



Full wwPDB X-ray Structure Validation Report ⓘ

May 14, 2020 – 10:38 pm BST

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 X-ray Structure Validation Report for a publicly released PDB entry.

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A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

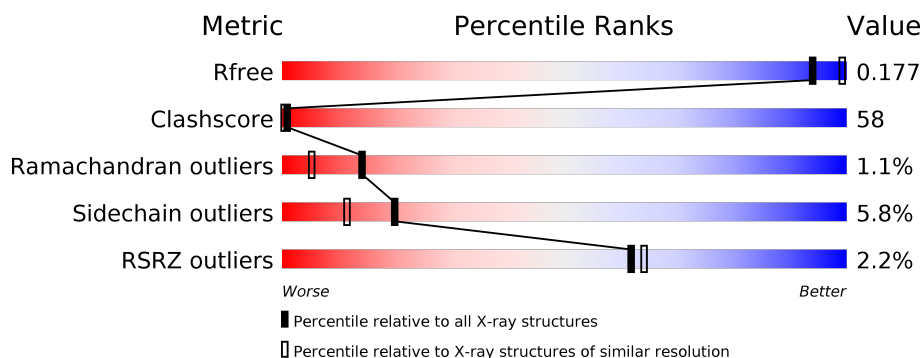
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.06 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}	130704	2684 (2.08-2.04)
Clashscore	141614	2801 (2.08-2.04)
Ramachandran outliers	138981	2768 (2.08-2.04)
Sidechain outliers	138945	2768 (2.08-2.04)
RSRZ outliers	127900	2646 (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 respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	1-A	103	<div> <div></div> <div> <div></div> <div>29%</div> <div>45%</div> <div>15%</div> <div>12%</div> </div> </div>
1	2-A	103	<div> <div></div> <div> <div></div> <div>22%</div> <div>50%</div> <div>16%</div> <div>12%</div> </div> </div>
1	3-A	103	<div> <div></div> <div> <div></div> <div>29%</div> <div>45%</div> <div>12%</div> <div>12%</div> </div> </div>
1	4-A	103	<div> <div></div> <div> <div></div> <div>25%</div> <div>48%</div> <div>15%</div> <div>12%</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit crite-

ria:

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

2 Entry composition [i](#)

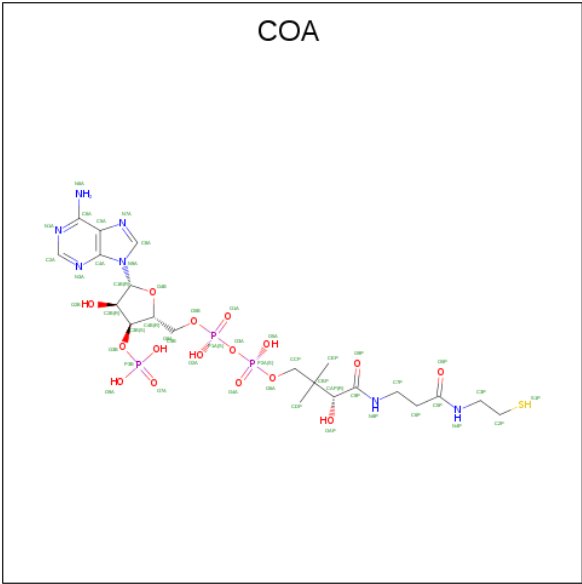
There are 3 unique types of molecules in this entry. The entry contains 3404 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called 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: C₂₁H₃₆N₇O₁₆P₃S).



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		

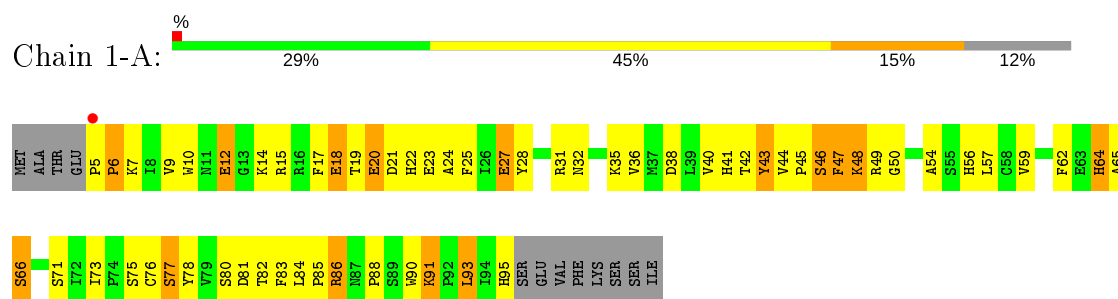
- Molecule 3 is water.

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		

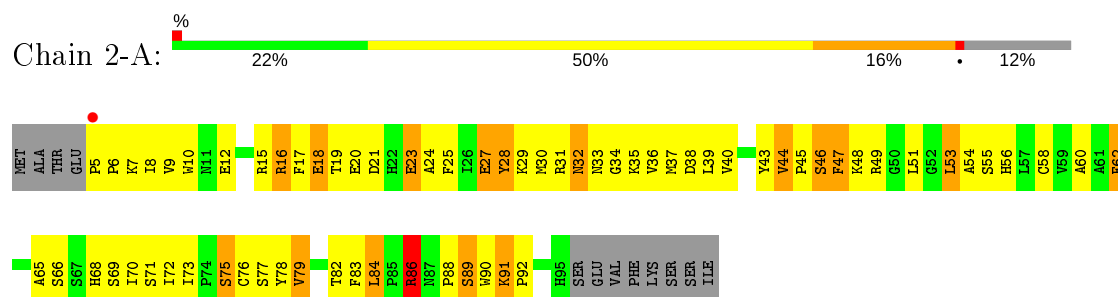
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

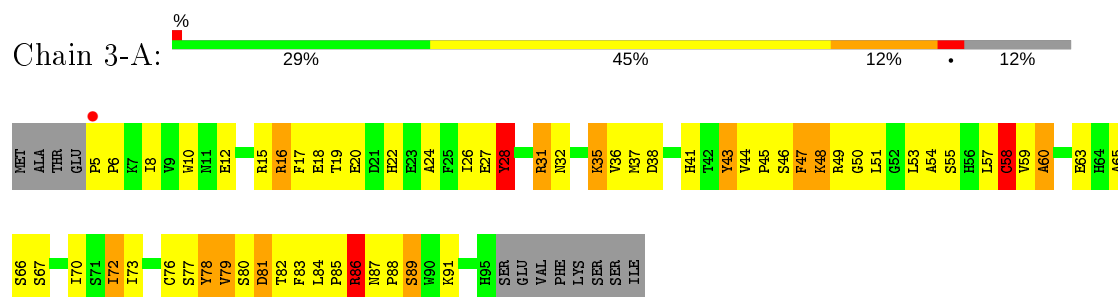
• Molecule 1: Uncharacterized protein At1g77540



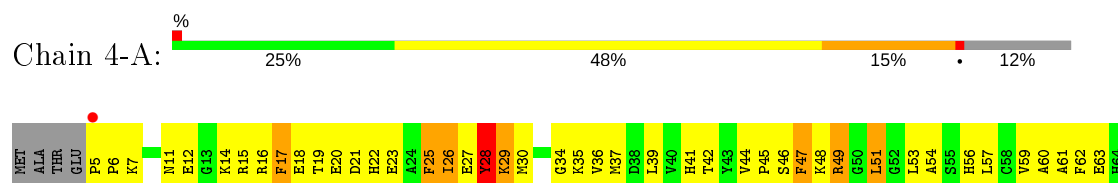
• Molecule 1: Uncharacterized protein At1g77540



• Molecule 1: Uncharacterized protein At1g77540



• Molecule 1: Uncharacterized protein At1g77540



A65	S66	S67	H68	S69		I72	I73	P74	S75	C76	S77	Y78	V79	D81	T82	F83	L84	P86	R86	N87	P88	S89	I90	K91	P92	L93	I94	H95	SER	GLU	VAL	PHE	LYS	SER	SER	ILE
-----	-----	-----	-----	-----	--	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----

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 0.177 , 0.177	Depositor DCC
R_{free} test set	636 reflections (9.85%)	wwPDB-VP
Wilson B-factor (Å ²)	29.0	Xtriage
Anisotropy	0.036	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 64.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.066 for h,-k,-l	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.

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: 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 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	1-A	735	0	720	76	0
1	2-A	735	0	720	86	0
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	0
3	2-A	68	0	0	11	0
3	3-A	68	0	0	11	0
3	4-A	68	0	0	22	0
All	All	3404	0	3005	353	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 58.

All (353) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:42:THR:H	2:A:104:COA:H142	1.01	1.13
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:HB2	1.41	1.01
1:A:45:PRO:HG2	1:A:48:LYS:HD3	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
1:A:31:ARG:HG2	1:A:31:ARG:HH11	1.46	0.80
2:A:104:COA:CDP	2:A:104:COA:HN8	1.95	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:42:THR:HG22	2:A:104:COA:OAP	1.83	0.78
1:A:49:ARG:HD3	2:A:104:COA:H71	1.66	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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: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:37:MET:SD	1:A:72:ILE:HD11	2.27	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:90:TRP:CE3	1:A:93:LEU:HD23	2.24	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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:30:MET:CE	1:A:34:GLY:HA2	2.26	0.66
2:A:104:COA:HN8	2:A:104:COA:H131	1.60	0.66
1:A:46:SER:CA	1:A:49:ARG:NH2	2.59	0.66
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
1:A:56:HIS:HE1	3:A:158:HOH:O	1.81	0.63
2:A:104:COA:H133	2:A:104:COA:N8P	2.13	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:24:ALA:HA	1:A:43:TYR:O	2.00	0.62
1:A:29:LYS:HG3	3:A:161:HOH:O	1.99	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:O5A	2:A:104:COA:O1A	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:23:GLU:OE2	1:A:45:PRO:CG	2.39	0.61
1:A:80:SER:HA	1:A:84:LEU:HD12	1.82	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:CG2	2:A:104:COA:H122	2.32	0.59
1:A:44:VAL:HB	2:A:104:COA:H122	1.83	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:25:PHE:CD1	1:A:27:GLU:HG3	2.37	0.59
1:A:29:LYS:CE	3:A:162:HOH:O	2.23	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:84:LEU:HB2	1:A:85:PRO:HD3	1.85	0.58
1:A:81:ASP:O	1:A:85:PRO:HG2	2.02	0.58
1:A:7:LYS:HE2	1:A:9:VAL:HG12	1.85	0.58
1:A:45:PRO:HG2	1:A:48:LYS:HB2	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
1:A:14:LYS:O	1:A:15:ARG:HB2	2.03	0.58
1:A:29:LYS:NZ	3:A:162:HOH:O	2.37	0.58
1:A:54:ALA:HB2	2:A:104:COA:H143	1.85	0.57
1:A:42:THR:HG21	2:A:104:COA:H142	1.86	0.57
1:A:5:PRO:HG2	3:A:149:HOH:O	2.04	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:49:ARG:O	1:A:51:LEU:HG	2.05	0.56
1:A:86:ARG:NE	2:A:104:COA:O9A	2.38	0.56
1:A:54:ALA:HB2	2:A:104:COA:H143	1.86	0.56
1:A:19:THR:CG2	1:A:24:ALA:HB3	2.35	0.56
1:A:25:PHE:CE1	1:A:41:HIS:CD2	2.93	0.56
1:A:15:ARG:NH1	1:A:29:LYS:HG2	2.21	0.56
1:A:54:ALA:HB2	2:A:104:COA:O6A	2.06	0.56
1:A:31:ARG:HB2	1:A:36:VAL:HB	1.87	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:10:TRP:HZ2	1:A:28:TYR:HH	1.53	0.55
1:A:21:ASP:OD2	1:A:23:GLU:HB2	2.07	0.55
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:32:ASN:ND2	1:A:35:LYS:HD2	2.22	0.54
1:A:65:ALA:CB	1:A:72:ILE:HD11	2.37	0.54
1:A:41:HIS:CE1	3:A:163:HOH:O	2.60	0.54
1:A:25:PHE:CE2	1:A:43:TYR:HB3	2.42	0.54
1:A:30:MET:HE2	1:A:34:GLY:HA2	1.90	0.54
1:A:38:ASP:O	1:A:40:VAL:HG13	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:34:GLY:O	3:A:170:HOH:O	2.19	0.54
2:A:104:COA:H8A	3:A:144:HOH:O	2.07	0.54
1:A:75:SER:O	1:A:76:CYS:C	2.46	0.54
1:A:31:ARG:C	1:A:36:VAL:HB	2.28	0.54
1:A:73:ILE:HG22	1:A:75:SER:OG	2.08	0.54
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
2:A:104:COA:N8P	2:A:104:COA:H133	2.13	0.53
1:A:30:MET:HE3	1:A:34:GLY:HA2	1.90	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:54:ALA:HB2	2:A:104:COA:CEP	2.39	0.52
1:A:42:THR:O	2:A:104:COA:H122	2.09	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:19:THR:HB	3:A:123:HOH:O	2.09	0.52
1:A:28:TYR:HA	1:A:38:ASP: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:18:GLU:OE1	1:A:22:HIS:HA	2.09	0.52
1:A:42:THR:CG2	2:A:104:COA:H142	2.40	0.52
1:A:38:ASP:HA	1:A:73:ILE:O	2.10	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:23:GLU:CD	1:A:48:LYS:HZ3	2.12	0.50
1:A:24:ALA:CB	1:A:53:LEU:HD21	2.41	0.50
1:A:7:LYS:HE2	1:A:9:VAL:HG13	1.92	0.50
1:A:65:ALA:HB2	1:A:72:ILE:HD11	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:37:MET:HB3	1:A:72:ILE:HG13	1.92	0.50
1:A:46:SER:OG	1:A:49:ARG:NH2	2.45	0.50
1:A:45:PRO:HB2	3:A:124:HOH:O	2.11	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:7:LYS:HB3	3:A:149:HOH:O	2.12	0.50
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: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:54:ALA:O	1:A:58:CYS:HB3	2.12	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: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:54:ALA:HB3	1:A:78:TYR:CZ	2.48	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:54:ALA:HB1	2:A:104:COA:CEP	2.39	0.48
1:A:45:PRO:CG	1:A:48:LYS:HD3	2.44	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:15:ARG:HD2	3:A:106:HOH:O	2.13	0.47
1:A:50:GLY:C	1:A:51:LEU:HD23	2.34	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: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
1:A:47:PHE:CG	1:A:47:PHE:O	2.67	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:84:LEU:N	1:A:85:PRO:CD	2.78	0.46
1:A:25:PHE:C	1:A:57:LEU:HD13	2.36	0.46
1:A:8:ILE:HD11	1:A:56:HIS:HB3	1.97	0.46
1:A:25:PHE:CZ	1:A:41:HIS:CD2	3.03	0.46
1:A:78:TYR:O	1:A:79:VAL:C	2.49	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
1:A:80:SER:OG	1:A:81:ASP:N	2.48	0.46
1:A:32:ASN:ND2	1:A:35:LYS:HB2	2.26	0.46
1:A:19:THR:CG2	1:A:24:ALA:HB3	2.46	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:25:PHE:HE1	1:A:27:GLU:OE2	2.00	0.45
1:A:45:PRO:CG	1:A:48:LYS:CD	2.82	0.45
1:A:84:LEU:N	1:A:85:PRO:CD	2.80	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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:CE2	1:A:12:GLU:HG2	2.53	0.44
1:A:10:TRP:CZ2	1:A:12:GLU:HG2	2.53	0.43
2:A:104:COA:CDP	2:A:104:COA:HN8	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
1:A:59:VAL:O	1:A:59:VAL:CG1	2.67	0.43
1:A:46:SER:O	1:A:47:PHE:C	2.55	0.43
1:A:15:ARG:HB3	1:A:28:TYR:O	2.18	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:6:PRO:HB2	1:A:7:LYS:H	1.44	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:84:LEU:CB	1:A:85:PRO:HD3	2.48	0.42
1:A:16:ARG:NH2	1:A:27:GLU:CD	2.73	0.42
1:A:41:HIS:CE1	3:A:163:HOH:O	2.71	0.42
1:A:49:ARG:O	1:A:51:LEU:HG	2.19	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: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:15:ARG:HA	1:A:28:TYR:OH	2.20	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:77:SER:O	1:A:81:ASP:HB2	2.19	0.42
1:A:65:ALA:CB	1:A:72:ILE:HD12	2.50	0.42
1:A:20:GLU:C	1:A:22:HIS:N	2.73	0.42
1:A:48:LYS:HE2	1:A:53:LEU:CD2	2.50	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:36:VAL:HG11	1:A:73:ILE:HD11	2.02	0.41
1:A:39:LEU:CD1	1:A:79:VAL:HG21	2.50	0.41
1:A:91:LYS:NZ	3:A:139:HOH:O	2.50	0.41
1:A:54:ALA:HA	1:A:57:LEU:HD12	2.01	0.41
1:A:82:THR:HG21	2:A:104:COA:H1B	2.02	0.41
1:A:47:PHE:HD1	1:A:47:PHE:H	1.57	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:49:ARG:HG2	2:A:104:COA:H121	2.03	0.41
1:A:20:GLU:HG3	3:A:165:HOH:O	2.20	0.41
1:A:10:TRP:CE2	1:A:12:GLU:HG2	2.56	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: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
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:80:SER:HA	1:A:84:LEU:HD12	2.03	0.40
1:A:65:ALA:HA	1:A:70:ILE:HB	2.03	0.40
1:A:78:TYR:HA	2:A:104:COA:H2A	2.03	0.40
1:A:65:ALA:CB	1:A:93:LEU:HD11	2.52	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:44:VAL:HG23	2:A:104:COA:H122	2.03	0.40
1:A:59:VAL:O	1:A:63:GLU:CG	2.68	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	1-A	89/103 (86%)	79 (89%)	6 (7%)	4 (4%)	2	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%)	14	5

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
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%)	49	43
1	2-A	82/93 (88%)	75 (92%)	7 (8%)	10	4
1	3-A	82/93 (88%)	76 (93%)	6 (7%)	14	6
1	4-A	82/93 (88%)	78 (95%)	4 (5%)	25	17
All	All	328/372 (88%)	309 (94%)	19 (6%)	20	11

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

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Mol	Chain	Res	Type
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 [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

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	-	41,50,50	2.37	13 (31%)	52,75,75	2.79	22 (42%)
2	COA	4-A	104	-	41,50,50	3.84	20 (48%)	52,75,75	3.32	24 (46%)
2	COA	3-A	104	-	41,50,50	2.82	18 (43%)	52,75,75	2.85	25 (48%)
2	COA	2-A	104	-	41,50,50	3.28	23 (56%)	52,75,75	2.43	18 (34%)

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	-	-	6/44/64/64	0/3/3/3
2	COA	4-A	104	-	-	10/44/64/64	0/3/3/3
2	COA	3-A	104	-	-	8/44/64/64	0/3/3/3
2	COA	2-A	104	-	-	6/44/64/64	0/3/3/3

All (74) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	2-A	104	COA	C2B-C1B	11.11	1.70	1.53
2	4-A	104	COA	C2B-C1B	10.05	1.69	1.53
2	1-A	104	COA	C2B-C1B	8.31	1.66	1.53
2	4-A	104	COA	C9P-N8P	8.10	1.51	1.33
2	3-A	104	COA	C2B-C1B	8.06	1.66	1.53
2	4-A	104	COA	C3P-N4P	-7.57	1.28	1.46
2	4-A	104	COA	C5P-N4P	-7.29	1.16	1.33
2	4-A	104	COA	P3B-O3B	6.74	1.72	1.59
2	2-A	104	COA	C7P-N8P	6.68	1.61	1.46
2	3-A	104	COA	O5B-C5B	-6.30	1.20	1.44
2	4-A	104	COA	O9P-C9P	6.09	1.35	1.23
2	3-A	104	COA	C5P-N4P	5.50	1.45	1.33
2	4-A	104	COA	O6A-CCP	5.48	1.61	1.43
2	2-A	104	COA	C2A-N1A	-5.46	1.23	1.33
2	1-A	104	COA	P1A-O5B	5.29	1.80	1.59
2	3-A	104	COA	O4B-C1B	-5.17	1.33	1.41
2	3-A	104	COA	OAP-CAP	-5.17	1.32	1.42
2	4-A	104	COA	P1A-O1A	5.13	1.69	1.50
2	4-A	104	COA	C2P-S1P	-5.11	1.63	1.80
2	4-A	104	COA	C6P-C5P	5.03	1.60	1.51
2	4-A	104	COA	O4B-C4B	-4.76	1.34	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	2-A	104	COA	O5B-C5B	-4.64	1.26	1.44
2	1-A	104	COA	O9P-C9P	-4.55	1.14	1.23
2	2-A	104	COA	P1A-O5B	-4.55	1.40	1.59
2	2-A	104	COA	C2B-C3B	4.55	1.63	1.52
2	4-A	104	COA	P2A-O5A	4.54	1.76	1.55
2	2-A	104	COA	C9P-N8P	4.49	1.43	1.33
2	2-A	104	COA	OAP-CAP	-4.44	1.34	1.42
2	2-A	104	COA	O6A-CCP	4.02	1.56	1.43
2	3-A	104	COA	P1A-O5B	-3.98	1.43	1.59
2	4-A	104	COA	C5B-C4B	3.87	1.63	1.51
2	3-A	104	COA	C3P-N4P	3.68	1.54	1.46
2	2-A	104	COA	P3B-O8A	3.67	1.69	1.54
2	2-A	104	COA	O4B-C1B	3.64	1.46	1.41
2	1-A	104	COA	C7P-N8P	-3.62	1.37	1.46
2	1-A	104	COA	P3B-O3B	-3.59	1.52	1.59
2	1-A	104	COA	O4B-C1B	-3.57	1.36	1.41
2	4-A	104	COA	CEP-CBP	3.51	1.61	1.53
2	4-A	104	COA	C5A-C4A	3.44	1.50	1.40
2	2-A	104	COA	C3B-C4B	3.38	1.62	1.52
2	3-A	104	COA	O9P-C9P	3.35	1.30	1.23
2	2-A	104	COA	C6A-N6A	3.09	1.45	1.34
2	3-A	104	COA	P3B-O9A	3.04	1.66	1.54
2	4-A	104	COA	P1A-O5B	2.94	1.71	1.59
2	3-A	104	COA	C6A-N6A	2.92	1.44	1.34
2	3-A	104	COA	O6A-CCP	2.91	1.53	1.43
2	2-A	104	COA	P2A-O5A	-2.88	1.41	1.55
2	3-A	104	COA	C2A-N3A	2.86	1.36	1.32
2	1-A	104	COA	C2B-C3B	2.77	1.59	1.52
2	3-A	104	COA	P1A-O1A	-2.71	1.41	1.50
2	2-A	104	COA	P3B-O9A	2.66	1.65	1.54
2	4-A	104	COA	C3B-C4B	2.64	1.60	1.52
2	3-A	104	COA	C9P-N8P	2.63	1.39	1.33
2	1-A	104	COA	O4B-C4B	2.52	1.50	1.45
2	3-A	104	COA	C2P-S1P	2.50	1.89	1.80
2	1-A	104	COA	C3P-N4P	-2.50	1.40	1.46
2	1-A	104	COA	P3B-O9A	2.41	1.64	1.54
2	3-A	104	COA	P2A-O5A	-2.37	1.44	1.55
2	2-A	104	COA	C6A-C5A	2.33	1.51	1.43
2	1-A	104	COA	C9P-N8P	-2.33	1.28	1.33
2	4-A	104	COA	P3B-O8A	2.30	1.63	1.54
2	2-A	104	COA	C8A-N7A	-2.27	1.30	1.34
2	1-A	104	COA	O2B-C2B	2.26	1.48	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	1-A	104	COA	CDP-CBP	-2.22	1.48	1.53
2	3-A	104	COA	P1A-O2A	-2.20	1.45	1.55
2	2-A	104	COA	P1A-O2A	-2.17	1.45	1.55
2	4-A	104	COA	OAP-CAP	-2.16	1.38	1.42
2	3-A	104	COA	C7P-N8P	-2.15	1.41	1.46
2	2-A	104	COA	C6P-C5P	-2.14	1.47	1.51
2	2-A	104	COA	C2P-S1P	2.12	1.88	1.80
2	2-A	104	COA	C5A-N7A	2.11	1.47	1.39
2	4-A	104	COA	P2A-O6A	2.10	1.67	1.59
2	2-A	104	COA	P2A-O6A	2.04	1.67	1.59
2	2-A	104	COA	P2A-O4A	2.02	1.58	1.50

All (89) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	4-A	104	COA	O6A-CCP-CBP	10.00	126.62	110.55
2	4-A	104	COA	O9A-P3B-O8A	-9.24	72.32	107.64
2	1-A	104	COA	P2A-O3A-P1A	-7.84	105.94	132.83
2	4-A	104	COA	CEP-CBP-CCP	7.36	120.23	108.23
2	4-A	104	COA	CEP-CBP-CAP	-7.00	96.68	108.82
2	1-A	104	COA	O9A-P3B-O7A	-6.87	83.80	110.68
2	1-A	104	COA	O9A-P3B-O8A	-6.59	82.44	107.64
2	2-A	104	COA	O9A-P3B-O3B	-6.56	76.61	105.99
2	3-A	104	COA	O9A-P3B-O7A	-6.01	87.16	110.68
2	3-A	104	COA	O9A-P3B-O8A	-5.77	85.57	107.64
2	4-A	104	COA	P2A-O3A-P1A	-5.76	113.06	132.83
2	2-A	104	COA	CDP-CBP-CCP	5.74	117.59	108.23
2	3-A	104	COA	C7P-N8P-C9P	5.48	132.37	122.59
2	2-A	104	COA	O9A-P3B-O8A	-5.46	86.77	107.64
2	3-A	104	COA	N3A-C2A-N1A	-5.44	120.18	128.68
2	3-A	104	COA	O6A-CCP-CBP	-5.40	101.86	110.55
2	1-A	104	COA	C6P-C7P-N8P	-5.36	101.08	111.90
2	4-A	104	COA	O4B-C4B-C3B	5.36	116.35	104.87
2	4-A	104	COA	O9A-P3B-O7A	-5.31	89.89	110.68
2	4-A	104	COA	O9A-P3B-O3B	-5.26	82.44	105.99
2	2-A	104	COA	O9A-P3B-O7A	-5.19	90.36	110.68
2	3-A	104	COA	O4B-C4B-C3B	5.09	115.78	104.87
2	1-A	104	COA	O9A-P3B-O3B	-4.86	84.20	105.99
2	3-A	104	COA	O9A-P3B-O3B	-4.72	84.85	105.99
2	1-A	104	COA	CEP-CBP-CAP	-4.69	100.68	108.82
2	2-A	104	COA	CEP-CBP-CCP	-4.62	100.70	108.23
2	1-A	104	COA	N3A-C2A-N1A	-4.55	121.57	128.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	3-A	104	COA	CAP-C9P-N8P	-4.42	107.77	116.58
2	2-A	104	COA	O8A-P3B-O3B	4.25	125.02	105.99
2	2-A	104	COA	O4B-C4B-C3B	4.19	113.85	104.87
2	2-A	104	COA	O3B-P3B-O7A	4.19	125.57	109.39
2	1-A	104	COA	O3B-P3B-O7A	4.17	125.50	109.39
2	3-A	104	COA	CDP-CBP-CAP	-4.12	101.68	108.82
2	1-A	104	COA	O5B-C5B-C4B	-4.07	94.99	108.99
2	3-A	104	COA	P2A-O3A-P1A	-4.05	118.91	132.83
2	1-A	104	COA	C4A-C5A-N7A	-3.96	105.27	109.40
2	4-A	104	COA	N3A-C2A-N1A	-3.84	122.68	128.68
2	3-A	104	COA	C7P-C6P-C5P	-3.72	106.16	112.36
2	1-A	104	COA	O9P-C9P-N8P	-3.70	115.05	122.99
2	3-A	104	COA	C1B-N9A-C4A	-3.69	120.15	126.64
2	3-A	104	COA	O5B-C5B-C4B	-3.67	96.36	108.99
2	4-A	104	COA	O9P-C9P-CAP	-3.67	109.87	121.06
2	2-A	104	COA	N3A-C2A-N1A	-3.46	123.28	128.68
2	3-A	104	COA	O8A-P3B-O3B	3.36	121.05	105.99
2	4-A	104	COA	O5P-C5P-C6P	3.32	128.09	122.02
2	2-A	104	COA	O6A-P2A-O4A	3.17	121.44	109.07
2	3-A	104	COA	OAP-CAP-CBP	3.15	117.67	110.25
2	3-A	104	COA	CEP-CBP-CCP	-3.15	103.10	108.23
2	4-A	104	COA	C6P-C7P-N8P	3.11	118.18	111.90
2	1-A	104	COA	OAP-CAP-CBP	-3.11	102.94	110.25
2	4-A	104	COA	C3P-N4P-C5P	-3.07	117.13	122.84
2	4-A	104	COA	O9P-C9P-N8P	3.06	129.56	122.99
2	4-A	104	COA	O5P-C5P-N4P	-2.98	117.39	123.01
2	4-A	104	COA	O5B-P1A-O1A	2.91	120.45	109.07
2	4-A	104	COA	C2B-C3B-C4B	-2.87	98.13	103.22
2	2-A	104	COA	O5B-C5B-C4B	-2.84	99.21	108.99
2	3-A	104	COA	C2A-N1A-C6A	2.82	123.58	118.75
2	3-A	104	COA	O3B-P3B-O7A	2.81	120.25	109.39
2	3-A	104	COA	CDP-CBP-CCP	-2.81	103.65	108.23
2	1-A	104	COA	C3B-C2B-C1B	2.76	106.00	99.89
2	4-A	104	COA	O8A-P3B-O3B	2.76	118.34	105.99
2	1-A	104	COA	O6A-CCP-CBP	-2.75	106.12	110.55
2	3-A	104	COA	C4A-C5A-N7A	-2.74	106.54	109.40
2	3-A	104	COA	C5A-C6A-N6A	2.67	124.41	120.35
2	2-A	104	COA	O9P-C9P-CAP	-2.65	112.97	121.06
2	1-A	104	COA	O4B-C4B-C3B	2.64	110.53	104.87
2	2-A	104	COA	O9P-C9P-N8P	2.64	128.67	122.99
2	1-A	104	COA	O8A-P3B-O7A	2.62	120.93	110.68
2	4-A	104	COA	C2P-C3P-N4P	-2.60	106.36	112.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	3-A	104	COA	C6P-C5P-N4P	2.56	120.73	116.42
2	1-A	104	COA	CAP-C9P-N8P	2.56	121.67	116.58
2	4-A	104	COA	OAP-CAP-CBP	2.50	116.13	110.25
2	1-A	104	COA	C3P-N4P-C5P	-2.47	118.25	122.84
2	1-A	104	COA	C2P-C3P-N4P	-2.41	106.79	112.31
2	4-A	104	COA	C3B-C2B-C1B	2.39	105.19	99.89
2	4-A	104	COA	O3B-P3B-O7A	2.32	118.35	109.39
2	2-A	104	COA	O5B-P1A-O1A	2.28	117.99	109.07
2	1-A	104	COA	O4B-C4B-C5B	2.27	116.86	109.37
2	4-A	104	COA	O4B-C1B-C2B	-2.27	103.61	106.93
2	1-A	104	COA	P1A-O5B-C5B	-2.26	108.43	121.68
2	2-A	104	COA	O3B-C3B-C4B	2.25	118.23	110.08
2	2-A	104	COA	CEP-CBP-CAP	-2.24	104.94	108.82
2	3-A	104	COA	O5P-C5P-C6P	-2.13	118.13	122.02
2	2-A	104	COA	O2B-C2B-C1B	2.12	118.67	110.85
2	4-A	104	COA	P2A-O6A-CCP	-2.12	109.36	121.56
2	1-A	104	COA	CEP-CBP-CDP	2.08	113.40	109.17
2	3-A	104	COA	CEP-CBP-CDP	2.01	113.27	109.17
2	2-A	104	COA	P2A-O3A-P1A	-2.01	125.92	132.83
2	3-A	104	COA	O6A-P2A-O4A	2.00	116.90	109.07

There are no chirality outliers.

All (30) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	1-A	104	COA	C5B-O5B-P1A-O3A
2	1-A	104	COA	O9P-C9P-CAP-CBP
2	1-A	104	COA	N8P-C9P-CAP-CBP
2	4-A	104	COA	C2B-C3B-O3B-P3B
2	4-A	104	COA	O9P-C9P-CAP-CBP
2	4-A	104	COA	N8P-C9P-CAP-CBP
2	4-A	104	COA	O9P-C9P-CAP-OAP
2	4-A	104	COA	N8P-C9P-CAP-OAP
2	3-A	104	COA	C5B-O5B-P1A-O3A
2	2-A	104	COA	O9P-C9P-CAP-CBP
2	2-A	104	COA	N8P-C9P-CAP-CBP
2	4-A	104	COA	O4B-C4B-C5B-O5B
2	3-A	104	COA	O4B-C4B-C5B-O5B
2	2-A	104	COA	O9P-C9P-CAP-OAP
2	3-A	104	COA	O9P-C9P-CAP-CBP
2	3-A	104	COA	N8P-C9P-CAP-CBP
2	2-A	104	COA	N8P-C9P-CAP-OAP

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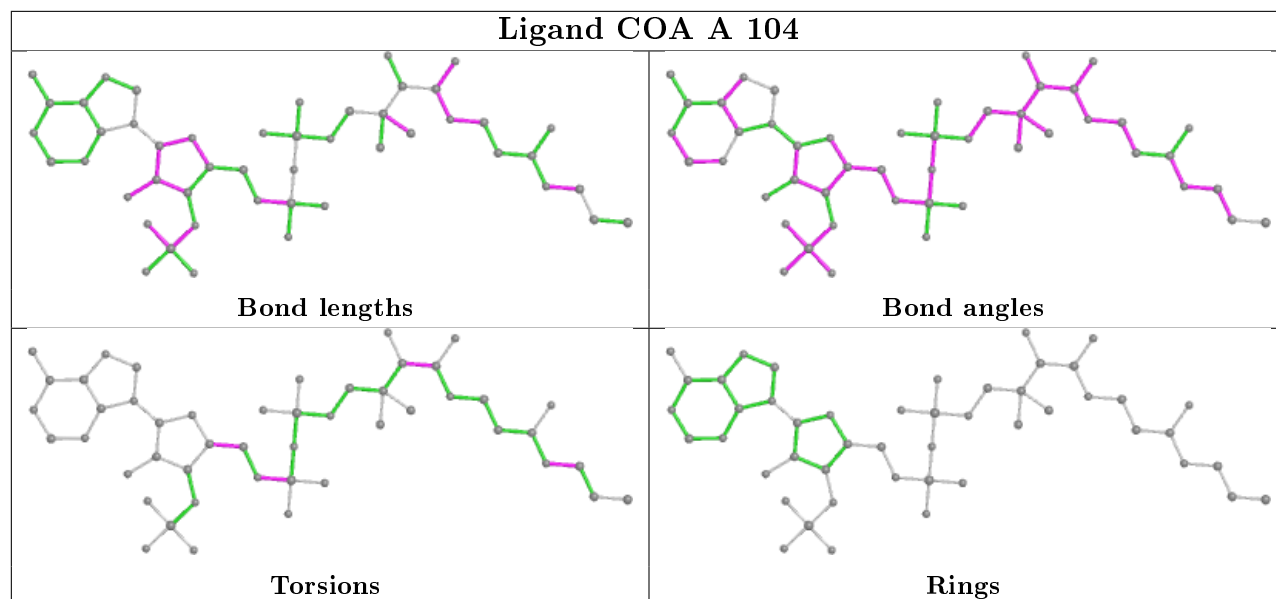
Mol	Chain	Res	Type	Atoms
2	1-A	104	COA	C5B-O5B-P1A-O2A
2	3-A	104	COA	C5B-O5B-P1A-O2A
2	1-A	104	COA	C2P-C3P-N4P-C5P
2	3-A	104	COA	C2P-C3P-N4P-C5P
2	1-A	104	COA	O4B-C4B-C5B-O5B
2	4-A	104	COA	C2P-C3P-N4P-C5P
2	3-A	104	COA	C3B-O3B-P3B-O7A
2	3-A	104	COA	N8P-C9P-CAP-OAP
2	2-A	104	COA	C2P-C3P-N4P-C5P
2	4-A	104	COA	C5B-O5B-P1A-O3A
2	4-A	104	COA	C4B-C3B-O3B-P3B
2	4-A	104	COA	CCP-O6A-P2A-O4A
2	2-A	104	COA	CCP-O6A-P2A-O4A

There are no ring outliers.

4 monomers are involved in 80 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	1-A	104	COA	24	0
2	4-A	104	COA	22	0
2	3-A	104	COA	21	0
2	2-A	104	COA	13	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [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	1-A	91/103 (88%)	-0.08	1 (1%) 80 81	18, 26, 42, 74	91 (100%)
1	2-A	91/103 (88%)	-0.08	1 (1%) 80 81	18, 26, 42, 74	91 (100%)
1	3-A	91/103 (88%)	-0.08	1 (1%) 80 81	18, 26, 42, 74	91 (100%)
1	4-A	91/103 (88%)	-0.08	1 (1%) 80 81	18, 26, 42, 74	91 (100%)
All	All	364/412 (88%)	-0.08	4 (1%) 62 81	18, 27, 47, 74	364 (100%)

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	1-A	5	PRO	8.5
1	2-A	5	PRO	8.5
1	3-A	5	PRO	8.5
1	4-A	5	PRO	8.5

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

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

6.3 Carbohydrates [i](#)

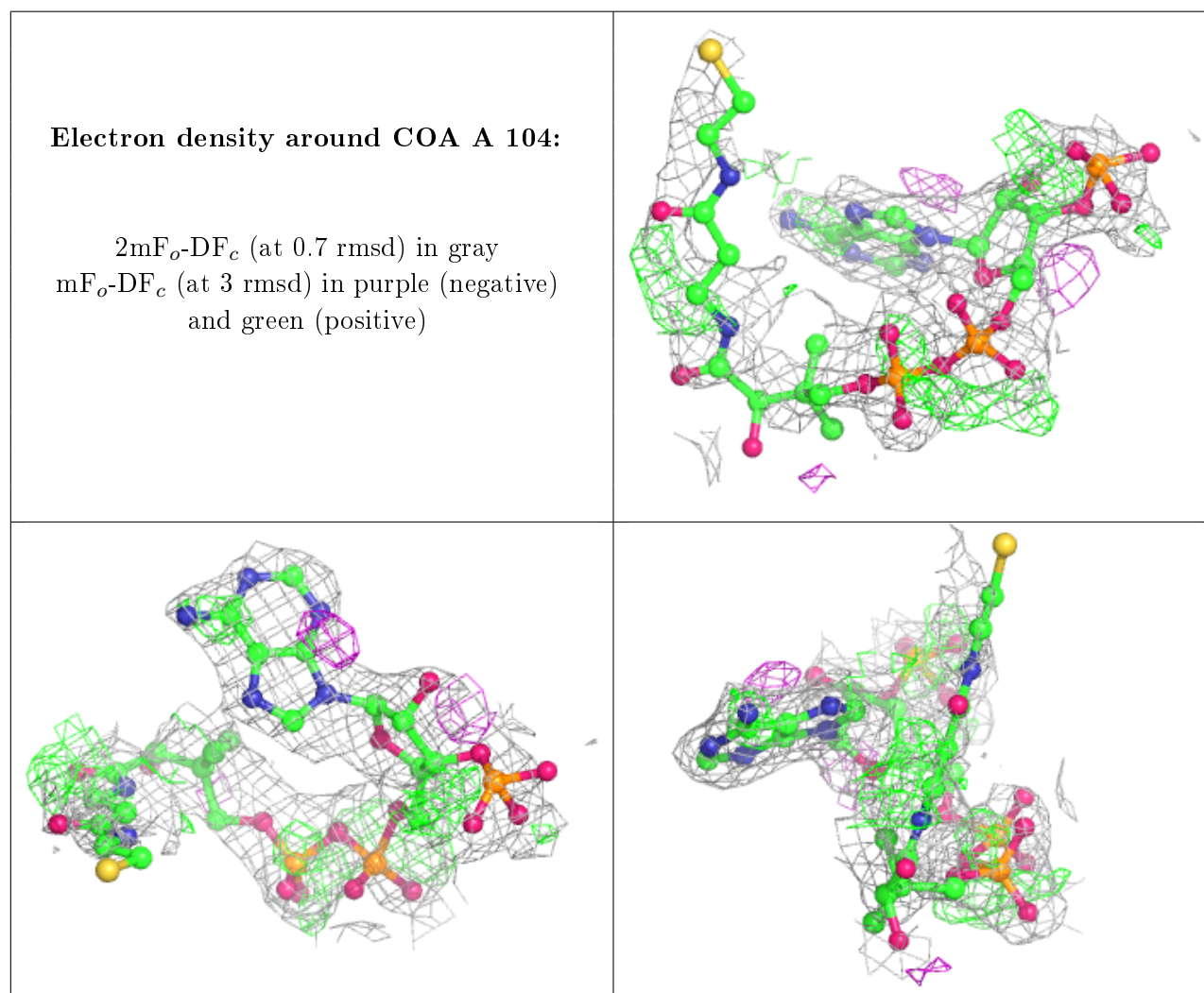
There are no carbohydrates in this entry.

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(\AA^2)	Q<0.9
2	COA	1-A	104	48/48	0.73	0.41	19,34,50,57	48
2	COA	4-A	104	48/48	0.73	0.41	24,42,53,57	48
2	COA	3-A	104	48/48	0.73	0.41	9,40,58,59	48
2	COA	2-A	104	48/48	0.73	0.41	12,43,56,60	48

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers [i](#)

There are no such residues in this entry.