



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 04:59 am GMT

PDB ID : 4UBU  
Title : Structure of a modified C93S variant of the 3-ketoacyl-CoA thiolase FadA5 from *M. tuberculosis* in complex with CoA  
Authors : Schaefer, C.M.; Kisker, C.  
Deposited on : 2014-08-13  
Resolution : 3.00 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 : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : recalc28949

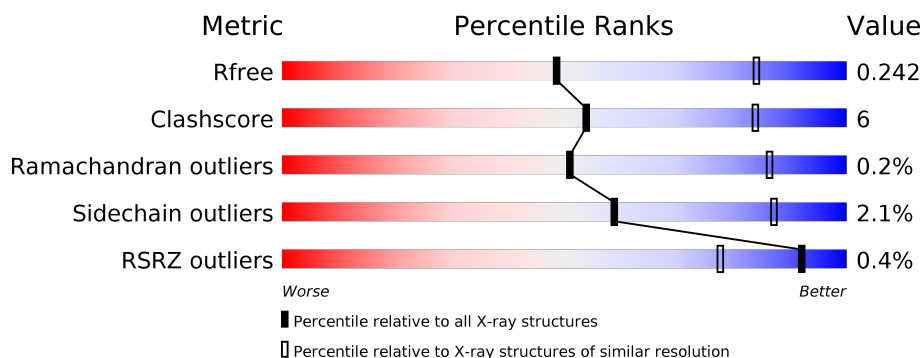
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.00 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))
$R_{free}$	100719	1692 (3.00-3.00)
Clashscore	112137	2037 (3.00-3.00)
Ramachandran outliers	110173	1973 (3.00-3.00)
Sidechain outliers	110143	1976 (3.00-3.00)
RSRZ outliers	101464	1716 (3.00-3.00)

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	399	<div> <div>84%</div> <div>12%</div> <div>••</div> </div>
1	B	399	<div> <div>84%</div> <div>13%</div> <div>••</div> </div>
1	C	399	<div> <div>84%</div> <div>13%</div> <div>••</div> </div>
1	D	399	<div> <div>88%</div> <div>10%</div> <div>••</div> </div>
1	E	399	<div> <div>85%</div> <div>12%</div> <div>••</div> </div>
1	F	399	<div> <div>86%</div> <div>11%</div> <div>••</div> </div>

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Mol	Chain	Length	Quality of chain
1	G	399	<div> <div>%</div> <div> </div> <div>86% 11% ..</div> </div>
1	H	399	<div> <div>%</div> <div> </div> <div>86% 12% ..</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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	COA	A	400	-	-	-	X
2	COA	B	400	-	-	-	X
2	COA	C	400	-	-	-	X
2	COA	D	400	-	-	-	X
2	COA	E	401	-	-	-	X
2	COA	F	400	-	-	-	X
2	COA	G	401	-	-	-	X
2	COA	H	401	-	-	-	X
3	GOL	E	402	-	-	-	X
3	GOL	G	403	-	-	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 23784 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 Acetyl-CoA acetyltransferase FadA5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	391	Total	C	N	O	S	0	1	0
			2911	1801	539	559	12			
1	B	391	Total	C	N	O	S	0	2	0
			2918	1806	541	559	12			
1	C	391	Total	C	N	O	S	0	0	0
			2903	1796	536	559	12			
1	D	391	Total	C	N	O	S	0	1	0
			2911	1801	539	559	12			
1	E	392	Total	C	N	O	S	0	1	0
			2917	1804	540	561	12			
1	F	391	Total	C	N	O	S	0	2	0
			2917	1805	539	561	12			
1	G	393	Total	C	N	O	S	0	1	0
			2921	1806	541	562	12			
1	H	394	Total	C	N	O	S	0	2	0
			2939	1820	543	564	12			

There are 64 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	HIS	-	expression tag	UNP I6XHI4
A	-6	HIS	-	expression tag	UNP I6XHI4
A	-5	HIS	-	expression tag	UNP I6XHI4
A	-4	HIS	-	expression tag	UNP I6XHI4
A	-3	HIS	-	expression tag	UNP I6XHI4
A	-2	HIS	-	expression tag	UNP I6XHI4
A	-1	GLY	-	expression tag	UNP I6XHI4
A	0	SER	-	expression tag	UNP I6XHI4
B	-7	HIS	-	expression tag	UNP I6XHI4
B	-6	HIS	-	expression tag	UNP I6XHI4
B	-5	HIS	-	expression tag	UNP I6XHI4
B	-4	HIS	-	expression tag	UNP I6XHI4
B	-3	HIS	-	expression tag	UNP I6XHI4

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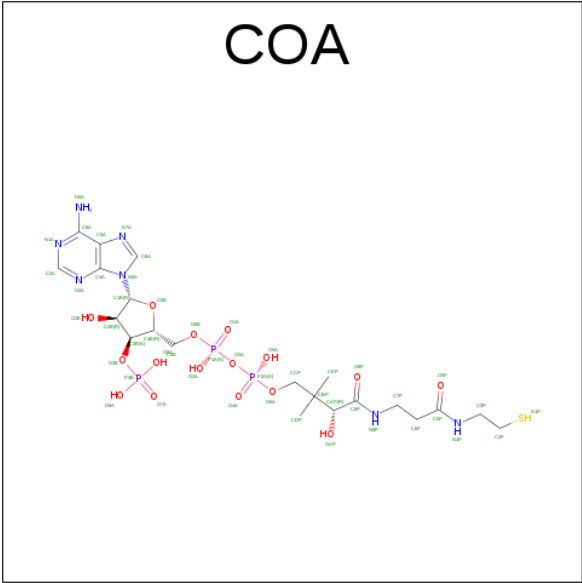
Chain	Residue	Modelled	Actual	Comment	Reference
B	-2	HIS	-	expression tag	UNP I6XHI4
B	-1	GLY	-	expression tag	UNP I6XHI4
B	0	SER	-	expression tag	UNP I6XHI4
C	-7	HIS	-	expression tag	UNP I6XHI4
C	-6	HIS	-	expression tag	UNP I6XHI4
C	-5	HIS	-	expression tag	UNP I6XHI4
C	-4	HIS	-	expression tag	UNP I6XHI4
C	-3	HIS	-	expression tag	UNP I6XHI4
C	-2	HIS	-	expression tag	UNP I6XHI4
C	-1	GLY	-	expression tag	UNP I6XHI4
C	0	SER	-	expression tag	UNP I6XHI4
D	-7	HIS	-	expression tag	UNP I6XHI4
D	-6	HIS	-	expression tag	UNP I6XHI4
D	-5	HIS	-	expression tag	UNP I6XHI4
D	-4	HIS	-	expression tag	UNP I6XHI4
D	-3	HIS	-	expression tag	UNP I6XHI4
D	-2	HIS	-	expression tag	UNP I6XHI4
D	-1	GLY	-	expression tag	UNP I6XHI4
D	0	SER	-	expression tag	UNP I6XHI4
E	-7	HIS	-	expression tag	UNP I6XHI4
E	-6	HIS	-	expression tag	UNP I6XHI4
E	-5	HIS	-	expression tag	UNP I6XHI4
E	-4	HIS	-	expression tag	UNP I6XHI4
E	-3	HIS	-	expression tag	UNP I6XHI4
E	-2	HIS	-	expression tag	UNP I6XHI4
E	-1	GLY	-	expression tag	UNP I6XHI4
E	0	SER	-	expression tag	UNP I6XHI4
F	-7	HIS	-	expression tag	UNP I6XHI4
F	-6	HIS	-	expression tag	UNP I6XHI4
F	-5	HIS	-	expression tag	UNP I6XHI4
F	-4	HIS	-	expression tag	UNP I6XHI4
F	-3	HIS	-	expression tag	UNP I6XHI4
F	-2	HIS	-	expression tag	UNP I6XHI4
F	-1	GLY	-	expression tag	UNP I6XHI4
F	0	SER	-	expression tag	UNP I6XHI4
G	-7	HIS	-	expression tag	UNP I6XHI4
G	-6	HIS	-	expression tag	UNP I6XHI4
G	-5	HIS	-	expression tag	UNP I6XHI4
G	-4	HIS	-	expression tag	UNP I6XHI4
G	-3	HIS	-	expression tag	UNP I6XHI4
G	-2	HIS	-	expression tag	UNP I6XHI4
G	-1	GLY	-	expression tag	UNP I6XHI4

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Chain	Residue	Modelled	Actual	Comment	Reference
G	0	SER	-	expression tag	UNP I6XHI4
H	-7	HIS	-	expression tag	UNP I6XHI4
H	-6	HIS	-	expression tag	UNP I6XHI4
H	-5	HIS	-	expression tag	UNP I6XHI4
H	-4	HIS	-	expression tag	UNP I6XHI4
H	-3	HIS	-	expression tag	UNP I6XHI4
H	-2	HIS	-	expression tag	UNP I6XHI4
H	-1	GLY	-	expression tag	UNP I6XHI4
H	0	SER	-	expression tag	UNP I6XHI4

- 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 48	C 21	N 7	O 16	P 3	S 1	0	0
2	B	1	Total 48	C 21	N 7	O 16	P 3	S 1	0	0
2	C	1	Total 48	C 21	N 7	O 16	P 3	S 1	0	0
2	D	1	Total 48	C 21	N 7	O 16	P 3	S 1	0	0
2	E	1	Total 48	C 21	N 7	O 16	P 3	S 1	0	0
2	F	1	Total 48	C 21	N 7	O 16	P 3	S 1	0	0
2	G	1	Total 48	C 21	N 7	O 16	P 3	S 1	0	0

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

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	E	1	Total	C	O	0	0
			6	3	3		
3	G	1	Total	C	O	0	0
			6	3	3		
3	G	1	Total	C	O	0	0
			6	3	3		
3	H	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	9	Total	O	0	0
			9	9		
4	B	6	Total	O	0	0
			6	6		
4	C	5	Total	O	0	0
			5	5		
4	E	3	Total	O	0	0
			3	3		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	F	13	Total 13	O 13	0	0
4	G	1	Total 1	O 1	0	0
4	H	2	Total 2	O 2	0	0

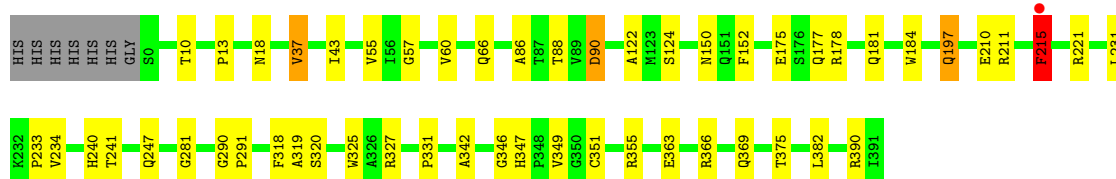






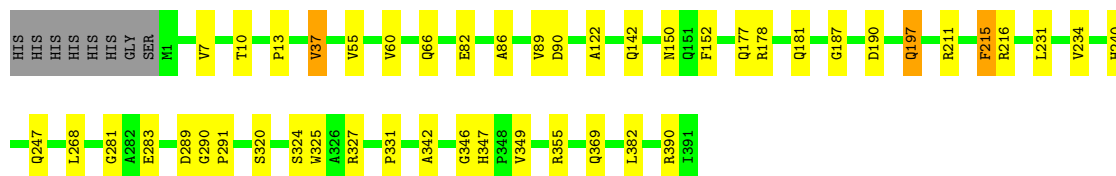
• Molecule 1: Acetyl-CoA acetyltransferase FadA5

Chain E: 85% 12% ..



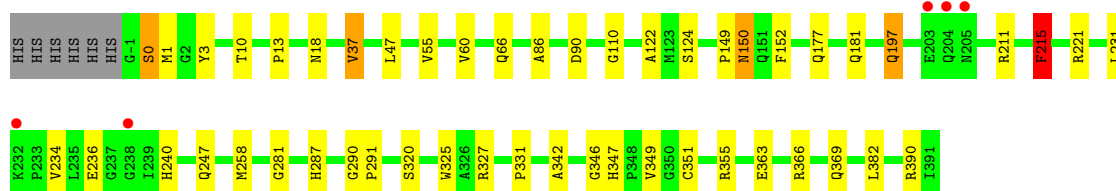
• Molecule 1: Acetyl-CoA acetyltransferase FadA5

Chain F: 86% 11% ..



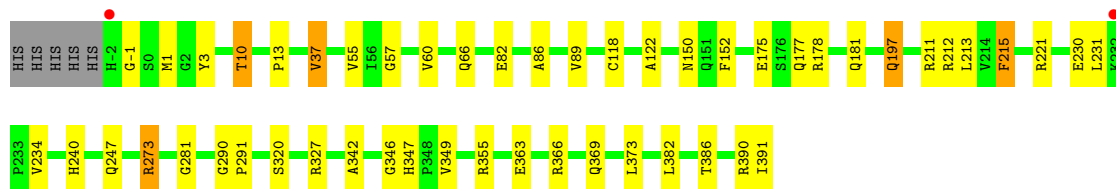
• Molecule 1: Acetyl-CoA acetyltransferase FadA5

Chain G: 86% 11% ..



• Molecule 1: Acetyl-CoA acetyltransferase FadA5

Chain H: 86% 12% ..



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	127.61Å 105.16Å 147.66Å 90.00° 107.33° 90.00°	Depositor
Resolution (Å)	58.82 – 3.00 58.75 – 3.00	Depositor EDS
% Data completeness (in resolution range)	94.7 (58.82-3.00) 94.7 (58.75-3.00)	Depositor EDS
$R_{merge}$	0.17	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.60 (at 3.01Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, $R_{free}$	0.247 , 0.287 0.212 , 0.242	Depositor DCC
$R_{free}$ test set	3551 reflections (5.27%)	DCC
Wilson B-factor (Å <sup>2</sup> )	41.4	Xtriage
Anisotropy	0.546	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 68.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.44$ , $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	23784	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 37.91 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 3.9909e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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, 0.333 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, GOL, OAS

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	0.65	0/2948	0.74	2/3996 (0.1%)
1	B	0.64	0/2959	0.73	2/4011 (0.0%)
1	C	0.66	0/2937	0.74	2/3982 (0.1%)
1	D	0.64	0/2948	0.74	1/3996 (0.0%)
1	E	0.65	0/2954	0.74	3/4004 (0.1%)
1	F	0.63	0/2957	0.74	2/4008 (0.0%)
1	G	0.64	0/2958	0.73	2/4009 (0.0%)
1	H	0.58	0/2982	0.74	2/4044 (0.0%)
All	All	0.64	0/23643	0.74	16/32050 (0.0%)

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	A	0	1
1	B	0	1
1	E	0	1
All	All	0	3

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	355	ARG	NE-CZ-NH2	-7.63	116.49	120.30
1	C	355	ARG	NE-CZ-NH2	-6.80	116.90	120.30
1	C	355	ARG	NE-CZ-NH1	6.61	123.61	120.30
1	H	355	ARG	NE-CZ-NH2	-6.46	117.07	120.30
1	G	355	ARG	NE-CZ-NH2	-6.45	117.07	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	355	ARG	NE-CZ-NH2	-6.19	117.20	120.30
1	E	355	ARG	NE-CZ-NH2	-5.97	117.31	120.30
1	F	355	ARG	NE-CZ-NH1	5.87	123.23	120.30
1	H	355	ARG	NE-CZ-NH1	5.74	123.17	120.30
1	D	355	ARG	NE-CZ-NH2	-5.73	117.44	120.30
1	A	178	ARG	NE-CZ-NH1	5.59	123.09	120.30
1	B	355	ARG	NE-CZ-NH1	5.59	123.09	120.30
1	G	215	PHE	CB-CG-CD2	-5.21	117.15	120.80
1	E	215	PHE	CB-CG-CD2	-5.09	117.24	120.80
1	E	215	PHE	CB-CG-CD1	5.07	124.35	120.80
1	B	215	PHE	CB-CG-CD2	-5.01	117.30	120.80

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	90	ASP	Peptide
1	B	90	ASP	Peptide
1	E	90	ASP	Peptide

## 5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2911	0	2910	57	0
1	B	2918	0	2917	41	0
1	C	2903	0	2897	32	0
1	D	2911	0	2910	26	0
1	E	2917	0	2915	53	0
1	F	2917	0	2916	43	0
1	G	2921	0	2918	41	0
1	H	2939	0	2928	31	0
2	A	48	0	32	3	0
2	B	48	0	32	2	0
2	C	48	0	32	3	0
2	D	48	0	32	2	0
2	E	48	0	32	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	F	48	0	32	1	0
2	G	48	0	32	3	0
2	H	48	0	32	3	0
3	E	6	0	8	1	0
3	G	12	0	16	2	0
3	H	6	0	8	1	0
4	A	9	0	0	0	0
4	B	6	0	0	0	0
4	C	5	0	0	1	0
4	E	3	0	0	0	0
4	F	13	0	0	3	0
4	G	1	0	0	0	0
4	H	2	0	0	0	0
All	All	23784	0	23599	289	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 6.

All (289) 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:B:190:ASP:HB3	1:F:215:PHE:CE2	1.86	1.11
1:A:195:PRO:HG2	1:E:215:PHE:HZ	1.05	1.10
1:B:190:ASP:HB3	1:F:215:PHE:CZ	1.87	1.09
1:E:178[B]:ARG:HH11	1:E:178[B]:ARG:HG2	0.95	1.08
1:A:195:PRO:CG	1:E:215:PHE:HZ	1.69	1.06
1:B:215:PHE:CE2	1:F:190:ASP:HB3	1.94	1.03
1:A:195:PRO:HG2	1:E:215:PHE:CZ	1.94	1.01
1:B:215:PHE:CZ	1:F:190:ASP:HB3	1.96	1.00
1:F:178[B]:ARG:CG	1:F:178[B]:ARG:HH11	1.79	0.95
1:E:178[B]:ARG:HG2	1:E:178[B]:ARG:NH1	1.72	0.88
1:F:178[B]:ARG:HH11	1:F:178[B]:ARG:HG2	1.43	0.83
1:E:178[B]:ARG:HH11	1:E:178[B]:ARG:CG	1.86	0.82
1:A:195:PRO:CG	1:E:215:PHE:CZ	2.59	0.82
1:B:190:ASP:CB	1:F:215:PHE:CZ	2.66	0.78
1:F:142:GLN:HG3	4:F:506:HOH:O	1.82	0.77
1:B:187:GLY:HA2	1:F:216:ARG:NH2	1.99	0.77
1:B:216:ARG:NH2	1:F:187:GLY:HA2	2.01	0.76
1:A:215:PHE:CE2	1:E:215:PHE:CD2	2.75	0.74
1:B:215:PHE:HD1	1:B:215:PHE:N	1.87	0.71
1:B:215:PHE:N	1:B:215:PHE:CD1	2.59	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:215:PHE:HD1	1:F:215:PHE:N	1.90	0.70
1:G:215:PHE:HD1	1:G:215:PHE:N	1.89	0.69
1:A:215:PHE:HD1	1:A:215:PHE:N	1.90	0.69
1:D:215:PHE:CD1	1:D:215:PHE:N	2.60	0.69
1:D:215:PHE:HD1	1:D:215:PHE:N	1.89	0.69
1:B:215:PHE:CZ	1:F:190:ASP:CB	2.75	0.68
1:A:215:PHE:CD1	1:A:215:PHE:N	2.61	0.68
1:H:215:PHE:HD1	1:H:215:PHE:N	1.91	0.68
1:F:215:PHE:CD1	1:F:215:PHE:N	2.61	0.68
1:A:265:ALA:N	1:G:1:MET:CE	2.57	0.67
1:C:240:HIS:HE1	1:C:320:SER:H	1.43	0.67
1:A:216:ARG:NH2	1:E:184:TRP:O	2.29	0.66
1:E:215:PHE:N	1:E:215:PHE:HD1	1.94	0.66
1:B:240:HIS:HE1	1:B:320:SER:H	1.44	0.66
1:H:215:PHE:CD1	1:H:215:PHE:N	2.63	0.66
1:C:215:PHE:CD1	1:C:215:PHE:N	2.64	0.66
1:C:215:PHE:HD1	1:C:215:PHE:N	1.93	0.65
1:A:240:HIS:HE1	1:A:320:SER:H	1.43	0.65
1:G:215:PHE:CD1	1:G:215:PHE:N	2.60	0.65
1:E:215:PHE:N	1:E:215:PHE:CD1	2.64	0.65
1:H:240:HIS:HE1	1:H:320:SER:H	1.45	0.64
1:A:215:PHE:CD2	1:E:215:PHE:CE2	2.86	0.64
1:E:240:HIS:HE1	1:E:320:SER:H	1.44	0.63
1:A:265:ALA:N	1:G:1:MET:HE1	2.14	0.63
1:H:273:ARG:HD3	3:H:402:GOL:H2	1.80	0.63
1:F:178[B]:ARG:NH1	1:F:178[B]:ARG:HG2	2.11	0.62
1:F:215:PHE:HD1	1:F:215:PHE:H	1.47	0.62
1:F:240:HIS:HE1	1:F:320:SER:H	1.44	0.62
1:D:240:HIS:HE1	1:D:320:SER:H	1.47	0.62
1:G:240:HIS:HE1	1:G:320:SER:H	1.46	0.62
1:A:215:PHE:HD1	1:A:215:PHE:H	1.46	0.61
1:F:283:GLU:HG3	4:F:511:HOH:O	1.98	0.61
1:B:215:PHE:HD1	1:B:215:PHE:H	1.47	0.60
1:G:215:PHE:H	1:G:215:PHE:HD1	1.47	0.60
1:E:215:PHE:H	1:E:215:PHE:HD1	1.48	0.60
1:D:215:PHE:H	1:D:215:PHE:HD1	1.48	0.60
1:F:281:GLY:HA2	1:F:382:LEU:HD23	1.82	0.60
1:A:213:LEU:HD23	1:E:215:PHE:CD2	2.36	0.60
1:H:215:PHE:HD1	1:H:215:PHE:H	1.48	0.59
1:C:215:PHE:HD1	1:C:215:PHE:H	1.49	0.59
1:C:281:GLY:HA2	1:C:382:LEU:HD23	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:175:GLU:OE1	1:H:178:ARG:HD3	2.03	0.59
1:F:90:ASP:OD1	1:F:90:ASP:C	2.40	0.58
1:D:90:ASP:C	1:D:90:ASP:OD1	2.41	0.58
1:D:281:GLY:HA2	1:D:382:LEU:HD23	1.86	0.58
1:D:55:VAL:O	1:D:86:ALA:HA	2.04	0.58
1:B:281:GLY:HA2	1:B:382:LEU:HD23	1.85	0.58
1:B:89:VAL:HG12	1:C:89:VAL:HG12	1.86	0.57
1:A:43:ILE:HG13	1:E:210:GLU:OE2	2.04	0.57
1:A:265:ALA:N	1:G:1:MET:HE3	2.20	0.57
1:H:363:GLU:OE2	1:H:366:ARG:NH2	2.38	0.57
1:A:150:ASN:OD1	1:A:150:ASN:C	2.43	0.56
1:F:178[B]:ARG:HH11	1:F:178[B]:ARG:HG3	1.66	0.56
1:G:18:ASN:N	1:G:124:SER:OG	2.36	0.56
1:F:55:VAL:O	1:F:86:ALA:HA	2.06	0.56
1:G:55:VAL:O	1:G:86:ALA:HA	2.05	0.55
1:A:281:GLY:HA2	1:A:382:LEU:HD23	1.88	0.55
1:G:342:ALA:HB1	1:G:347:HIS:HB2	1.88	0.55
1:G:281:GLY:HA2	1:G:382:LEU:HD23	1.87	0.55
1:A:89:VAL:HG12	1:D:89:VAL:HG12	1.88	0.54
1:F:82:GLU:HG3	4:F:503:HOH:O	2.07	0.54
1:C:342:ALA:HB1	1:C:347:HIS:HB2	1.88	0.54
1:G:90:ASP:OD1	1:G:90:ASP:C	2.45	0.54
1:H:150:ASN:OD1	1:H:150:ASN:C	2.46	0.54
1:A:342:ALA:HB1	1:A:347:HIS:HB2	1.90	0.54
1:E:281:GLY:HA2	1:E:382:LEU:HD23	1.89	0.54
1:D:342:ALA:HB1	1:D:347:HIS:HB2	1.90	0.54
1:A:197:GLN:HG2	1:A:211:ARG:HG2	1.90	0.53
1:C:90:ASP:OD1	1:C:90:ASP:C	2.44	0.53
1:A:265:ALA:HB2	1:G:1:MET:HE1	1.89	0.53
1:A:3:TYR:HD2	1:G:3:TYR:CE2	2.27	0.53
1:E:342:ALA:HB1	1:E:347:HIS:HB2	1.90	0.53
1:F:342:ALA:HB1	1:F:347:HIS:HB2	1.90	0.53
1:F:89:VAL:HG12	1:H:89:VAL:HG12	1.90	0.53
1:A:264:ARG:C	1:G:1:MET:HE3	2.29	0.52
1:D:231:LEU:HD21	2:D:400:COA:H1B	1.91	0.52
1:H:281:GLY:HA2	1:H:382:LEU:HD23	1.91	0.52
1:F:152:PHE:CD1	1:F:234:VAL:HG11	2.44	0.52
1:C:60:VAL:HG21	1:C:349:VAL:CG1	2.39	0.52
1:A:3:TYR:HD2	1:G:3:TYR:CZ	2.27	0.52
1:D:60:VAL:HG21	1:D:349:VAL:CG1	2.39	0.52
1:F:281:GLY:HA3	1:H:82:GLU:O	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:231:LEU:HD21	2:H:401:COA:H1B	1.91	0.52
1:B:231:LEU:HD21	2:B:400:COA:H1B	1.90	0.52
1:A:67:SER:OG	1:A:68:ASN:ND2	2.42	0.52
1:H:55:VAL:O	1:H:86:ALA:HA	2.10	0.52
1:F:60:VAL:HG21	1:F:349:VAL:CG1	2.40	0.51
1:D:13:PRO:HG3	1:D:215:PHE:HA	1.93	0.51
1:D:363:GLU:OE2	1:D:366:ARG:NH2	2.43	0.51
1:E:60:VAL:HG21	1:E:349:VAL:CG1	2.40	0.51
1:H:197:GLN:HG2	1:H:211:ARG:HG2	1.91	0.51
1:H:373:LEU:HD12	1:H:386:THR:O	2.10	0.51
1:E:231:LEU:HD21	2:E:401:COA:H1B	1.92	0.51
1:H:342:ALA:HB1	1:H:347:HIS:HB2	1.91	0.51
1:B:197:GLN:HG2	1:B:211:ARG:HG2	1.92	0.51
1:A:210:GLU:HG3	1:E:43:ILE:CG1	2.41	0.51
1:F:197:GLN:HG2	1:F:211:ARG:HG2	1.93	0.51
1:H:212:ARG:HG2	1:H:213:LEU:N	2.25	0.51
1:E:197:GLN:HG2	1:E:211:ARG:HG2	1.93	0.51
1:D:197:GLN:HG2	1:D:211:ARG:HG2	1.93	0.50
1:A:210:GLU:CG	1:E:43:ILE:HG12	2.41	0.50
1:G:197:GLN:HG2	1:G:211:ARG:HG2	1.94	0.50
1:A:265:ALA:CA	1:G:1:MET:CE	2.89	0.50
1:A:60:VAL:HG21	1:A:349:VAL:CG1	2.41	0.49
1:C:67:SER:OG	1:C:68:ASN:ND2	2.44	0.49
1:G:60:VAL:HG21	1:G:349:VAL:CG1	2.42	0.49
1:C:197:GLN:HG2	1:C:211:ARG:HG2	1.94	0.49
1:C:231:LEU:HD21	2:C:400:COA:H1B	1.93	0.49
1:G:363:GLU:OE2	1:G:366:ARG:NH2	2.44	0.49
1:A:215:PHE:CE2	1:E:215:PHE:CE2	3.01	0.49
1:H:60:VAL:HG21	1:H:349:VAL:CG1	2.43	0.49
1:A:247:GLN:HB2	1:A:346:GLY:HA2	1.95	0.49
1:G:231:LEU:HD21	2:G:401:COA:H1B	1.95	0.49
1:H:273:ARG:HG2	1:H:391:ILE:HD11	1.93	0.48
1:B:55:VAL:O	1:B:86:ALA:HA	2.13	0.48
1:E:369:GLN:O	1:E:390:ARG:HD3	2.13	0.48
1:B:13:PRO:HG3	1:B:215:PHE:HA	1.96	0.48
1:B:281:GLY:HA3	1:C:82:GLU:O	2.13	0.48
1:G:13:PRO:HG3	1:G:215:PHE:HA	1.95	0.48
1:B:342:ALA:HB1	1:B:347:HIS:HB2	1.96	0.48
1:E:150:ASN:OD1	1:E:150:ASN:C	2.52	0.47
1:C:13:PRO:HG3	1:C:215:PHE:HA	1.97	0.47
1:E:178[B]:ARG:NH1	1:E:178[B]:ARG:CG	2.54	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:60:VAL:HG21	1:B:349:VAL:CG1	2.44	0.47
1:C:247:GLN:HB2	1:C:346:GLY:HA2	1.96	0.47
1:F:177:GLN:OE1	1:F:181:GLN:NE2	2.44	0.47
1:F:231:LEU:HD21	2:F:400:COA:H1B	1.96	0.47
1:H:10:THR:HG22	1:H:37:VAL:HG13	1.97	0.47
1:F:152:PHE:CE1	1:F:234:VAL:HG11	2.50	0.47
1:E:13:PRO:HG3	1:E:215:PHE:HA	1.96	0.47
1:H:13:PRO:HG3	1:H:215:PHE:HA	1.97	0.47
1:H:247:GLN:HB2	1:H:346:GLY:HA2	1.97	0.47
1:C:101:HIS:HB3	1:C:277:GLN:OE1	2.15	0.46
1:E:152:PHE:CD1	1:E:234:VAL:HG11	2.51	0.46
1:F:10:THR:HG22	1:F:37:VAL:HG13	1.97	0.46
1:G:10:THR:HG22	1:G:37:VAL:HG13	1.97	0.46
1:D:150:ASN:OD1	1:D:150:ASN:C	2.53	0.46
1:E:175:GLU:OE1	1:E:178[A]:ARG:HD3	2.15	0.46
1:E:66:GLN:HE22	1:E:122:ALA:H	1.64	0.46
1:A:13:PRO:HG3	1:A:215:PHE:HA	1.97	0.46
1:E:10:THR:HG22	1:E:37:VAL:HG13	1.98	0.46
1:F:13:PRO:HG3	1:F:215:PHE:HA	1.97	0.46
1:A:325:TRP:CZ3	1:A:331:PRO:HG3	2.51	0.46
1:B:10:THR:HG22	1:B:37:VAL:HG13	1.97	0.46
1:B:18:ASN:N	1:B:124:SER:OG	2.39	0.46
1:G:152:PHE:CD1	1:G:234:VAL:HG11	2.51	0.46
1:G:325:TRP:CZ3	1:G:331:PRO:HG3	2.51	0.46
1:D:369:GLN:O	1:D:390:ARG:HD3	2.16	0.46
1:B:67:SER:OG	1:B:68:ASN:ND2	2.48	0.45
1:A:265:ALA:CA	1:G:1:MET:HE1	2.46	0.45
1:H:230:GLU:OE1	1:H:230:GLU:HA	2.17	0.45
1:A:369:GLN:O	1:A:390:ARG:HD3	2.16	0.45
1:G:66:GLN:HE22	1:G:122:ALA:H	1.64	0.45
1:A:210:GLU:CG	1:E:43:ILE:CG1	2.94	0.45
1:B:101:HIS:HB3	1:B:277:GLN:OE1	2.17	0.45
1:G:236:GLU:O	3:G:403:GOL:H32	2.16	0.45
1:A:231:LEU:HD21	2:A:400:COA:H1B	1.98	0.45
1:A:177:GLN:OE1	1:A:181:GLN:NE2	2.46	0.45
1:D:197:GLN:HG2	1:D:211:ARG:CZ	2.47	0.45
1:H:57:GLY:HA2	1:H:118:CYS:O	2.16	0.45
1:E:18:ASN:N	1:E:124:SER:OG	2.44	0.45
1:H:290:GLY:N	1:H:291:PRO:CD	2.80	0.45
1:B:90:ASP:OD1	1:B:90:ASP:C	2.55	0.45
2:E:401:COA:O9P	2:E:401:COA:H62	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:210:GLU:HG3	1:E:43:ILE:HG13	1.97	0.45
1:D:373:LEU:HD12	1:D:386:THR:O	2.16	0.45
1:H:66:GLN:HE22	1:H:122:ALA:H	1.65	0.44
1:C:150:ASN:C	1:C:150:ASN:OD1	2.54	0.44
1:C:177:GLN:OE1	1:C:181:GLN:NE2	2.46	0.44
1:E:325:TRP:CZ3	1:E:331:PRO:HG3	2.52	0.44
1:G:150:ASN:OD1	1:G:150:ASN:C	2.55	0.44
1:A:82:GLU:O	1:D:281:GLY:HA3	2.17	0.44
1:B:369:GLN:O	1:B:390:ARG:HD3	2.16	0.44
1:B:48[A]:HIS:CE1	1:B:51:ASP:OD1	2.70	0.44
1:B:260:GLU:OE1	1:B:273:ARG:NH2	2.51	0.44
1:A:213:LEU:HD23	1:E:215:PHE:CE2	2.52	0.44
1:C:197:GLN:HG2	1:C:211:ARG:CZ	2.48	0.44
1:C:150:ASN:OD1	1:C:152:PHE:N	2.51	0.44
1:H:369:GLN:O	1:H:390:ARG:HD3	2.17	0.44
1:C:325:TRP:CZ3	1:C:331:PRO:HG3	2.53	0.44
1:A:213:LEU:CD2	1:E:215:PHE:CE2	3.00	0.44
1:B:150:ASN:OD1	1:B:150:ASN:C	2.56	0.43
1:B:347:HIS:CG	1:B:347:HIS:O	2.68	0.43
1:F:150:ASN:C	1:F:150:ASN:OD1	2.56	0.43
1:A:10:THR:HG22	1:A:37:VAL:HG13	2.00	0.43
1:C:66:GLN:HE22	1:C:122:ALA:H	1.66	0.43
1:D:10:THR:HG22	1:D:37:VAL:HG13	1.99	0.43
1:D:290:GLY:N	1:D:291:PRO:CD	2.81	0.43
1:F:369:GLN:O	1:F:390:ARG:HD3	2.18	0.43
1:F:289:ASP:OD1	1:F:324:SER:OG	2.20	0.43
1:A:178:ARG:HH22	1:A:182:ARG:NH1	2.17	0.43
1:A:90:ASP:C	1:A:90:ASP:OD1	2.56	0.43
1:E:290:GLY:N	1:E:291:PRO:CD	2.81	0.43
1:E:247:GLN:HB2	1:E:346:GLY:HA2	2.00	0.43
1:G:0:SER:O	1:G:110:GLY:HA3	2.19	0.43
1:A:210:GLU:HG3	1:E:43:ILE:HG12	2.00	0.43
1:E:177:GLN:OE1	1:E:181:GLN:NE2	2.48	0.43
1:F:197:GLN:HG2	1:F:211:ARG:CZ	2.48	0.43
1:A:363:GLU:OE2	1:A:366:ARG:NH2	2.48	0.43
1:C:152:PHE:CD1	1:C:234:VAL:HG11	2.54	0.43
1:E:90:ASP:C	1:E:90:ASP:OD1	2.57	0.43
1:G:247:GLN:HB2	1:G:346:GLY:HA2	2.01	0.43
1:H:152:PHE:CD1	1:H:234:VAL:HG11	2.53	0.43
1:A:55:VAL:O	1:A:86:ALA:HA	2.18	0.43
1:E:184:TRP:HB3	3:E:402:GOL:H2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:152:PHE:CE1	1:G:234:VAL:HG11	2.53	0.43
1:F:290:GLY:N	1:F:291:PRO:CD	2.82	0.43
1:D:325:TRP:CZ3	1:D:331:PRO:HG3	2.53	0.42
1:B:152:PHE:CD1	1:B:234:VAL:HG11	2.55	0.42
1:B:227:GLY:O	1:B:230:GLU:HB2	2.20	0.42
1:C:57:GLY:HA2	1:C:118:CYS:O	2.20	0.42
1:E:197:GLN:HG2	1:E:211:ARG:CZ	2.49	0.42
1:E:55:VAL:O	1:E:86:ALA:HA	2.20	0.42
1:B:363:GLU:OE2	1:B:366:ARG:NH2	2.52	0.42
1:D:273[A]:ARG:HB3	1:D:391:ILE:HD11	2.01	0.42
1:E:363:GLU:OE2	1:E:366:ARG:NH2	2.52	0.42
1:A:195:PRO:HG3	1:E:215:PHE:CZ	2.52	0.42
1:G:369:GLN:O	1:G:390:ARG:HD3	2.19	0.42
1:C:55:VAL:O	1:C:86:ALA:HA	2.19	0.42
1:E:57:GLY:O	1:E:88:THR:HA	2.19	0.42
1:F:66:GLN:HE22	1:F:122:ALA:H	1.66	0.42
1:A:197:GLN:HG2	1:A:211:ARG:CZ	2.50	0.42
2:A:400:COA:O9P	2:A:400:COA:H141	2.19	0.42
2:H:401:COA:O9P	2:H:401:COA:H62	2.20	0.42
1:B:221:ARG:CZ	2:B:400:COA:H2B	2.50	0.42
1:D:194:SER:O	1:D:195:PRO:C	2.58	0.42
1:F:178[B]:ARG:NH1	1:F:178[B]:ARG:CG	2.52	0.42
1:A:47:LEU:HD13	1:A:258:MET:HE1	2.01	0.41
1:A:273[B]:ARG:NH1	1:A:274:ILE:O	2.54	0.41
1:D:60:VAL:HG21	1:D:349:VAL:HG12	2.02	0.41
1:C:290:GLY:N	1:C:291:PRO:CD	2.84	0.41
1:F:325:TRP:CZ3	1:F:331:PRO:HG3	2.55	0.41
2:G:401:COA:O9P	2:G:401:COA:H62	2.20	0.41
1:C:369:GLN:O	1:C:390:ARG:HD3	2.19	0.41
1:C:216:ARG:NH2	4:C:501:HOH:O	2.44	0.41
1:F:7:VAL:HG11	1:F:268:LEU:HD13	2.02	0.41
1:B:7:VAL:HG11	1:B:268:LEU:HD13	2.01	0.41
1:C:10:THR:HG22	1:C:37:VAL:HG13	2.01	0.41
1:C:233:PRO:HA	1:C:241:THR:HG22	2.02	0.41
1:E:240:HIS:CE1	1:E:320:SER:H	2.31	0.41
1:G:150:ASN:OD1	1:G:152:PHE:N	2.54	0.41
1:A:221:ARG:CZ	2:A:400:COA:H2B	2.50	0.41
1:B:149:PRO:O	1:B:287:HIS:ND1	2.53	0.41
1:B:290:GLY:N	1:B:291:PRO:CD	2.84	0.41
1:G:177:GLN:OE1	1:G:181:GLN:NE2	2.48	0.41
1:E:233:PRO:HA	1:E:241:THR:HG22	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:177:GLN:OE1	1:H:181:GLN:NE2	2.50	0.41
1:B:197:GLN:HG2	1:B:211:ARG:CZ	2.51	0.41
1:G:221:ARG:CZ	2:G:401:COA:H2B	2.51	0.41
1:A:150:ASN:OD1	1:A:152:PHE:N	2.53	0.41
1:A:290:GLY:O	1:A:293:GLN:N	2.54	0.41
1:B:233:PRO:HA	1:B:241:THR:HG22	2.03	0.41
1:C:221:ARG:CZ	2:C:400:COA:H2B	2.51	0.41
2:E:401:COA:O9P	2:E:401:COA:C6P	2.69	0.41
1:G:290:GLY:N	1:G:291:PRO:CD	2.84	0.41
1:A:152:PHE:CD1	1:A:234:VAL:HG11	2.56	0.40
1:F:247:GLN:HB2	1:F:346:GLY:HA2	2.03	0.40
1:G:197:GLN:HG2	1:G:211:ARG:CZ	2.51	0.40
1:H:221:ARG:CZ	2:H:401:COA:H2B	2.51	0.40
1:B:247:GLN:HB2	1:B:346:GLY:HA2	2.03	0.40
2:C:400:COA:O9P	2:C:400:COA:H62	2.20	0.40
1:C:46:GLY:O	1:C:47:LEU:C	2.60	0.40
1:D:221:ARG:CZ	2:D:400:COA:H2B	2.51	0.40
1:E:221:ARG:CZ	2:E:401:COA:H2B	2.51	0.40
1:E:318:PHE:O	1:E:319:ALA:C	2.59	0.40
1:G:149:PRO:O	1:G:287:HIS:ND1	2.52	0.40
1:G:236:GLU:O	3:G:403:GOL:C3	2.70	0.40
1:G:47:LEU:HD13	1:G:258:MET:HE1	2.03	0.40
1:H:150:ASN:OD1	1:H:152:PHE:N	2.54	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	A	389/399 (98%)	375 (96%)	14 (4%)	0	100	100
1	B	390/399 (98%)	376 (96%)	13 (3%)	1 (0%)	44	81

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	388/399 (97%)	373 (96%)	14 (4%)	1 (0%)	44	81
1	D	389/399 (98%)	377 (97%)	11 (3%)	1 (0%)	44	81
1	E	390/399 (98%)	377 (97%)	13 (3%)	0	100	100
1	F	390/399 (98%)	376 (96%)	14 (4%)	0	100	100
1	G	391/399 (98%)	376 (96%)	14 (4%)	1 (0%)	44	81
1	H	393/399 (98%)	375 (95%)	16 (4%)	2 (0%)	32	74
All	All	3120/3192 (98%)	3005 (96%)	109 (4%)	6 (0%)	51	86

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	0	SER
1	H	1	MET
1	D	130	ALA
1	B	130	ALA
1	C	130	ALA
1	H	-1	GLY

### 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	296/302 (98%)	289 (98%)	7 (2%)	54	85
1	B	297/302 (98%)	291 (98%)	6 (2%)	60	87
1	C	295/302 (98%)	287 (97%)	8 (3%)	50	82
1	D	296/302 (98%)	289 (98%)	7 (2%)	54	85
1	E	297/302 (98%)	291 (98%)	6 (2%)	60	87
1	F	297/302 (98%)	293 (99%)	4 (1%)	73	92
1	G	297/302 (98%)	291 (98%)	6 (2%)	60	87
1	H	299/302 (99%)	291 (97%)	8 (3%)	50	82
All	All	2374/2416 (98%)	2322 (98%)	52 (2%)	59	86

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	VAL
1	A	197	GLN
1	A	215	PHE
1	A	273[A]	ARG
1	A	273[B]	ARG
1	A	327	ARG
1	A	351	CYS
1	B	37	VAL
1	B	153	GLU
1	B	197	GLN
1	B	215	PHE
1	B	327	ARG
1	B	351	CYS
1	C	37	VAL
1	C	153	GLU
1	C	178	ARG
1	C	197	GLN
1	C	215	PHE
1	C	327	ARG
1	C	351	CYS
1	C	375	THR
1	D	37	VAL
1	D	197	GLN
1	D	215	PHE
1	D	327	ARG
1	D	351	CYS
1	D	354	SER
1	D	375	THR
1	E	37	VAL
1	E	197	GLN
1	E	215	PHE
1	E	327	ARG
1	E	351	CYS
1	E	375	THR
1	F	37	VAL
1	F	197	GLN
1	F	215	PHE
1	F	327	ARG
1	G	37	VAL
1	G	150	ASN
1	G	197	GLN
1	G	215	PHE

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Mol	Chain	Res	Type
1	G	327	ARG
1	G	351	CYS
1	H	3[A]	TYR
1	H	3[B]	TYR
1	H	10	THR
1	H	37	VAL
1	H	197	GLN
1	H	215	PHE
1	H	273	ARG
1	H	327	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (41) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	GLN
1	A	68	ASN
1	A	100	ASN
1	A	240	HIS
1	A	369	GLN
1	B	66	GLN
1	B	68	ASN
1	B	100	ASN
1	B	240	HIS
1	B	369	GLN
1	C	66	GLN
1	C	68	ASN
1	C	100	ASN
1	C	240	HIS
1	C	369	GLN
1	D	66	GLN
1	D	68	ASN
1	D	100	ASN
1	D	240	HIS
1	D	369	GLN
1	E	44	GLN
1	E	66	GLN
1	E	68	ASN
1	E	100	ASN
1	E	240	HIS
1	E	369	GLN
1	F	66	GLN
1	F	68	ASN

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Mol	Chain	Res	Type
1	F	100	ASN
1	F	240	HIS
1	F	369	GLN
1	G	66	GLN
1	G	68	ASN
1	G	100	ASN
1	G	240	HIS
1	G	369	GLN
1	H	66	GLN
1	H	68	ASN
1	H	100	ASN
1	H	240	HIS
1	H	369	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

8 non-standard protein/DNA/RNA residues are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
1	OAS	A	93	1	8,8,9	1.53	2 (25%)	6,9,11	2.06	2 (33%)
1	OAS	B	93	1	8,8,9	1.34	1 (12%)	6,9,11	1.94	1 (16%)
1	OAS	C	93	1	8,8,9	1.38	1 (12%)	6,9,11	1.73	1 (16%)
1	OAS	D	93	1	8,8,9	1.29	1 (12%)	6,9,11	1.67	1 (16%)
1	OAS	E	93	1	8,8,9	1.27	1 (12%)	6,9,11	2.26	1 (16%)
1	OAS	F	93	1	8,8,9	1.44	1 (12%)	6,9,11	2.02	2 (33%)
1	OAS	G	93	1	8,8,9	1.44	1 (12%)	6,9,11	2.12	2 (33%)
1	OAS	H	93	1	8,8,9	1.47	2 (25%)	6,9,11	2.13	3 (50%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	OAS	A	93	1	-	0/5/7/9	0/0/0/0
1	OAS	B	93	1	-	0/5/7/9	0/0/0/0
1	OAS	C	93	1	-	0/5/7/9	0/0/0/0
1	OAS	D	93	1	-	0/5/7/9	0/0/0/0
1	OAS	E	93	1	-	0/5/7/9	0/0/0/0
1	OAS	F	93	1	-	0/5/7/9	0/0/0/0
1	OAS	G	93	1	-	0/5/7/9	0/0/0/0
1	OAS	H	93	1	-	0/5/7/9	0/0/0/0

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	93	OAS	CA-C	-2.67	1.46	1.50
1	A	93	OAS	CA-C	-2.29	1.47	1.50
1	H	93	OAS	OG-C1A	2.94	1.48	1.33
1	E	93	OAS	OG-C1A	3.01	1.48	1.33
1	D	93	OAS	OG-C1A	3.16	1.49	1.33
1	A	93	OAS	OG-C1A	3.30	1.49	1.33
1	B	93	OAS	OG-C1A	3.36	1.50	1.33
1	C	93	OAS	OG-C1A	3.42	1.50	1.33
1	F	93	OAS	OG-C1A	3.55	1.51	1.33
1	G	93	OAS	OG-C1A	3.73	1.51	1.33

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	93	OAS	OAC-C1A-C2A	-2.26	116.51	124.82
1	A	93	OAS	O-C-CA	-2.24	118.83	125.02
1	H	93	OAS	O-C-CA	-2.14	119.11	125.02
1	G	93	OAS	OAC-C1A-C2A	-2.10	117.09	124.82
1	H	93	OAS	OG-C1A-C2A	2.07	121.73	112.44
1	C	93	OAS	CB-OG-C1A	2.99	124.78	117.17
1	D	93	OAS	CB-OG-C1A	3.29	125.55	117.17
1	F	93	OAS	CB-OG-C1A	3.76	126.73	117.17
1	B	93	OAS	CB-OG-C1A	3.99	127.31	117.17
1	H	93	OAS	CB-OG-C1A	3.99	127.32	117.17
1	A	93	OAS	CB-OG-C1A	4.00	127.36	117.17
1	G	93	OAS	CB-OG-C1A	4.18	127.80	117.17
1	E	93	OAS	CB-OG-C1A	5.13	130.23	117.17

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

12 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	400	-	43,50,50	0.84	1 (2%)	48,75,75	1.95	6 (12%)
2	COA	B	400	-	43,50,50	0.88	1 (2%)	48,75,75	1.93	5 (10%)
2	COA	C	400	-	43,50,50	0.86	1 (2%)	48,75,75	1.89	5 (10%)
2	COA	D	400	-	43,50,50	0.89	1 (2%)	48,75,75	1.99	6 (12%)
2	COA	E	401	-	43,50,50	0.87	1 (2%)	48,75,75	1.99	8 (16%)
3	GOL	E	402	-	5,5,5	0.60	0	5,5,5	1.76	1 (20%)
2	COA	F	400	-	43,50,50	0.87	1 (2%)	48,75,75	1.91	5 (10%)
2	COA	G	401	-	43,50,50	0.83	1 (2%)	48,75,75	1.93	6 (12%)
3	GOL	G	402	-	5,5,5	0.59	0	5,5,5	1.63	1 (20%)
3	GOL	G	403	-	5,5,5	0.81	0	5,5,5	1.17	1 (20%)
2	COA	H	401	-	43,50,50	0.85	1 (2%)	48,75,75	1.90	5 (10%)
3	GOL	H	402	-	5,5,5	0.66	0	5,5,5	1.39	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	COA	A	400	-	-	0/44/64/64	0/3/3/3
2	COA	B	400	-	-	0/44/64/64	0/3/3/3
2	COA	C	400	-	-	0/44/64/64	0/3/3/3
2	COA	D	400	-	-	0/44/64/64	0/3/3/3
2	COA	E	401	-	-	0/44/64/64	0/3/3/3
3	GOL	E	402	-	-	0/4/4/4	0/0/0/0
2	COA	F	400	-	-	0/44/64/64	0/3/3/3
2	COA	G	401	-	-	0/44/64/64	0/3/3/3
3	GOL	G	402	-	-	0/4/4/4	0/0/0/0
3	GOL	G	403	-	-	0/4/4/4	0/0/0/0
2	COA	H	401	-	-	0/44/64/64	0/3/3/3
3	GOL	H	402	-	-	0/4/4/4	0/0/0/0

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	401	COA	C5A-C4A	2.94	1.47	1.40
2	B	400	COA	C5A-C4A	3.00	1.47	1.40
2	G	401	COA	C5A-C4A	3.01	1.47	1.40
2	A	400	COA	C5A-C4A	3.02	1.47	1.40
2	D	400	COA	C5A-C4A	3.07	1.47	1.40
2	H	401	COA	C5A-C4A	3.12	1.47	1.40
2	F	400	COA	C5A-C4A	3.15	1.47	1.40
2	C	400	COA	C5A-C4A	3.16	1.47	1.40

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	401	COA	N3A-C2A-N1A	-8.10	121.80	128.86
2	D	400	COA	C4B-O4B-C1B	-7.87	101.39	109.77
2	A	400	COA	C4B-O4B-C1B	-7.73	101.54	109.77
2	H	401	COA	N3A-C2A-N1A	-7.61	122.23	128.86
2	F	400	COA	C4B-O4B-C1B	-7.58	101.70	109.77
2	D	400	COA	N3A-C2A-N1A	-7.53	122.30	128.86
2	A	400	COA	N3A-C2A-N1A	-7.53	122.30	128.86
2	C	400	COA	C4B-O4B-C1B	-7.50	101.78	109.77
2	G	401	COA	C4B-O4B-C1B	-7.42	101.88	109.77
2	B	400	COA	N3A-C2A-N1A	-7.39	122.42	128.86
2	B	400	COA	C4B-O4B-C1B	-7.38	101.92	109.77
2	F	400	COA	N3A-C2A-N1A	-7.33	122.48	128.86
2	E	401	COA	C4B-O4B-C1B	-7.32	101.98	109.77
2	H	401	COA	C4B-O4B-C1B	-7.31	101.98	109.77
2	C	400	COA	N3A-C2A-N1A	-7.26	122.53	128.86

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	401	COA	N3A-C2A-N1A	-7.20	122.59	128.86
2	B	400	COA	C4A-C5A-N7A	-3.92	105.62	109.41
2	D	400	COA	C4A-C5A-N7A	-3.89	105.65	109.41
2	G	401	COA	C4A-C5A-N7A	-3.74	105.80	109.41
2	F	400	COA	C4A-C5A-N7A	-3.64	105.89	109.41
2	A	400	COA	C4A-C5A-N7A	-3.56	105.97	109.41
2	C	400	COA	C4A-C5A-N7A	-3.52	106.01	109.41
2	H	401	COA	C4A-C5A-N7A	-3.40	106.12	109.41
2	E	401	COA	C4A-C5A-N7A	-3.32	106.21	109.41
2	B	400	COA	C1B-N9A-C4A	-3.20	121.11	126.64
2	F	400	COA	C1B-N9A-C4A	-3.17	121.15	126.64
2	D	400	COA	C1B-N9A-C4A	-3.16	121.17	126.64
2	G	401	COA	C1B-N9A-C4A	-3.13	121.23	126.64
2	E	401	COA	C1B-N9A-C4A	-3.09	121.29	126.64
2	H	401	COA	C1B-N9A-C4A	-3.04	121.38	126.64
2	C	400	COA	C1B-N9A-C4A	-3.04	121.39	126.64
2	A	400	COA	C1B-N9A-C4A	-3.03	121.41	126.64
2	E	401	COA	O5P-C5P-C6P	-2.57	117.19	122.01
2	G	401	COA	OAP-CAP-CBP	-2.25	104.95	110.25
2	B	400	COA	O5P-C5P-C6P	-2.22	117.85	122.01
2	A	400	COA	O5P-C5P-C6P	-2.18	117.91	122.01
2	G	401	COA	O5P-C5P-C6P	-2.14	118.00	122.01
3	G	403	GOL	O3-C3-C2	2.04	120.34	110.07
2	D	400	COA	O6A-CCP-CBP	2.08	113.89	110.55
2	E	401	COA	CEP-CBP-CCP	2.10	111.45	108.37
2	C	400	COA	C2A-N1A-C6A	2.17	122.56	118.77
2	E	401	COA	O5P-C5P-N4P	2.20	127.17	122.97
2	A	400	COA	C2A-N1A-C6A	2.20	122.62	118.77
2	F	400	COA	C2A-N1A-C6A	2.22	122.65	118.77
2	E	401	COA	C2A-N1A-C6A	2.22	122.65	118.77
2	H	401	COA	C2A-N1A-C6A	2.25	122.70	118.77
2	D	400	COA	C2A-N1A-C6A	2.30	122.79	118.77
3	G	402	GOL	O3-C3-C2	2.36	121.95	110.07
3	E	402	GOL	O3-C3-C2	2.70	123.65	110.07

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

11 monomers are involved in 25 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	400	COA	3	0
2	B	400	COA	2	0
2	C	400	COA	3	0
2	D	400	COA	2	0
2	E	401	COA	4	0
3	E	402	GOL	1	0
2	F	400	COA	1	0
2	G	401	COA	3	0
3	G	403	GOL	2	0
2	H	401	COA	3	0
3	H	402	GOL	1	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	390/399 (97%)	-0.56	1 (0%) 93 82	22, 35, 66, 185	1 (0%)
1	B	390/399 (97%)	-0.42	1 (0%) 93 82	22, 37, 67, 147	1 (0%)
1	C	390/399 (97%)	-0.53	1 (0%) 93 82	20, 39, 67, 173	1 (0%)
1	D	390/399 (97%)	-0.44	3 (0%) 86 64	20, 36, 61, 167	1 (0%)
1	E	391/399 (97%)	-0.44	1 (0%) 93 82	21, 36, 68, 132	1 (0%)
1	F	390/399 (97%)	-0.35	0 100 100	20, 37, 65, 157	1 (0%)
1	G	392/399 (98%)	-0.36	5 (1%) 77 51	20, 36, 71, 179	1 (0%)
1	H	393/399 (98%)	-0.33	2 (0%) 90 74	21, 46, 74, 124	1 (0%)
All	All	3126/3192 (97%)	-0.43	14 (0%) 92 77	20, 37, 70, 185	8 (0%)

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	-2	HIS	2.6
1	A	204	GLN	2.6
1	G	204	GLN	2.5
1	G	238	GLY	2.4
1	B	204	GLN	2.4
1	D	204	GLN	2.3
1	G	232	LYS	2.3
1	E	215	PHE	2.3
1	G	203	GLU	2.2
1	G	205	ASN	2.1
1	D	205	ASN	2.1
1	D	208	THR	2.1
1	C	208	THR	2.1
1	H	232	LYS	2.0

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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
1	OAS	C	93	9/10	0.95	0.16	-	23,24,28,30	0
1	OAS	D	93	9/10	0.97	0.20	-	26,27,31,33	0
1	OAS	E	93	9/10	0.95	0.15	-	26,30,32,32	0
1	OAS	F	93	9/10	0.93	0.20	-	27,29,34,36	0
1	OAS	A	93	9/10	0.93	0.25	-	29,30,34,34	0
1	OAS	B	93	9/10	0.94	0.22	-	25,26,30,33	0
1	OAS	G	93	9/10	0.89	0.24	-	24,28,33,34	0
1	OAS	H	93	9/10	0.95	0.15	-	30,33,35,36	0

## 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. 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
2	COA	B	400	48/48	0.69	0.52	7.75	90,115,175,188	0
2	COA	A	400	48/48	0.71	0.38	6.44	88,125,172,182	0
2	COA	C	400	48/48	0.73	0.39	5.03	82,112,197,204	0
2	COA	D	400	48/48	0.73	0.47	4.66	77,94,138,142	0
2	COA	H	401	48/48	0.59	0.46	4.12	70,148,201,210	0
2	COA	G	401	48/48	0.70	0.51	4.01	79,132,172,184	0
2	COA	E	401	48/48	0.64	0.38	3.63	74,113,176,183	0
2	COA	F	400	48/48	0.78	0.43	2.74	63,106,147,159	0
3	GOL	E	402	6/6	0.70	0.38	2.38	47,57,60,61	0
3	GOL	G	403	6/6	0.90	0.50	2.05	58,66,70,72	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	GOL	H	402	6/6	0.84	0.32	1.72	70,72,75,78	0
3	GOL	G	402	6/6	0.75	0.33	-	57,64,67,73	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.