



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 22, 2016 – 06:29 PM EDT

PDB ID : 5HMN  
Title : Crystal structure of an aminoglycoside acetyltransferase HMB0005 from an uncultured soil metagenomic sample, unknown active site density modeled as polyethylene glycol  
Authors : Xu, Z.; Stogios, P.J.; Wawrzak, Z.; Skarina, T.; Yim, V.; Savchenko, A.; Anderson, W.F.; Center for Structural Genomics of Infectious Diseases (CSGID)  
Deposited on : 2016-01-16  
Resolution : 2.02 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7.1 (RC1), CSD as537be (2016)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20027107  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0122  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20027107

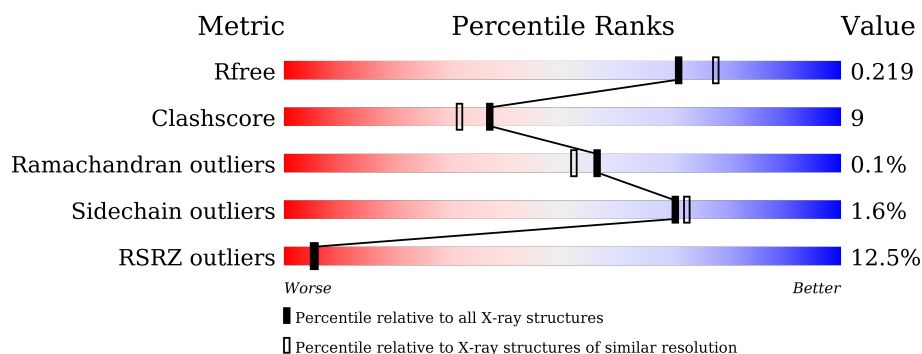
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.02 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))
$R_{free}$	91344	7858 (2.04-2.00)
Clashscore	102246	9060 (2.04-2.00)
Ramachandran outliers	100387	8952 (2.04-2.00)
Sidechain outliers	100360	8951 (2.04-2.00)
RSRZ outliers	91569	7873 (2.04-2.00)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	159	<div> <div>87%</div> <div>8%</div> <div>5%</div> </div>
1	B	159	<div>6%</div> <div>87%</div> <div>12%</div> <div>.</div>
1	C	159	<div>9%</div> <div>80%</div> <div>16%</div> <div>.</div>
1	D	159	<div>12%</div> <div>82%</div> <div>13%</div> <div>.</div>
1	E	159	<div>19%</div> <div>74%</div> <div>21%</div> <div>5%</div>
1	F	159	<div>26%</div> <div>65%</div> <div>28%</div> <div>.</div> <div>5%</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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	COA	B	500	-	-	-	X
3	PG4	A	202[A]	-	-	X	X
3	PG4	A	202[B]	-	-	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8407 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.

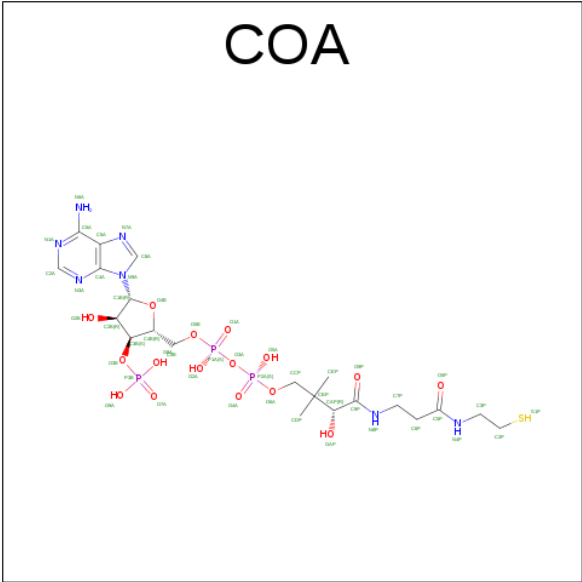
- Molecule 1 is a protein called AAC3-I.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	151	Total	C	N	O	S	0	3	0
			1207	774	203	228	2			
1	B	159	Total	C	N	O	S	0	4	0
			1270	814	211	242	3			
1	C	152	Total	C	N	O	S	0	1	0
			1201	770	200	228	3			
1	D	154	Total	C	N	O	S	0	1	0
			1218	780	203	232	3			
1	E	151	Total	C	N	O	S	0	1	0
			1193	763	199	229	2			
1	F	151	Total	C	N	O	S	0	0	0
			1188	760	199	227	2			

There are 12 discrepancies between the modelled and reference sequences:

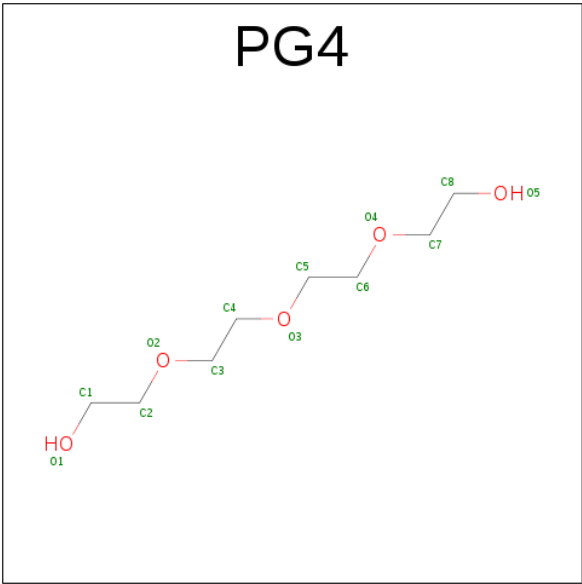
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLN	-	expression tag	UNP A0A059WV44
A	0	SER	-	expression tag	UNP A0A059WV44
B	-1	GLN	-	expression tag	UNP A0A059WV44
B	0	SER	-	expression tag	UNP A0A059WV44
C	-1	GLN	-	expression tag	UNP A0A059WV44
C	0	SER	-	expression tag	UNP A0A059WV44
D	-1	GLN	-	expression tag	UNP A0A059WV44
D	0	SER	-	expression tag	UNP A0A059WV44
E	-1	GLN	-	expression tag	UNP A0A059WV44
E	0	SER	-	expression tag	UNP A0A059WV44
F	-1	GLN	-	expression tag	UNP A0A059WV44
F	0	SER	-	expression tag	UNP A0A059WV44

- Molecule 2 is COENZYME A (three-letter code: COA) (formula: C<sub>21</sub>H<sub>36</sub>N<sub>7</sub>O<sub>16</sub>P<sub>3</sub>S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	S	0	1
			96	42	14	32	6	2		
2	B	1	Total	C	N	O	P	S	0	0
			48	21	7	16	3	1		
2	C	1	Total	C	N	O	P	S	0	1
			96	42	14	32	6	2		
2	E	1	Total	C	N	O	P	S	0	0
			48	21	7	16	3	1		
2	F	1	Total	C	N	O	P	S	0	0
			48	21	7	16	3	1		

- Molecule 3 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: C<sub>8</sub>H<sub>18</sub>O<sub>5</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	1
			26	16	10		

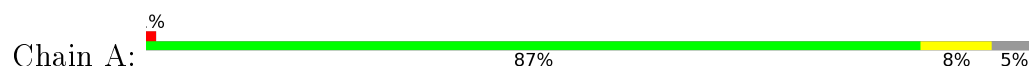
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	176	Total	O	0	10
			186	186		
4	B	185	Total	O	0	9
			194	194		
4	C	143	Total	O	0	2
			145	145		
4	D	109	Total	O	0	2
			111	111		
4	E	84	Total	O	0	0
			84	84		
4	F	47	Total	O	0	1
			48	48		

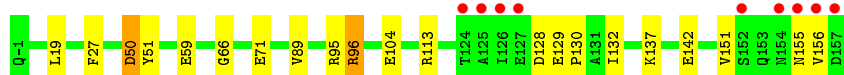
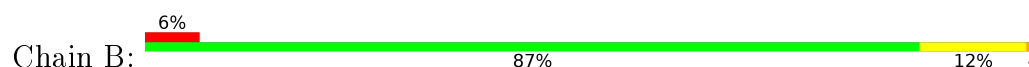
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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 ( $\text{RSRZ} > 2$ ). 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.

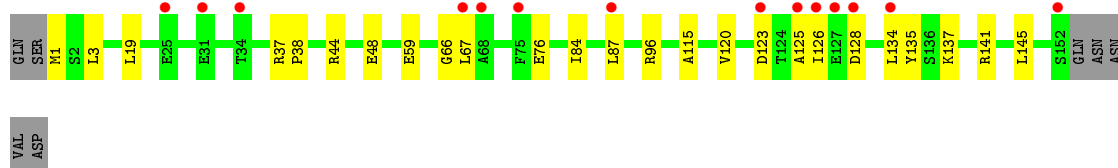
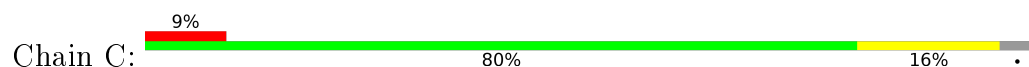
#### • Molecule 1: AAC3-I



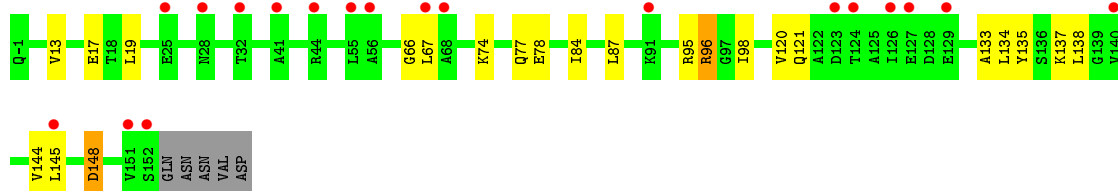
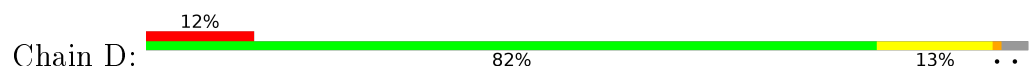
#### • Molecule 1: AAC3-I



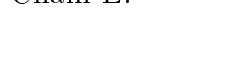
#### • Molecule 1: AAC3-I

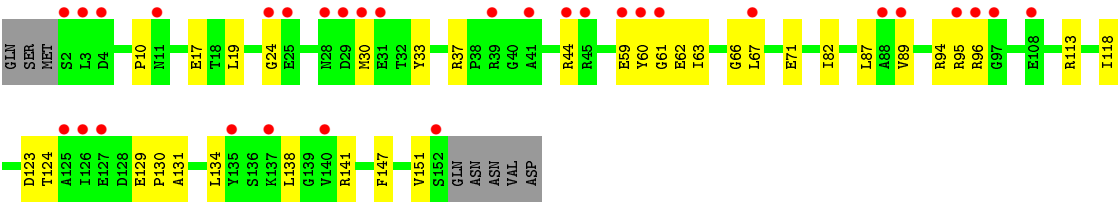


#### • Molecule 1: AAC3-I

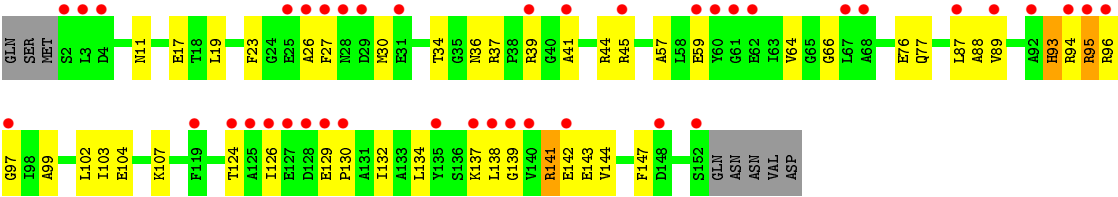


#### • Molecule 1: AAC3-I





• Molecule 1: AAC3-I





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	58.27Å 62.76Å 253.61Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.87 – 2.02 19.87 – 2.02	Depositor EDS
% Data completeness (in resolution range)	99.6 (19.87-2.02) 96.7 (19.87-2.02)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.41 (at 2.02Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, $R_{free}$	0.191 , 0.221 0.190 , 0.219	Depositor DCC
$R_{free}$ test set	1920 reflections (3.19%)	DCC
Wilson B-factor (Å <sup>2</sup> )	34.9	Xtriage
Anisotropy	0.137	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 55.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 62029 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	8407	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 19.78% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: COA, PG4

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	0.43	0/1237	0.51	0/1672
1	B	0.40	0/1303	0.52	0/1763
1	C	0.28	0/1225	0.47	0/1656
1	D	0.42	0/1243	0.58	0/1680
1	E	0.38	0/1217	0.53	0/1646
1	F	0.48	0/1209	0.66	0/1635
All	All	0.40	0/7434	0.55	0/10052

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1207	0	1214	12	0
1	B	1270	0	1282	16	0
1	C	1201	0	1205	19	0
1	D	1218	0	1212	18	0
1	E	1193	0	1186	34	0
1	F	1188	0	1182	50	0
2	A	96	0	62	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	48	0	32	8	0
2	C	96	0	62	1	0
2	E	48	0	31	11	0
2	F	48	0	32	16	0
3	A	26	0	36	7	0
4	A	186	0	0	1	0
4	B	194	0	0	6	0
4	C	145	0	0	0	0
4	D	111	0	0	0	0
4	E	84	0	0	2	0
4	F	48	0	0	0	0
All	All	8407	0	7536	142	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 9.

All (142) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:95:ARG:NH1	2:F:500:COA:O8A	1.65	1.27
2:B:500:COA:O4B	2:B:500:COA:C1B	1.65	1.21
2:F:500:COA:O4B	2:F:500:COA:C1B	1.65	1.16
1:C:135:TYR:HH	2:C:500[A]:COA:HS1	0.98	0.90
1:E:94:ARG:NH1	2:E:500:COA:H72	1.87	0.90
1:F:95:ARG:HD3	2:F:500:COA:O8A	1.74	0.87
1:E:95:ARG:NH2	2:E:500:COA:O9A	2.07	0.87
3:A:202[A]:PG4:H31	3:A:202[A]:PG4:H61	1.59	0.85
1:E:17:GLU:HB3	1:E:37:ARG:NH1	1.91	0.84
1:F:76:GLU:HG2	1:F:77:GLN:HG3	1.65	0.79
1:C:59:GLU:OE2	1:C:96:ARG:NH2	2.18	0.77
2:B:500:COA:S1P	4:B:613:HOH:O	2.43	0.76
3:A:202[A]:PG4:H61	3:A:202[A]:PG4:C3	2.15	0.75
1:F:124:THR:HG22	1:F:124:THR:O	1.84	0.75
1:A:85:TYR:O	3:A:202[A]:PG4:H32	1.88	0.74
1:F:141:ARG:O	1:F:142:GLU:HG3	1.88	0.73
1:E:17:GLU:HB3	1:E:37:ARG:HH11	1.53	0.73
1:A:85:TYR:O	3:A:202[A]:PG4:C3	2.37	0.72
1:B:59:GLU:OE1	1:B:96:ARG:NH1	2.20	0.70
1:B:19:LEU:HD11	1:B:66:GLY:HA3	1.74	0.70
1:B:130:PRO:HB3	2:B:500:COA:N6A	2.08	0.68
1:F:95:ARG:HH11	2:F:500:COA:P3B	2.15	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:134:LEU:O	1:C:137:LYS:HG2	1.96	0.65
1:E:19:LEU:HD11	1:E:66:GLY:HA3	1.78	0.64
1:F:19:LEU:HD11	1:F:66:GLY:HA3	1.78	0.64
1:F:129:GLU:N	1:F:130:PRO:HD2	2.13	0.64
1:D:95:ARG:NH2	1:E:113:ARG:O	2.31	0.63
1:D:67:LEU:HD12	1:D:87:LEU:HD12	1.81	0.61
1:F:17:GLU:O	1:F:37:ARG:NH2	2.33	0.60
1:F:94:ARG:O	1:F:95:ARG:HB2	2.02	0.60
1:B:89:VAL:HG23	2:B:500:COA:H132	1.85	0.59
1:F:124:THR:CG2	1:F:124:THR:O	2.49	0.59
1:B:137:LYS:NZ	4:B:608:HOH:O	2.35	0.59
1:D:148:ASP:O	1:F:139:GLY:HA3	2.02	0.59
1:F:96:ARG:O	1:F:96:ARG:HG2	2.03	0.59
1:F:89:VAL:CG2	2:F:500:COA:H142	2.33	0.59
1:E:10:PRO:O	1:E:44:ARG:HG3	2.02	0.59
1:E:71:GLU:OE1	1:E:113:ARG:NH1	2.34	0.58
1:A:19:LEU:HD11	1:A:66:GLY:HA3	1.86	0.57
1:C:44:ARG:NH1	1:C:48:GLU:OE1	2.38	0.57
2:E:500:COA:O5P	4:E:601:HOH:O	2.17	0.57
1:E:95:ARG:NH2	2:E:500:COA:P3B	2.78	0.56
1:F:95:ARG:NH2	2:F:500:COA:C2A	2.69	0.56
1:E:124:THR:HG22	1:E:124:THR:O	2.06	0.56
1:F:129:GLU:O	1:F:132:ILE:N	2.39	0.56
1:E:67:LEU:HD12	1:E:87:LEU:HD13	1.88	0.55
1:B:130:PRO:HB3	2:B:500:COA:H62A	1.70	0.55
1:A:36:ASN:ND2	4:B:606:HOH:O	2.33	0.55
1:F:57:ALA:HB2	1:F:102:LEU:HD21	1.89	0.54
1:F:89:VAL:HG21	2:F:500:COA:H142	1.90	0.54
1:F:27:PHE:HE1	2:F:500:COA:C6P	2.21	0.54
1:C:67:LEU:HB2	1:C:87[B]:LEU:HD23	1.89	0.54
1:E:95:ARG:HH22	2:E:500:COA:P3B	2.30	0.54
1:A:85:TYR:O	3:A:202[A]:PG4:H31	2.07	0.53
1:A:141:ARG:NH1	4:A:306:HOH:O	2.33	0.52
1:C:134:LEU:HD12	1:C:137:LYS:HE2	1.92	0.52
1:C:59:GLU:CD	1:C:96:ARG:HH22	2.13	0.52
1:E:94:ARG:O	1:E:95:ARG:HB2	2.10	0.51
1:B:95:ARG:N	2:B:500:COA:O5A	2.39	0.51
1:D:13:VAL:O	1:D:17:GLU:HG3	2.11	0.51
1:F:97:GLY:N	2:F:500:COA:O1A	2.38	0.51
1:F:104:GLU:HA	1:F:107:LYS:HG2	1.92	0.51
1:D:96:ARG:HD3	1:D:98:ILE:HG13	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:59:GLU:O	1:E:60:TYR:HB2	2.11	0.51
1:E:61:GLY:O	1:E:62:GLU:HG2	2.10	0.51
1:F:99:ALA:HB1	1:F:134:LEU:HD21	1.94	0.50
1:F:137:LYS:C	1:F:138:LEU:HD12	2.32	0.50
1:F:94:ARG:HD2	2:F:500:COA:H10	1.94	0.50
1:C:123:ASP:HB3	1:C:128:ASP:OD1	2.12	0.50
1:C:19:LEU:HD11	1:C:66:GLY:HA3	1.94	0.49
1:F:93:HIS:ND1	1:F:93:HIS:N	2.60	0.49
1:F:34:THR:O	1:F:34:THR:HG22	2.12	0.49
1:E:82:ILE:HB	1:E:118:ILE:HG12	1.95	0.49
1:F:126:ILE:HG22	1:F:126:ILE:O	2.12	0.49
1:D:133:ALA:O	1:D:137:LYS:HG3	2.12	0.49
1:D:78:GLU:OE2	1:F:39:ARG:HG3	2.11	0.49
1:B:71:GLU:OE1	1:B:113:ARG:NH1	2.43	0.48
1:B:104:GLU:OE1	4:B:602:HOH:O	2.20	0.48
1:B:27:PHE:CE1	2:B:500:COA:H22	2.49	0.48
1:C:84:ILE:HB	1:C:120:VAL:HB	1.94	0.48
2:A:201[A]:COA:S1P	3:A:202[A]:PG4:H42	2.53	0.48
1:C:145:LEU:HD22	1:E:141:ARG:HD3	1.96	0.48
1:D:145:LEU:HD23	1:F:143:GLU:HG2	1.96	0.48
1:F:89:VAL:HB	2:F:500:COA:O9P	2.12	0.48
2:B:500:COA:H2B	4:B:643:HOH:O	2.14	0.47
1:F:27:PHE:CZ	1:F:88:ALA:HB2	2.50	0.47
1:C:67:LEU:HD12	1:C:87[A]:LEU:HD13	1.96	0.47
1:F:138:LEU:N	1:F:138:LEU:HD12	2.30	0.47
1:F:141:ARG:O	1:F:142:GLU:CG	2.60	0.46
1:F:87:LEU:O	2:F:500:COA:N4P	2.34	0.46
1:D:77:GLN:HG2	1:F:36:ASN:O	2.16	0.46
1:F:23:PHE:HD1	1:F:88:ALA:HB3	1.81	0.46
1:F:41:ALA:O	1:F:45:ARG:HG2	2.16	0.46
1:C:59:GLU:CD	1:C:96:ARG:NH2	2.69	0.46
1:F:59:GLU:HB2	1:F:64:VAL:HG11	1.98	0.45
1:A:123:ASP:HB2	1:A:128:ASP:OD2	2.17	0.45
1:F:26:ALA:HB1	2:F:500:COA:C7P	2.47	0.45
1:D:84:ILE:HB	1:D:120:VAL:HG23	1.99	0.45
1:D:19:LEU:HD11	1:D:66:GLY:HA3	1.99	0.45
1:E:95:ARG:C	1:E:96:ARG:HD2	2.37	0.44
1:B:129:GLU:N	1:B:130:PRO:HD2	2.33	0.44
1:F:138:LEU:N	1:F:138:LEU:CD1	2.81	0.44
1:B:155:ASN:HB3	1:B:156:VAL:H	1.65	0.44
1:F:103:ILE:HG22	1:F:107:LYS:HE3	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:121:GLN:HB3	1:F:144:VAL:CG1	2.47	0.44
1:B:50:ASP:OD1	4:B:603:HOH:O	2.21	0.44
1:E:134:LEU:O	1:E:138:LEU:HD12	2.17	0.44
1:C:1:MET:HE2	1:C:3:LEU:HB2	1.99	0.44
1:D:145:LEU:CD2	1:F:143:GLU:HG2	2.47	0.44
1:A:130:PRO:HB3	2:A:201[B]:COA:N6A	2.33	0.43
1:F:26:ALA:HB1	2:F:500:COA:H72	2.00	0.43
1:A:86:ASP:OD1	3:A:202[A]:PG4:H21	2.18	0.43
1:D:95:ARG:NH2	1:E:71:GLU:OE2	2.42	0.43
1:F:27:PHE:CE1	2:F:500:COA:C6P	3.01	0.43
1:F:95:ARG:O	2:F:500:COA:O9A	2.36	0.43
1:D:74:LYS:HD2	1:D:74:LYS:HA	1.68	0.43
1:E:59:GLU:O	1:E:60:TYR:CB	2.67	0.43
1:C:125:ALA:O	1:C:128:ASP:HB2	2.19	0.43
1:C:141:ARG:HB2	1:E:147:PHE:CE1	2.53	0.42
1:E:63:ILE:HG12	4:E:634:HOH:O	2.19	0.42
1:D:135:TYR:HB3	1:F:147:PHE:CG	2.54	0.42
1:A:115:ALA:O	1:B:151[A]:VAL:HG21	2.19	0.42
1:A:130:PRO:HB3	2:A:201[A]:COA:N6A	2.35	0.42
1:A:146:HIS:HB3	1:B:142:GLU:HG3	2.02	0.41
1:E:94:ARG:NH1	2:E:500:COA:C7P	2.71	0.41
1:D:96:ARG:HB2	1:D:96:ARG:HE	1.78	0.41
1:E:134:LEU:HD12	2:E:500:COA:H52A	2.02	0.41
1:E:131:ALA:HB2	2:E:500:COA:H32	2.00	0.41
1:F:141:ARG:HH11	1:F:141:ARG:HG3	1.85	0.41
1:C:76:GLU:HB2	1:E:33:TYR:HA	2.02	0.41
1:D:134:LEU:HD23	1:D:138:LEU:HD12	2.02	0.41
1:E:129:GLU:N	1:E:130:PRO:HD2	2.36	0.41
1:E:89:VAL:HG23	2:E:500:COA:H132	2.03	0.41
1:E:24:GLY:O	1:E:30:MET:HG2	2.21	0.41
1:E:130:PRO:HB2	2:E:500:COA:H143	2.02	0.41
1:C:115:ALA:O	1:E:151:VAL:HG21	2.20	0.41
1:E:96:ARG:HD2	1:E:96:ARG:N	2.35	0.41
1:B:128:ASP:O	1:B:132:ILE:HG13	2.22	0.40
1:C:37:ARG:HA	1:C:38:PRO:HD3	1.91	0.40
1:E:94:ARG:HH11	2:E:500:COA:H72	1.77	0.40
1:F:129:GLU:N	1:F:130:PRO:CD	2.84	0.40
1:F:11:ASN:ND2	1:F:44:ARG:HH22	2.19	0.40

There are no symmetry-related clashes.

## 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	152/159 (96%)	150 (99%)	2 (1%)	0	100	100
1	B	161/159 (101%)	156 (97%)	5 (3%)	0	100	100
1	C	151/159 (95%)	146 (97%)	5 (3%)	0	100	100
1	D	153/159 (96%)	146 (95%)	7 (5%)	0	100	100
1	E	150/159 (94%)	138 (92%)	10 (7%)	2 (1%)	15	7
1	F	149/159 (94%)	145 (97%)	4 (3%)	0	100	100
All	All	916/954 (96%)	881 (96%)	33 (4%)	2 (0%)	56	47

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	123[A]	ASP
1	E	123[B]	ASP

### 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	123/128 (96%)	122 (99%)	1 (1%)	86	89
1	B	132/128 (103%)	129 (98%)	3 (2%)	58	58
1	C	122/128 (95%)	121 (99%)	1 (1%)	86	89
1	D	123/128 (96%)	120 (98%)	3 (2%)	57	56
1	E	121/128 (94%)	121 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	F	120/128 (94%)	116 (97%)	4 (3%)	45	42
All	All	741/768 (96%)	729 (98%)	12 (2%)	70	72

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	126	ILE
1	B	50	ASP
1	B	51	TYR
1	B	96	ARG
1	C	126	ILE
1	D	96	ARG
1	D	144	VAL
1	D	148	ASP
1	F	30	MET
1	F	93	HIS
1	F	95	ARG
1	F	141	ARG

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

Mol	Chain	Res	Type
1	A	146	HIS

### 5.3.3 RNA ⓘ

There are no RNA molecules 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

9 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	A	201[A]	-	41,50,50	3.96	19 (46%)	49,75,75	1.86	4 (8%)
2	COA	A	201[B]	-	41,50,50	4.01	19 (46%)	49,75,75	1.87	4 (8%)
3	PG4	A	202[A]	-	12,12,12	0.53	0	11,11,11	0.52	0
3	PG4	A	202[B]	-	12,12,12	0.53	0	11,11,11	0.53	0
2	COA	B	500	-	41,50,50	3.94	10 (24%)	49,75,75	2.09	4 (8%)
2	COA	C	500[A]	-	41,50,50	3.84	16 (39%)	49,75,75	2.01	10 (20%)
2	COA	C	500[B]	-	41,50,50	3.85	16 (39%)	49,75,75	2.09	12 (24%)
2	COA	E	500	-	41,50,50	4.00	19 (46%)	49,75,75	1.87	4 (8%)
2	COA	F	500	-	41,50,50	3.91	10 (24%)	49,75,75	2.12	9 (18%)

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	A	201[A]	-	-	0/44/64/64	0/3/3/3
2	COA	A	201[B]	-	-	0/44/64/64	0/3/3/3
3	PG4	A	202[A]	-	-	0/10/10/10	0/0/0/0
3	PG4	A	202[B]	-	-	0/10/10/10	0/0/0/0
2	COA	B	500	-	-	0/44/64/64	0/3/3/3
2	COA	C	500[A]	-	-	0/44/64/64	0/3/3/3
2	COA	C	500[B]	-	-	0/44/64/64	0/3/3/3
2	COA	E	500	-	-	0/44/64/64	0/3/3/3
2	COA	F	500	-	-	0/44/64/64	0/3/3/3

All (109) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	500[B]	COA	C2B-C1B	-14.10	1.31	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	500[A]	COA	C2B-C1B	-13.99	1.31	1.53
2	A	201[B]	COA	C2B-C1B	-13.66	1.31	1.53
2	E	500	COA	C2B-C1B	-13.62	1.32	1.53
2	F	500	COA	C2B-C1B	-13.58	1.32	1.53
2	B	500	COA	C2B-C1B	-13.48	1.32	1.53
2	A	201[A]	COA	C2B-C1B	-13.43	1.32	1.53
2	A	201[B]	COA	O4B-C4B	-8.10	1.26	1.45
2	E	500	COA	O4B-C4B	-8.09	1.26	1.45
2	A	201[A]	COA	O4B-C4B	-7.97	1.26	1.45
2	C	500[B]	COA	O4B-C4B	-7.16	1.28	1.45
2	C	500[A]	COA	O4B-C4B	-7.12	1.28	1.45
2	F	500	COA	O4B-C4B	-6.93	1.29	1.45
2	B	500	COA	O4B-C4B	-6.54	1.30	1.45
2	A	201[B]	COA	O3B-C3B	-4.01	1.31	1.44
2	E	500	COA	O3B-C3B	-3.98	1.31	1.44
2	C	500[A]	COA	O3B-C3B	-3.97	1.31	1.44
2	A	201[A]	COA	O3B-C3B	-3.94	1.31	1.44
2	C	500[B]	COA	O3B-C3B	-3.93	1.31	1.44
2	F	500	COA	O3B-C3B	-3.87	1.31	1.44
2	A	201[B]	COA	OAP-CAP	-3.74	1.34	1.42
2	E	500	COA	OAP-CAP	-3.74	1.34	1.42
2	A	201[A]	COA	OAP-CAP	-3.68	1.34	1.42
2	A	201[B]	COA	P2A-O5A	-3.50	1.40	1.55
2	E	500	COA	P2A-O5A	-3.48	1.40	1.55
2	A	201[A]	COA	P2A-O5A	-3.38	1.40	1.55
2	B	500	COA	O3B-C3B	-3.30	1.33	1.44
2	C	500[B]	COA	O9P-C9P	-3.24	1.17	1.23
2	C	500[B]	COA	P2A-O5A	-3.24	1.41	1.55
2	C	500[A]	COA	P2A-O5A	-3.21	1.41	1.55
2	C	500[A]	COA	O9P-C9P	-3.17	1.17	1.23
2	C	500[B]	COA	P1A-O2A	-3.03	1.42	1.55
2	A	201[A]	COA	O9P-C9P	-2.96	1.17	1.23
2	E	500	COA	O9P-C9P	-2.95	1.17	1.23
2	A	201[B]	COA	O9P-C9P	-2.95	1.17	1.23
2	C	500[B]	COA	OAP-CAP	-2.94	1.36	1.42
2	C	500[A]	COA	P1A-O2A	-2.90	1.42	1.55
2	C	500[B]	COA	P2A-O4A	-2.89	1.40	1.51
2	C	500[A]	COA	OAP-CAP	-2.82	1.36	1.42
2	E	500	COA	C5A-C4A	-2.82	1.34	1.40
2	A	201[B]	COA	C5A-C4A	-2.79	1.34	1.40
2	C	500[A]	COA	P2A-O4A	-2.79	1.41	1.51
2	A	201[B]	COA	C7P-N8P	-2.77	1.39	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	500	COA	C7P-N8P	-2.75	1.39	1.46
2	A	201[A]	COA	C7P-N8P	-2.71	1.39	1.46
2	C	500[A]	COA	C5A-C4A	-2.71	1.34	1.40
2	C	500[B]	COA	C5A-C4A	-2.71	1.34	1.40
2	A	201[A]	COA	C5A-C4A	-2.70	1.34	1.40
2	E	500	COA	P1A-O2A	-2.69	1.43	1.55
2	A	201[B]	COA	P1A-O2A	-2.69	1.43	1.55
2	A	201[A]	COA	P1A-O2A	-2.68	1.43	1.55
2	C	500[B]	COA	P3B-O8A	-2.60	1.45	1.54
2	A	201[A]	COA	P2A-O4A	-2.59	1.41	1.51
2	A	201[B]	COA	P2A-O4A	-2.59	1.41	1.51
2	E	500	COA	P2A-O4A	-2.58	1.41	1.51
2	C	500[A]	COA	P3B-O8A	-2.57	1.45	1.54
2	C	500[B]	COA	C7P-N8P	-2.52	1.40	1.46
2	F	500	COA	C5A-C4A	-2.48	1.34	1.40
2	E	500	COA	P3B-O8A	-2.40	1.46	1.54
2	A	201[B]	COA	P3B-O8A	-2.39	1.46	1.54
2	A	201[A]	COA	P3B-O9A	-2.39	1.46	1.54
2	A	201[A]	COA	P3B-O7A	-2.37	1.43	1.50
2	C	500[A]	COA	C7P-N8P	-2.36	1.40	1.46
2	A	201[A]	COA	P3B-O8A	-2.35	1.46	1.54
2	C	500[B]	COA	P1A-O1A	-2.28	1.42	1.51
2	A	201[B]	COA	P3B-O9A	-2.26	1.47	1.54
2	E	500	COA	P1A-O1A	-2.25	1.43	1.51
2	A	201[B]	COA	P1A-O1A	-2.25	1.43	1.51
2	E	500	COA	P3B-O9A	-2.24	1.47	1.54
2	A	201[B]	COA	P3B-O7A	-2.24	1.43	1.50
2	E	500	COA	P3B-O7A	-2.22	1.43	1.50
2	A	201[A]	COA	P1A-O1A	-2.20	1.43	1.51
2	E	500	COA	C5A-N7A	-2.17	1.31	1.39
2	C	500[A]	COA	P1A-O1A	-2.16	1.43	1.51
2	A	201[B]	COA	C5A-N7A	-2.16	1.31	1.39
2	E	500	COA	O6A-CCP	-2.09	1.36	1.43
2	A	201[A]	COA	C5A-N7A	-2.08	1.32	1.39
2	C	500[B]	COA	O6A-CCP	-2.07	1.36	1.43
2	A	201[A]	COA	O6A-CCP	-2.07	1.36	1.43
2	A	201[B]	COA	O6A-CCP	-2.07	1.36	1.43
2	B	500	COA	C5A-C4A	-2.07	1.35	1.40
2	C	500[A]	COA	O6A-CCP	-2.00	1.36	1.43
2	F	500	COA	C6A-N6A	2.23	1.43	1.34
2	F	500	COA	C2A-N3A	2.31	1.36	1.32
2	F	500	COA	O2B-C2B	2.31	1.48	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	500	COA	O2B-C2B	2.60	1.49	1.43
2	B	500	COA	C6A-N6A	2.60	1.44	1.34
2	B	500	COA	C2A-N3A	2.84	1.37	1.32
2	E	500	COA	C9P-N8P	3.00	1.39	1.33
2	A	201[B]	COA	C9P-N8P	3.00	1.39	1.33
2	A	201[A]	COA	C9P-N8P	3.03	1.39	1.33
2	F	500	COA	C9P-N8P	3.62	1.41	1.33
2	C	500[A]	COA	C9P-N8P	3.76	1.41	1.33
2	C	500[B]	COA	C9P-N8P	3.78	1.41	1.33
2	C	500[B]	COA	C5P-N4P	3.88	1.42	1.33
2	A	201[B]	COA	C5P-N4P	4.06	1.43	1.33
2	A	201[A]	COA	C5P-N4P	4.07	1.43	1.33
2	E	500	COA	C5P-N4P	4.09	1.43	1.33
2	C	500[A]	COA	C5P-N4P	4.28	1.43	1.33
2	B	500	COA	C5P-N4P	5.06	1.45	1.33
2	B	500	COA	C9P-N8P	5.26	1.44	1.33
2	F	500	COA	C5P-N4P	5.87	1.47	1.33
2	C	500[B]	COA	O4B-C1B	14.08	1.61	1.41
2	C	500[A]	COA	O4B-C1B	14.12	1.61	1.41
2	A	201[A]	COA	O4B-C1B	15.43	1.63	1.41
2	E	500	COA	O4B-C1B	15.48	1.63	1.41
2	A	201[B]	COA	O4B-C1B	15.51	1.63	1.41
2	F	500	COA	O4B-C1B	16.63	1.65	1.41
2	B	500	COA	O4B-C1B	17.22	1.65	1.41

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	500	COA	N3A-C2A-N1A	-11.05	120.19	128.87
2	F	500	COA	N3A-C2A-N1A	-9.65	121.29	128.87
2	A	201[A]	COA	N6A-C6A-N1A	-7.84	105.36	118.52
2	E	500	COA	N6A-C6A-N1A	-7.83	105.38	118.52
2	A	201[B]	COA	N6A-C6A-N1A	-7.81	105.41	118.52
2	A	201[A]	COA	N3A-C2A-N1A	-7.65	122.86	128.87
2	E	500	COA	N3A-C2A-N1A	-7.45	123.02	128.87
2	A	201[B]	COA	N3A-C2A-N1A	-7.43	123.04	128.87
2	C	500[B]	COA	N3A-C2A-N1A	-7.40	123.06	128.87
2	C	500[A]	COA	N3A-C2A-N1A	-7.32	123.12	128.87
2	C	500[B]	COA	N6A-C6A-N1A	-7.23	106.38	118.52
2	C	500[A]	COA	N6A-C6A-N1A	-7.14	106.53	118.52
2	F	500	COA	N6A-C6A-N1A	-6.94	106.87	118.52
2	B	500	COA	N6A-C6A-N1A	-6.74	107.20	118.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	500[B]	COA	C4B-O4B-C1B	-4.36	105.03	109.64
2	C	500[A]	COA	C4B-O4B-C1B	-4.20	105.19	109.64
2	C	500[B]	COA	C2P-C3P-N4P	-3.10	106.30	112.44
2	F	500	COA	C2B-C1B-N9A	-3.07	105.24	113.47
2	F	500	COA	OAP-CAP-C9P	-2.95	103.51	110.52
2	F	500	COA	O5P-C5P-C6P	-2.77	117.15	121.97
2	C	500[B]	COA	O9P-C9P-N8P	-2.77	117.54	123.04
2	C	500[A]	COA	O9P-C9P-N8P	-2.72	117.64	123.04
2	F	500	COA	C2B-C3B-C4B	-2.29	98.87	103.25
2	F	500	COA	O4B-C4B-C3B	-2.20	99.81	104.89
2	F	500	COA	C7P-N8P-C9P	-2.09	118.42	122.62
2	C	500[A]	COA	C1B-N9A-C4A	-2.08	124.49	126.81
2	C	500[B]	COA	O2B-C2B-C3B	-2.07	105.16	111.13
2	B	500	COA	C6P-C5P-N4P	2.05	120.02	116.46
2	C	500[B]	COA	CAP-C9P-N8P	2.07	121.19	116.31
2	F	500	COA	O8A-P3B-O3B	2.08	112.84	106.62
2	C	500[A]	COA	C6P-C7P-N8P	2.09	116.62	111.94
2	C	500[B]	COA	C3B-C2B-C1B	2.09	104.61	100.06
2	C	500[A]	COA	C3B-C2B-C1B	2.15	104.75	100.06
2	A	201[A]	COA	O2A-P1A-O3A	2.32	115.23	105.27
2	A	201[B]	COA	O2A-P1A-O3A	2.41	115.60	105.27
2	B	500	COA	C3B-C2B-C1B	2.41	105.31	100.06
2	E	500	COA	O2A-P1A-O3A	2.42	115.62	105.27
2	C	500[B]	COA	C6P-C7P-N8P	2.44	117.41	111.94
2	A	201[A]	COA	O9A-P3B-O3B	2.53	114.18	106.62
2	E	500	COA	O9A-P3B-O3B	2.53	114.20	106.62
2	A	201[B]	COA	O9A-P3B-O3B	2.54	114.20	106.62
2	C	500[B]	COA	O9A-P3B-O3B	2.58	114.34	106.62
2	C	500[A]	COA	O2A-P1A-O3A	2.62	116.48	105.27
2	C	500[A]	COA	O9A-P3B-O3B	2.63	114.48	106.62
2	C	500[B]	COA	C7P-C6P-C5P	2.77	116.97	112.22
2	C	500[B]	COA	O2A-P1A-O3A	2.83	117.38	105.27
2	C	500[A]	COA	C7P-C6P-C5P	3.16	117.63	112.22

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

7 monomers are involved in 45 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	201[A]	COA	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	201[B]	COA	1	0
3	A	202[A]	PG4	7	0
2	B	500	COA	8	0
2	C	500[A]	COA	1	0
2	E	500	COA	11	0
2	F	500	COA	16	0

## 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 ⓘ

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	151/159 (94%)	0.05	1 (0%)	89 89	21, 34, 65, 91	0
1	B	159/159 (100%)	0.26	9 (5%)	27 29	19, 37, 78, 134	0
1	C	152/159 (95%)	0.50	14 (9%)	11 12	23, 48, 82, 121	0
1	D	154/159 (96%)	0.67	19 (12%)	5 6	32, 56, 103, 149	0
1	E	151/159 (94%)	1.01	31 (20%)	1 1	36, 68, 101, 113	0
1	F	151/159 (94%)	1.33	41 (27%)	1 1	38, 72, 132, 162	0
All	All	918/954 (96%)	0.63	115 (12%)	5 5	19, 53, 107, 162	0

All (115) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	96	ARG	8.8
1	F	126	ILE	8.2
1	D	126	ILE	7.7
1	E	126	ILE	6.4
1	B	126	ILE	6.1
1	F	89	VAL	6.1
1	F	25	GLU	6.0
1	F	127	GLU	6.0
1	C	125	ALA	5.9
1	E	2	SER	5.8
1	F	95	ARG	5.7
1	E	30	MET	5.6
1	F	27	PHE	5.2
1	F	97	GLY	5.0
1	F	28	ASN	4.9
1	E	31	GLU	4.9
1	C	127	GLU	4.8
1	F	124	THR	4.8
1	E	125	ALA	4.4

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Mol	Chain	Res	Type	RSRZ
1	D	127	GLU	4.3
1	E	97	GLY	4.3
1	F	2	SER	4.2
1	B	156	VAL	4.2
1	F	125	ALA	4.2
1	C	126	ILE	4.0
1	D	145	LEU	4.0
1	F	142	GLU	4.0
1	D	124	THR	3.9
1	E	60	TYR	3.9
1	E	59	GLU	3.8
1	F	137	LYS	3.7
1	E	127	GLU	3.7
1	E	140	VAL	3.7
1	F	119	PHE	3.5
1	C	152	SER	3.5
1	B	125	ALA	3.5
1	F	61	GLY	3.4
1	D	67	LEU	3.4
1	E	96	ARG	3.4
1	E	3	LEU	3.4
1	F	94	ARG	3.3
1	F	140	VAL	3.3
1	F	39	ARG	3.3
1	E	61	GLY	3.2
1	B	124	THR	3.2
1	F	59	GLU	3.2
1	B	154	ASN	3.2
1	D	152	SER	3.1
1	B	127	GLU	3.1
1	C	67	LEU	3.1
1	F	135	TYR	3.1
1	E	4	ASP	3.1
1	E	28	ASN	3.0
1	C	87[A]	LEU	3.0
1	E	39	ARG	3.0
1	D	91	LYS	3.0
1	F	26	ALA	3.0
1	E	137	LYS	2.9
1	E	152	SER	2.9
1	E	44	ARG	2.9
1	E	108	GLU	2.9

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Mol	Chain	Res	Type	RSRZ
1	E	45	ARG	2.9
1	E	135	TYR	2.9
1	E	88	ALA	2.8
1	F	152	SER	2.8
1	F	3	LEU	2.8
1	E	89	VAL	2.8
1	E	95	ARG	2.8
1	E	25	GLU	2.8
1	F	29	ASP	2.8
1	C	123	ASP	2.7
1	F	4	ASP	2.7
1	F	60	TYR	2.6
1	C	31	GLU	2.6
1	F	31	GLU	2.6
1	A	127	GLU	2.6
1	D	28	ASN	2.6
1	F	129	GLU	2.6
1	D	55	LEU	2.6
1	B	155	ASN	2.6
1	D	140	VAL	2.5
1	F	87	LEU	2.4
1	B	152	SER	2.4
1	F	67	LEU	2.4
1	E	41	ALA	2.4
1	D	32	THR	2.3
1	B	157	ASP	2.3
1	D	123	ASP	2.3
1	D	68	ALA	2.3
1	C	25	GLU	2.3
1	D	129	GLU	2.3
1	D	56	ALA	2.2
1	F	138	LEU	2.2
1	F	62	GLU	2.2
1	F	45	ARG	2.2
1	F	139	GLY	2.2
1	F	148	ASP	2.2
1	D	25	GLU	2.2
1	E	67	LEU	2.1
1	E	11	ASN	2.1
1	F	41	ALA	2.1
1	F	68	ALA	2.1
1	E	29	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	75	PHE	2.1
1	C	34	THR	2.1
1	F	130	PRO	2.1
1	C	68	ALA	2.1
1	D	41	ALA	2.0
1	D	151	VAL	2.0
1	D	44	ARG	2.0
1	C	134	LEU	2.0
1	F	92	ALA	2.0
1	E	24	GLY	2.0
1	C	128	ASP	2.0
1	F	128	ASP	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, 95<sup>th</sup> 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	PG4	A	202[A]	13/13	0.51	0.34	6.94	66,79,84,88	13
3	PG4	A	202[B]	13/13	0.51	0.34	6.63	63,81,86,87	13
2	COA	B	500	48/48	0.74	0.42	3.98	75,177,238,250	0
2	COA	E	500	48/48	0.66	0.33	1.18	72,108,144,145	0
2	COA	F	500	48/48	0.69	0.30	0.07	89,124,153,215	0
2	COA	A	201[B]	48/48	0.94	0.12	-0.48	27,42,63,79	48
2	COA	A	201[A]	48/48	0.94	0.12	-0.51	27,42,63,79	48
2	COA	C	500[B]	48/48	0.93	0.13	-0.52	30,53,71,75	48
2	COA	C	500[A]	48/48	0.93	0.13	-0.56	30,53,71,75	48

## 6.5 Other polymers

There are no such residues in this entry.