N9A-C4A	-3.77	104.02	106.90
2	3-A	104	COA	C1B-N9A-C4A	-3.75	120.15	126.64
2	2-A	104	COA	C8A-N9A-C4A	3.61	109.66	106.90
2	4-A	104	COA	C5A-C4A-N3A	-3.58	117.92	125.70
2	3-A	104	COA	C7P-C6P-C5P	-3.58	106.16	112.25
2	3-A	104	COA	CDP-CBP-CCP	-3.54	103.65	108.76
2	1-A	104	COA	CBP-CAP-C9P	3.53	116.15	112.73
2	3-A	104	COA	C4A-C5A-N7A	-3.48	106.54	109.52
2	1-A	104	COA	O2A-P1A-O3A	3.47	121.60	105.14
2	3-A	104	COA	O5B-C5B-C4B	-3.43	96.36	108.94
2	3-A	104	COA	N3A-C4A-N9A	3.42	131.61	125.43
2	4-A	104	COA	O2A-P1A-O3A	3.36	121.08	105.14
2	4-A	104	COA	P3B-O3B-C3B	-3.35	114.90	121.96
2	1-A	104	COA	OAP-CAP-CBP	-3.34	102.94	110.71
2	3-A	104	COA	O9P-C9P-CAP	3.29	126.48	120.48
2	1-A	104	COA	O8A-P3B-O7A	3.21	120.93	110.44
2	4-A	104	COA	O9P-C9P-N8P	3.21	129.56	123.05
2	4-A	104	COA	O5P-C5P-C6P	3.10	128.09	121.92
2	2-A	104	COA	O6A-P2A-O4A	3.08	121.44	109.37
2	1-A	104	COA	CEP-CBP-CDP	3.03	113.40	109.08
2	3-A	104	COA	OAP-CAP-CBP	2.99	117.67	110.71
2	2-A	104	COA	O3A-P1A-O5B	-2.99	90.06	103.41
2	3-A	104	COA	CEP-CBP-CDP	2.94	113.27	109.08
2	4-A	104	COA	C6P-C7P-N8P	2.93	118.18	111.87
2	2-A	104	COA	CEP-CBP-CDP	2.90	113.21	109.08
2	4-A	104	COA	O5B-P1A-O1A	2.83	120.45	109.37
2	4-A	104	COA	C3P-N4P-C5P	-2.80	117.13	122.84
2	4-A	104	COA	O5P-C5P-N4P	-2.80	117.39	122.94
2	3-A	104	COA	C4B-O4B-C1B	-2.78	106.72	109.75
2	2-A	104	COA	O9P-C9P-N8P	2.77	128.67	123.05
2	4-A	104	COA	C2A-N3A-C4A	2.75	121.84	114.01
2	3-A	104	COA	O3A-P1A-O5B	-2.72	91.24	103.41
2	4-A	104	COA	C2P-C3P-N4P	-2.72	106.36	112.45
2	2-A	104	COA	O3A-P2A-O6A	-2.70	91.33	103.41
2	3-A	104	COA	C2A-N1A-C6A	2.67	123.58	118.77
2	1-A	104	COA	O6A-CCP-CBP	-2.66	106.12	110.57
2	2-A	104	COA	O5B-C5B-C4B	-2.65	99.21	108.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	1-A	104	COA	O3A-P2A-O6A	-2.65	91.57	103.41
2	1-A	104	COA	C2A-N3A-C4A	2.61	121.44	114.01
2	4-A	104	COA	C2B-C3B-C4B	-2.59	98.13	103.16
2	1-A	104	COA	C3B-C2B-C1B	2.54	106.00	99.98
2	1-A	104	COA	C2P-C3P-N4P	-2.52	106.79	112.45
2	2-A	104	COA	CCP-CBP-CAP	2.51	112.36	108.70
2	1-A	104	COA	O4B-C4B-C3B	2.49	110.53	104.88
2	2-A	104	COA	C2B-C1B-N9A	2.47	119.62	113.27
2	2-A	104	COA	O2B-C2B-C1B	2.46	118.67	111.23
2	1-A	104	COA	CAP-C9P-N8P	2.38	121.67	116.57
2	4-A	104	COA	OAP-CAP-CBP	2.33	116.13	110.71
2	3-A	104	COA	C6P-C5P-N4P	2.33	120.73	116.50
2	3-A	104	COA	C2A-N3A-C4A	2.29	120.52	114.01
2	2-A	104	COA	N7A-C8A-N9A	-2.28	107.91	114.36
2	1-A	104	COA	C3P-N4P-C5P	-2.25	118.25	122.84
2	2-A	104	COA	CEP-CBP-CAP	-2.24	104.94	108.82
2	1-A	104	COA	CDP-CBP-CCP	-2.21	105.57	108.76
2	2-A	104	COA	O5B-P1A-O1A	2.20	117.99	109.37
2	4-A	104	COA	C3B-C2B-C1B	2.19	105.19	99.98
2	4-A	104	COA	P2A-O6A-CCP	-2.18	109.36	121.66
2	1-A	104	COA	O4B-C4B-C5B	2.10	116.86	109.36
2	4-A	104	COA	O5A-P2A-O3A	2.09	115.07	105.14
2	2-A	104	COA	O3B-C3B-C4B	2.09	118.23	110.06
2	4-A	104	COA	O4B-C1B-C2B	-2.07	103.61	106.77
2	4-A	104	COA	C2B-C1B-N9A	2.05	118.54	113.27
2	3-A	104	COA	O8A-P3B-O7A	2.05	117.14	110.44

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	2-A	104	COA	O9P-C9P-CAP-CBP
2	2-A	104	COA	CBP-CAP-C9P-N8P
2	1-A	104	COA	O9P-C9P-CAP-CBP

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	1-A	91/103 (88%)	-0.27	1 (1%) 77 80	18, 26, 42, 74	91 (100%)
1	2-A	91/103 (88%)	-0.27	1 (1%) 77 80	18, 26, 42, 74	91 (100%)
1	3-A	91/103 (88%)	-0.27	1 (1%) 77 80	18, 26, 42, 74	91 (100%)
1	4-A	91/103 (88%)	-0.27	1 (1%) 77 80	18, 26, 42, 74	91 (100%)
All	All	364/412 (88%)	-0.27	4 (1%) 59 80	18, 27, 47, 74	364 (100%)

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	1-A	5	PRO	7.6
1	2-A	5	PRO	7.6
1	3-A	5	PRO	7.6
1	4-A	5	PRO	7.6

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(\AA^2)	Q<0.9
2	COA	3-A	104	48/48	0.39	10.79	9,40,58,59	48
2	COA	4-A	104	48/48	0.39	10.01	24,42,53,57	48
2	COA	1-A	104	48/48	0.39	9.54	19,34,50,57	48
2	COA	2-A	104	48/48	0.39	9.28	12,43,56,60	48

6.5 Other polymers ⓘ

There are no such residues in this entry.