



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:42 AM GMT

PDB ID : 2Y1W  
Title : CRYSTAL STRUCTURE OF COACTIVATOR ASSOCIATED ARGININE METHYLTRANSFERASE 1 (CARM1) IN COMPLEX WITH SINEFUNGIN AND INDOLE INHIBITOR  
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Deposited on : 2010-12-10  
Resolution : 2.10 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 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
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) : trunk26865

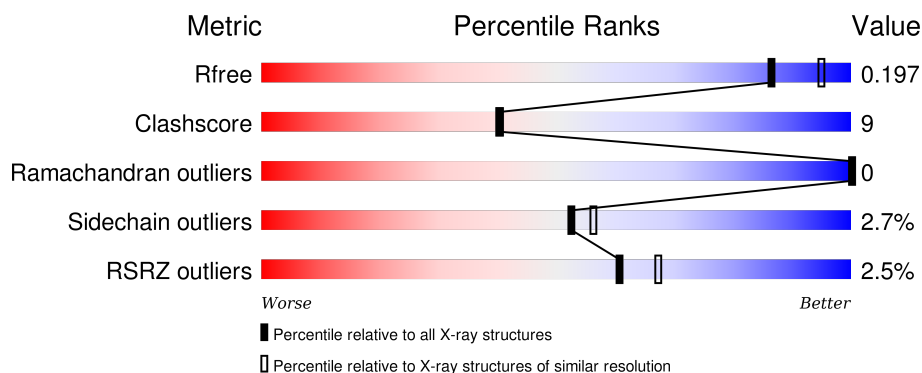
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.10 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	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (2.10-2.10)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	348	<div> <div>4%</div> <div>84%</div> <div>14%</div> <div>..</div> </div>
1	B	348	<div> <div>2%</div> <div>81%</div> <div>17%</div> <div>..</div> </div>
1	C	348	<div> <div>2%</div> <div>79%</div> <div>19%</div> <div>..</div> </div>
1	D	348	<div> <div>2%</div> <div>84%</div> <div>14%</div> <div>..</div> </div>

## 2 Entry composition [i](#)

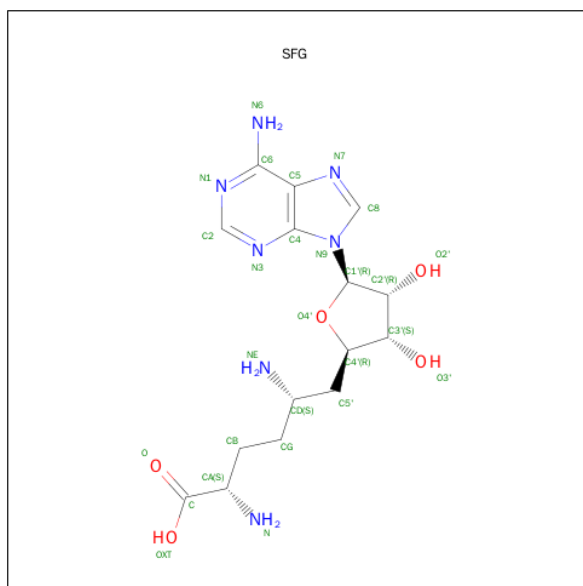
There are 4 unique types of molecules in this entry. The entry contains 11580 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 HISTONE-ARGININE METHYLTRANSFERASE CARM1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	344	Total	C	N	O	S	0	0	1
			2752	1776	455	507	14			
1	B	344	Total	C	N	O	S	0	0	1
			2752	1776	455	507	14			
1	C	344	Total	C	N	O	S	0	0	1
			2752	1776	455	507	14			
1	D	344	Total	C	N	O	S	0	0	1
			2752	1776	455	507	14			

- Molecule 2 is SINEFUNGIN (three-letter code: SFG) (formula: C<sub>15</sub>H<sub>23</sub>N<sub>7</sub>O<sub>5</sub>).



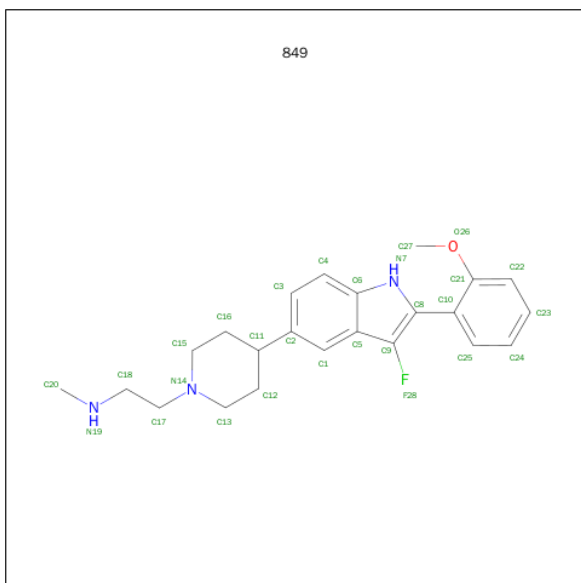
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			27	15	7	5		
2	B	1	Total	C	N	O	0	0
			27	15	7	5		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	C	1	Total	C	N	O	0	0
			27	15	7	5		
2	D	1	Total	C	N	O	0	0
			27	15	7	5		

- Molecule 3 is 2-{4-[3-FLUORO-2-(2-METHOXYPHENYL)-1H-INDOL-5-YL] PIPERIDIN-1-YL}-N-METHYLETHANAMINE (three-letter code: 849) (formula: C<sub>23</sub>H<sub>28</sub>FN<sub>3</sub>O).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	0	0
			28	23	1	3	1		
3	B	1	Total	C	F	N	O	0	0
			28	23	1	3	1		
3	C	1	Total	C	F	N	O	0	0
			28	23	1	3	1		
3	D	1	Total	C	F	N	O	0	0
			28	23	1	3	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	104	Total	O	0	0
			104	104		
4	B	70	Total	O	0	0
			70	70		
4	C	83	Total	O	0	0
			83	83		

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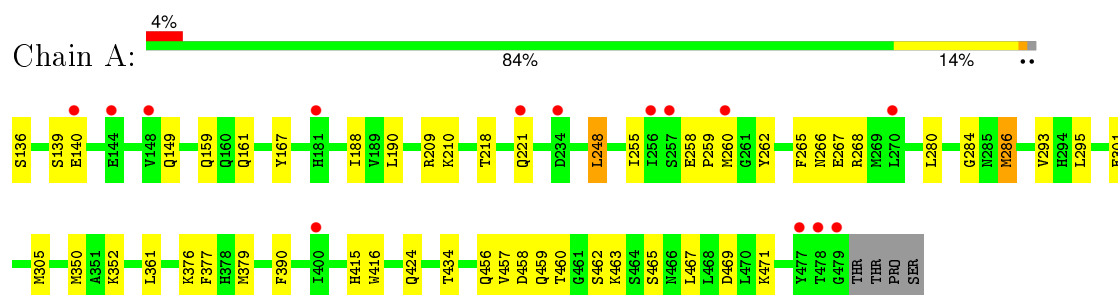
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	95	Total	O	0	0
			95	95		

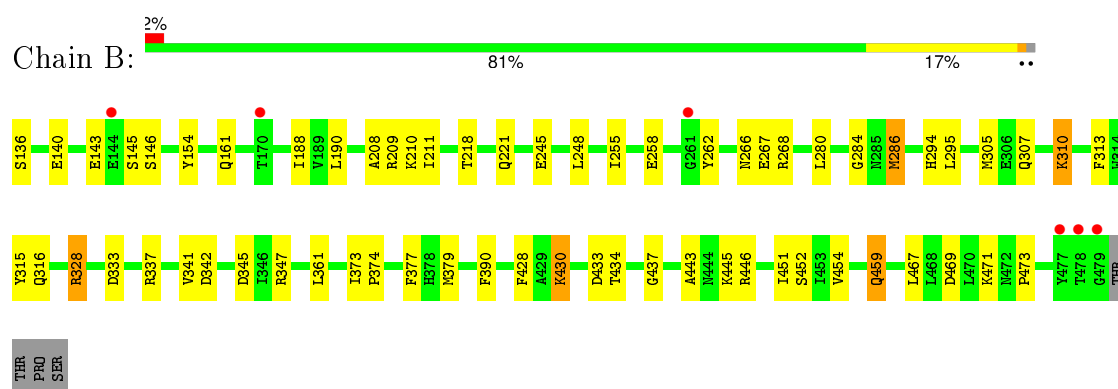
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

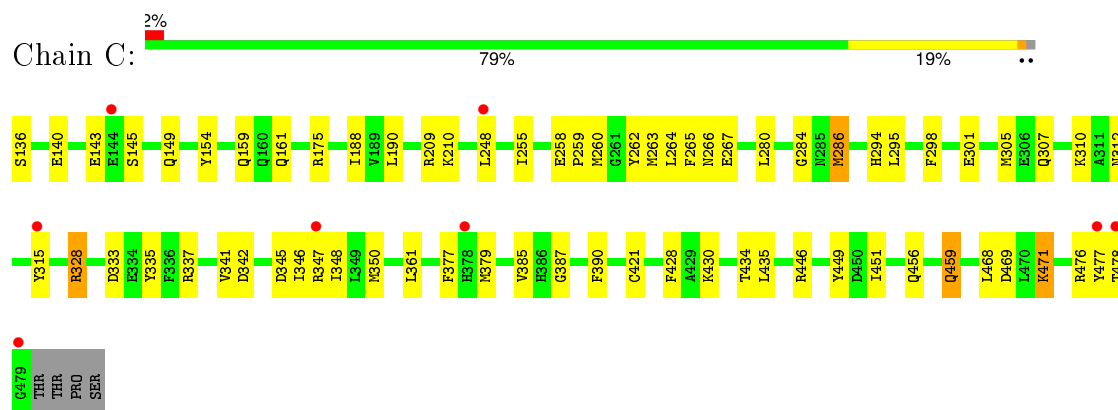
#### • Molecule 1: HISTONE-ARGININE METHYLTRANSFERASE CARM1



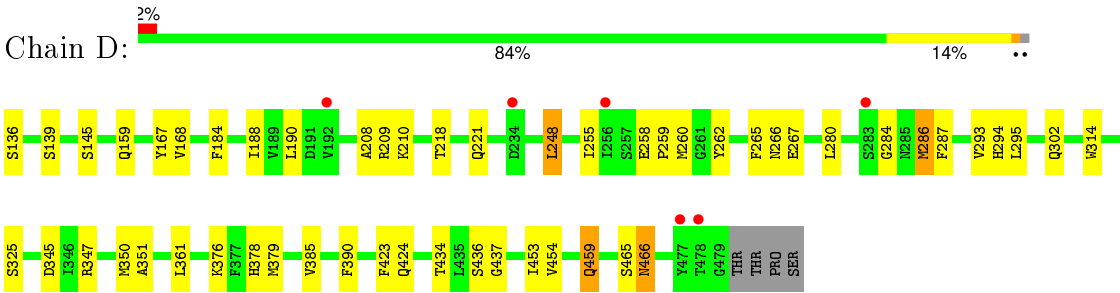
#### • Molecule 1: HISTONE-ARGININE METHYLTRANSFERASE CARM1



#### • Molecule 1: HISTONE-ARGININE METHYLTRANSFERASE CARM1



#### • Molecule 1: HISTONE-ARGININE METHYLTRANSFERASE CARM1



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 2 <sub>1</sub> 2 <sub>1</sub> 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	74.90 Å 98.47 Å 207.18 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.88 – 2.10 29.88 – 2.10	Depositor EDS
% Data completeness (in resolution range)	85.2 (29.88-2.10) 85.4 (29.88-2.10)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.60 (at 2.10 Å)	Xtriage
Refinement program	CNS NULL	Depositor
R, $R_{free}$	0.205 , 0.244 0.202 , 0.197	Depositor DCC
$R_{free}$ test set	3832 reflections (5.24%)	DCC
Wilson B-factor (Å <sup>2</sup> )	32.3	Xtriage
Anisotropy	0.474	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 43.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 76991 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	11580	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	36.0	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 53.58 % 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 4.1314e-05. 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, 0.375 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: SFG, 849

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	0.55	0/2822	0.62	0/3825
1	B	0.54	0/2822	0.60	0/3825
1	C	0.55	0/2822	0.60	0/3825
1	D	0.57	0/2822	0.60	0/3825
All	All	0.55	0/11288	0.61	0/15300

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	2752	0	2699	45	0
1	B	2752	0	2699	54	0
1	C	2752	0	2699	54	0
1	D	2752	0	2699	41	0
2	A	27	0	22	2	0
2	B	27	0	22	1	0
2	C	27	0	22	1	0
2	D	27	0	22	2	0
3	A	28	0	28	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	28	0	28	4	0
3	C	28	0	28	5	0
3	D	28	0	28	5	0
4	A	104	0	0	5	0
4	B	70	0	0	1	0
4	C	83	0	0	3	0
4	D	95	0	0	4	0
All	All	11580	0	10996	194	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 (194) 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:460:THR:HG23	1:A:462:SER:H	1.28	0.97
1:A:434:THR:H	1:A:459:GLN:HE22	1.11	0.96
1:D:459:GLN:H	1:D:459:GLN:NE2	1.74	0.84
1:A:457:VAL:HB	1:A:460:THR:HG22	1.60	0.82
1:B:307:GLN:HE22	1:B:310:LYS:NZ	1.77	0.82
1:C:459:GLN:NE2	1:C:459:GLN:H	1.79	0.80
1:C:459:GLN:HE21	1:C:459:GLN:H	1.30	0.78
1:A:434:THR:H	1:A:459:GLN:NE2	1.81	0.77
1:B:328:ARG:HH22	1:C:328:ARG:NH2	1.83	0.76
1:B:333:ASP:O	1:B:337:ARG:HG3	1.86	0.74
1:D:459:GLN:H	1:D:459:GLN:HE21	1.36	0.72
1:D:255:ILE:HB	1:D:286:MET:HE2	1.73	0.70
1:A:352:LYS:HE3	4:A:2056:HOH:O	1.93	0.69
1:A:457:VAL:HB	1:A:460:THR:CG2	2.23	0.69
1:B:280:LEU:HD11	1:B:284:GLY:HA3	1.75	0.68
1:D:255:ILE:HB	1:D:286:MET:CE	2.24	0.66
1:B:295:LEU:HG	1:B:390:PHE:CE2	2.31	0.66
1:A:286:MET:HG3	1:A:361:LEU:HD23	1.77	0.66
1:B:307:GLN:HE22	1:B:310:LYS:HZ1	1.44	0.65
1:A:136:SER:N	1:A:139:SER:HG	1.93	0.65
1:C:280:LEU:HD11	1:C:284:GLY:HA3	1.79	0.65
1:B:255:ILE:HB	1:B:286:MET:CE	2.27	0.65
1:B:345:ASP:OD2	1:B:347:ARG:HD3	1.98	0.64
1:C:255:ILE:HB	1:C:286:MET:CE	2.29	0.63
1:D:280:LEU:HD11	1:D:284:GLY:HA3	1.80	0.63
1:C:295:LEU:HG	1:C:390:PHE:CE2	2.33	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:286:MET:HG3	1:D:361:LEU:HD23	1.81	0.62
1:B:469:ASP:OD1	1:B:471:LYS:HE2	1.98	0.62
1:B:313:PHE:O	1:B:316:GLN:HG2	2.00	0.62
1:D:459:GLN:NE2	1:D:459:GLN:N	2.47	0.62
1:C:307:GLN:NE2	1:C:307:GLN:HA	2.15	0.62
2:C:1001:SFG:NE	3:C:1005:849:N19	2.47	0.62
1:A:457:VAL:CB	1:A:460:THR:HG22	2.29	0.61
1:B:451:ILE:O	1:B:467:LEU:HA	2.00	0.61
1:C:255:ILE:HB	1:C:286:MET:HE2	1.83	0.61
1:A:280:LEU:HD11	1:A:284:GLY:HA3	1.81	0.61
1:D:424:GLN:HB3	4:D:2088:HOH:O	2.00	0.61
1:C:188:ILE:HD11	1:C:209:ARG:NH2	2.15	0.60
1:D:255:ILE:HD12	1:D:286:MET:CE	2.32	0.59
1:A:262:TYR:CE2	3:A:1005:849:H3	2.36	0.59
1:D:255:ILE:HD12	1:D:286:MET:HE3	1.84	0.59
1:C:333:ASP:O	1:C:337:ARG:HG3	2.02	0.59
1:C:342:ASP:OD2	1:C:476:ARG:HD3	2.03	0.58
1:A:255:ILE:HD12	1:A:286:MET:CE	2.32	0.58
2:D:1001:SFG:NE	3:D:1005:849:N19	2.50	0.58
1:C:307:GLN:HE22	1:C:310:LYS:NZ	2.02	0.58
1:D:423:PHE:HA	1:D:466:ASN:OD1	2.03	0.58
1:D:345:ASP:OD1	1:D:347:ARG:HB2	2.04	0.58
1:A:188:ILE:HD11	1:A:209:ARG:NH2	2.18	0.58
1:B:188:ILE:HD11	1:B:209:ARG:NH2	2.19	0.58
1:A:268:ARG:HA	4:A:2092:HOH:O	2.03	0.57
1:D:466:ASN:ND2	1:D:466:ASN:N	2.51	0.57
1:A:255:ILE:HB	1:A:286:MET:CE	2.34	0.57
1:A:460:THR:HG21	4:A:2089:HOH:O	2.05	0.57
1:A:301:GLU:O	1:A:305:MET:HG3	2.05	0.57
1:D:188:ILE:HD11	1:D:209:ARG:NH2	2.19	0.57
1:B:428:PHE:CE2	1:B:430:LYS:HD3	2.39	0.57
1:C:459:GLN:NE2	1:C:459:GLN:N	2.50	0.56
1:C:136:SER:O	1:C:140:GLU:HG2	2.05	0.56
1:A:471:LYS:HG2	4:A:2095:HOH:O	2.05	0.56
1:B:377:PHE:O	1:B:434:THR:HA	2.06	0.56
1:B:255:ILE:HB	1:B:286:MET:HE2	1.88	0.55
1:C:451:ILE:HB	1:C:468:LEU:HB2	1.89	0.55
1:B:143:GLU:HB2	1:B:446:ARG:HH11	1.71	0.55
1:C:345:ASP:OD1	1:C:347:ARG:HG3	2.07	0.55
1:D:266:ASN:O	1:D:267:GLU:HB2	2.07	0.55
1:A:255:ILE:HB	1:A:286:MET:HE2	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:262:TYR:CE2	3:C:1005:849:H3	2.41	0.55
1:C:255:ILE:HD12	1:C:286:MET:CE	2.36	0.54
1:B:266:ASN:O	1:B:267:GLU:HB2	2.08	0.54
1:A:255:ILE:HD12	1:A:286:MET:HE3	1.90	0.54
1:D:262:TYR:CE2	3:D:1005:849:H3	2.43	0.54
1:A:424:GLN:HG3	1:A:465:SER:O	2.07	0.54
1:A:159:GLN:NE2	4:A:2005:HOH:O	2.40	0.53
1:B:305:MET:HE1	1:C:305:MET:HE1	1.90	0.53
1:B:262:TYR:CE2	3:B:1005:849:H3	2.44	0.53
1:C:377:PHE:O	1:C:434:THR:HA	2.09	0.53
1:C:421:CYS:SG	1:C:468:LEU:HD22	2.48	0.53
1:C:161:GLN:NE2	4:C:2005:HOH:O	2.41	0.52
1:A:136:SER:N	1:A:139:SER:OG	2.43	0.52
2:B:1001:SFG:NE	3:B:1005:849:N19	2.58	0.52
1:D:351:ALA:HB2	1:D:379:MET:HE1	1.91	0.52
1:A:456:GLN:HB3	1:A:463:LYS:HG2	1.91	0.51
1:A:377:PHE:HB2	1:A:379:MET:CE	2.40	0.51
1:D:159:GLN:NE2	4:D:2006:HOH:O	2.37	0.51
1:C:266:ASN:O	1:C:267:GLU:HB2	2.11	0.51
1:B:286:MET:HG3	1:B:361:LEU:HD23	1.92	0.51
1:A:265:PHE:CE2	1:A:293:VAL:HG21	2.46	0.51
1:B:307:GLN:HA	1:B:307:GLN:NE2	2.26	0.51
1:C:255:ILE:HD12	1:C:286:MET:HE3	1.93	0.50
1:D:424:GLN:N	4:D:2088:HOH:O	2.42	0.50
1:B:145:SER:HB3	1:D:145:SER:HA	1.92	0.50
1:B:452:SER:HA	1:B:467:LEU:HD23	1.93	0.50
1:D:295:LEU:HG	1:D:390:PHE:CE2	2.47	0.50
1:A:262:TYR:CZ	3:A:1005:849:H3	2.47	0.50
1:C:428:PHE:HZ	1:C:430:LYS:HE2	1.76	0.50
1:A:457:VAL:CG1	1:A:460:THR:HG22	2.43	0.49
1:C:469:ASP:OD1	1:C:471:LYS:HG2	2.11	0.49
1:A:259:PRO:HD2	1:A:260:MET:SD	2.53	0.49
1:D:190:LEU:HD13	1:D:248:LEU:HD11	1.94	0.49
1:C:286:MET:HG3	1:C:361:LEU:HD23	1.95	0.49
1:B:268:ARG:CZ	1:B:443:ALA:HB1	2.42	0.49
1:A:434:THR:OG1	1:A:458:ASP:HB2	2.13	0.49
1:B:315:TYR:HB2	1:B:328:ARG:HG3	1.95	0.48
1:C:259:PRO:HA	3:C:1005:849:H201	1.94	0.48
1:C:341:VAL:O	1:C:342:ASP:HB3	2.12	0.48
1:C:428:PHE:CZ	1:C:430:LYS:HE2	2.49	0.48
1:B:345:ASP:CG	1:B:347:ARG:HD3	2.33	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:416:TRP:NE1	3:A:1005:849:H151	2.28	0.48
1:C:477:TYR:O	1:C:478:THR:HG23	2.14	0.48
1:C:307:GLN:HE22	1:C:310:LYS:CE	2.27	0.48
1:D:280:LEU:CD1	1:D:284:GLY:HA3	2.44	0.47
1:B:154:TYR:OH	3:B:1005:849:H181	2.15	0.47
1:C:430:LYS:HE3	1:D:167:TYR:CD2	2.50	0.47
1:B:146:SER:HB2	1:B:446:ARG:CZ	2.45	0.47
1:B:143:GLU:HB2	1:B:446:ARG:NH1	2.30	0.47
1:D:376:LYS:HG3	1:D:436:SER:HB3	1.96	0.47
1:B:255:ILE:HD12	1:B:286:MET:CE	2.45	0.47
1:B:143:GLU:CB	1:B:446:ARG:HH11	2.28	0.46
1:B:161:GLN:NE2	4:B:2007:HOH:O	2.48	0.46
1:C:435:LEU:HD12	1:C:456:GLN:O	2.16	0.46
1:A:469:ASP:OD1	1:A:471:LYS:HG2	2.16	0.46
2:A:1001:SFG:NE	3:A:1005:849:N19	2.64	0.46
1:B:307:GLN:HE22	1:B:310:LYS:CE	2.28	0.45
1:B:315:TYR:CA	1:B:328:ARG:HG3	2.47	0.45
1:C:468:LEU:HD12	1:C:468:LEU:N	2.31	0.45
1:C:159:GLN:NE2	4:C:2004:HOH:O	2.37	0.45
1:B:190:LEU:HD13	1:B:248:LEU:HD11	1.98	0.45
1:D:314:TRP:O	1:D:325:SER:HA	2.16	0.45
1:B:471:LYS:C	1:B:473:PRO:HD3	2.37	0.45
1:D:350:MET:SD	1:D:385:VAL:HG22	2.57	0.44
1:C:312:ASN:O	1:C:315:TYR:HB3	2.17	0.44
1:B:377:PHE:HB2	1:B:379:MET:CE	2.47	0.44
1:B:315:TYR:HA	1:B:328:ARG:HG3	1.99	0.44
1:C:262:TYR:CZ	3:C:1005:849:H3	2.53	0.44
1:B:373:ILE:HA	1:B:374:PRO:HD3	1.84	0.44
1:B:218:THR:O	1:B:221:GLN:HG2	2.18	0.44
1:A:259:PRO:HA	3:A:1005:849:H201	2.00	0.43
1:B:459:GLN:H	1:B:459:GLN:NE2	2.15	0.43
1:A:266:ASN:O	1:A:267:GLU:HB2	2.18	0.43
1:D:168:VAL:HG11	4:D:2079:HOH:O	2.17	0.43
1:D:259:PRO:HD2	1:D:260:MET:SD	2.58	0.43
1:C:387:GLY:HA3	1:C:421:CYS:O	2.19	0.43
1:B:262:TYR:CZ	3:B:1005:849:H3	2.53	0.43
1:B:294:HIS:O	1:B:390:PHE:HA	2.19	0.43
1:A:255:ILE:HD12	1:A:286:MET:HE1	1.98	0.43
1:A:159:GLN:NE2	1:A:159:GLN:HA	2.34	0.43
1:B:341:VAL:O	1:B:342:ASP:HB3	2.19	0.43
1:A:190:LEU:HD13	1:A:248:LEU:HD11	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:377:PHE:HB2	1:B:379:MET:HE1	2.01	0.42
1:D:184:PHE:O	1:D:208:ALA:HA	2.19	0.42
1:C:263:MET:O	1:C:264:LEU:HB3	2.19	0.42
1:A:459:GLN:NE2	1:A:459:GLN:H	2.16	0.42
1:B:315:TYR:HB2	1:B:328:ARG:HE	1.83	0.42
1:C:280:LEU:CD1	1:C:284:GLY:HA3	2.48	0.42
1:B:471:LYS:O	1:B:473:PRO:HD3	2.19	0.42
1:B:437:GLY:HA3	1:B:454:VAL:O	2.20	0.42
1:D:262:TYR:CZ	3:D:1005:849:H3	2.54	0.42
1:C:259:PRO:HD2	1:C:260:MET:SD	2.60	0.42
1:B:433:ASP:HB3	1:B:459:GLN:NE2	2.35	0.42
1:C:265:PHE:HB3	1:C:449:TYR:CE1	2.54	0.42
1:D:294:HIS:O	1:D:390:PHE:HA	2.20	0.42
1:D:423:PHE:HE1	1:D:453:ILE:HG21	1.85	0.42
1:D:466:ASN:HD22	1:D:466:ASN:N	2.18	0.42
1:D:378:HIS:ND1	1:D:434:THR:HG22	2.34	0.42
1:A:415:HIS:NE2	3:A:1005:849:N14	2.68	0.41
1:D:453:ILE:O	1:D:465:SER:HA	2.20	0.41
1:A:167:TYR:CD2	1:B:430:LYS:HE3	2.55	0.41
1:D:437:GLY:HA3	1:D:454:VAL:O	2.19	0.41
1:B:208:ALA:HB3	1:B:211:ILE:HD11	2.02	0.41
1:C:294:HIS:O	1:C:390:PHE:HA	2.19	0.41
1:C:446:ARG:HG3	1:C:471:LYS:NZ	2.34	0.41
1:B:255:ILE:HD12	1:B:286:MET:HE1	2.01	0.41
1:C:154:TYR:OH	3:C:1005:849:H181	2.21	0.41
1:A:295:LEU:HG	1:A:390:PHE:CE2	2.56	0.41
1:D:259:PRO:HA	3:D:1005:849:H201	2.03	0.41
1:D:265:PHE:CE2	1:D:293:VAL:HG21	2.56	0.41
1:B:136:SER:HB3	1:B:245:GLU:HG2	2.02	0.41
1:C:175:ARG:HG3	4:C:2008:HOH:O	2.21	0.41
2:A:1001:SFG:HB1	2:A:1001:SFG:H4'	2.02	0.41
1:C:190:LEU:HD13	1:C:248:LEU:HD11	2.03	0.41
1:A:149:GLN:HG2	3:A:1005:849:H24	2.03	0.41
1:D:136:SER:N	1:D:139:SER:HG	2.18	0.41
1:A:218:THR:O	1:A:221:GLN:HG2	2.20	0.41
1:C:377:PHE:HB2	1:C:379:MET:CE	2.51	0.41
1:C:350:MET:SD	1:C:385:VAL:HG22	2.61	0.41
1:A:161:GLN:NE2	1:C:335:TYR:OH	2.54	0.41
1:D:218:THR:O	1:D:221:GLN:HG2	2.21	0.40
2:D:1001:SFG:HNE2	3:D:1005:849:C18	2.33	0.40
1:B:140:GLU:O	1:B:445:LYS:HE3	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:460:THR:HG23	1:A:462:SER:N	2.13	0.40
1:C:255:ILE:HD12	1:C:286:MET:HE1	2.02	0.40
1:C:145:SER:O	1:C:149:GLN:OE1	2.39	0.40
1:C:298:PHE:HB2	1:C:348:ILE:HD12	2.03	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	342/348 (98%)	329 (96%)	13 (4%)	0	100	100
1	B	342/348 (98%)	328 (96%)	14 (4%)	0	100	100
1	C	342/348 (98%)	327 (96%)	15 (4%)	0	100	100
1	D	342/348 (98%)	326 (95%)	16 (5%)	0	100	100
All	All	1368/1392 (98%)	1310 (96%)	58 (4%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains [i](#)

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	299/303 (99%)	291 (97%)	8 (3%)	52	56
1	B	299/303 (99%)	292 (98%)	7 (2%)	58	62

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	299/303 (99%)	290 (97%)	9 (3%)	48	51
1	D	299/303 (99%)	291 (97%)	8 (3%)	52	56
All	All	1196/1212 (99%)	1164 (97%)	32 (3%)	52	56

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	140	GLU
1	A	210	LYS
1	A	248	LEU
1	A	258	GLU
1	A	286	MET
1	A	350	MET
1	A	376	LYS
1	A	467	LEU
1	B	210	LYS
1	B	258	GLU
1	B	286	MET
1	B	310	LYS
1	B	328	ARG
1	B	430	LYS
1	B	459	GLN
1	C	143	GLU
1	C	210	LYS
1	C	258	GLU
1	C	286	MET
1	C	301	GLU
1	C	328	ARG
1	C	346	ILE
1	C	459	GLN
1	C	471	LYS
1	D	210	LYS
1	D	248	LEU
1	D	258	GLU
1	D	286	MET
1	D	287	PHE
1	D	302	GLN
1	D	459	GLN
1	D	466	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:



Mol	Chain	Res	Type
1	A	159	GLN
1	A	161	GLN
1	A	174	GLN
1	A	221	GLN
1	A	456	GLN
1	A	459	GLN
1	B	159	GLN
1	B	161	GLN
1	B	174	GLN
1	B	221	GLN
1	B	307	GLN
1	B	456	GLN
1	B	459	GLN
1	C	159	GLN
1	C	161	GLN
1	C	174	GLN
1	C	221	GLN
1	C	307	GLN
1	C	316	GLN
1	C	459	GLN
1	D	159	GLN
1	D	161	GLN
1	D	174	GLN
1	D	221	GLN
1	D	316	GLN
1	D	456	GLN
1	D	459	GLN

### 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

8 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	SFG	A	1001	-	21,29,29	1.36	1 (4%)	19,42,42	1.76	2 (10%)
3	849	A	1005	-	29,31,31	2.33	11 (37%)	36,43,43	1.93	5 (13%)
2	SFG	B	1001	-	21,29,29	1.42	3 (14%)	19,42,42	1.77	2 (10%)
3	849	B	1005	-	29,31,31	2.41	13 (44%)	36,43,43	1.93	7 (19%)
2	SFG	C	1001	-	21,29,29	1.45	1 (4%)	19,42,42	1.72	3 (15%)
3	849	C	1005	-	29,31,31	2.46	11 (37%)	36,43,43	1.98	7 (19%)
2	SFG	D	1001	-	21,29,29	1.46	3 (14%)	19,42,42	1.77	3 (15%)
3	849	D	1005	-	29,31,31	2.28	13 (44%)	36,43,43	1.96	7 (19%)

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	SFG	A	1001	-	-	0/9/33/33	0/3/3/3
3	849	A	1005	-	-	0/12/24/24	0/4/4/4
2	SFG	B	1001	-	-	0/9/33/33	0/3/3/3
3	849	B	1005	-	-	0/12/24/24	0/4/4/4
2	SFG	C	1001	-	-	0/9/33/33	0/3/3/3
3	849	C	1005	-	-	0/12/24/24	0/4/4/4
2	SFG	D	1001	-	-	0/9/33/33	0/3/3/3
3	849	D	1005	-	-	0/12/24/24	0/4/4/4

All (56) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1005	849	C10-C8	-5.56	1.43	1.49
3	D	1005	849	C10-C8	-5.41	1.43	1.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1005	849	C10-C8	-4.88	1.44	1.49
3	A	1005	849	C10-C8	-3.98	1.45	1.49
3	B	1005	849	C9-C5	-3.52	1.37	1.41
3	C	1005	849	C9-C5	-2.66	1.38	1.41
3	B	1005	849	C16-C11	2.01	1.58	1.53
3	D	1005	849	C16-C11	2.03	1.59	1.53
2	D	1001	SFG	C3'-C4'	2.06	1.58	1.53
3	D	1005	849	C2-C11	2.06	1.56	1.52
2	B	1001	SFG	C3'-C4'	2.13	1.58	1.53
3	D	1005	849	C22-C21	2.15	1.44	1.39
3	D	1005	849	C12-C11	2.16	1.59	1.53
3	B	1005	849	C12-C11	2.18	1.59	1.53
3	D	1005	849	C15-N14	2.20	1.52	1.46
3	C	1005	849	C13-N14	2.21	1.53	1.46
3	A	1005	849	C4-C3	2.31	1.41	1.36
3	B	1005	849	C17-N14	2.32	1.52	1.47
3	C	1005	849	C4-C3	2.35	1.41	1.36
3	B	1005	849	C13-N14	2.36	1.53	1.46
3	A	1005	849	C13-N14	2.50	1.53	1.46
3	C	1005	849	C22-C21	2.59	1.45	1.39
3	A	1005	849	C22-C21	2.59	1.45	1.39
3	B	1005	849	C22-C21	2.61	1.45	1.39
3	D	1005	849	C3-C2	2.64	1.43	1.39
3	C	1005	849	C17-N14	2.66	1.53	1.47
3	D	1005	849	C25-C10	2.69	1.44	1.39
3	B	1005	849	C3-C2	2.72	1.43	1.39
3	D	1005	849	C13-N14	2.74	1.54	1.46
3	C	1005	849	C25-C10	2.75	1.44	1.39
2	D	1001	SFG	CD-NE	2.87	1.55	1.47
3	D	1005	849	C17-N14	2.91	1.54	1.47
3	A	1005	849	C3-C2	2.91	1.43	1.39
3	A	1005	849	C2-C11	2.93	1.57	1.52
3	B	1005	849	C25-C10	2.96	1.45	1.39
2	B	1001	SFG	C2-N3	2.97	1.37	1.32
2	B	1001	SFG	C5'-C4'	3.01	1.57	1.52
3	A	1005	849	C17-N14	3.13	1.54	1.47
3	A	1005	849	C10-C21	3.24	1.45	1.40
3	A	1005	849	C25-C10	3.27	1.45	1.39
3	C	1005	849	C2-C11	3.32	1.58	1.52
3	B	1005	849	O26-C21	3.44	1.42	1.37
3	A	1005	849	O26-C21	3.63	1.43	1.37
3	B	1005	849	C10-C21	3.64	1.46	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	1005	849	O26-C21	3.64	1.43	1.37
3	D	1005	849	C10-C21	3.65	1.46	1.40
3	B	1005	849	C2-C11	3.75	1.59	1.52
2	A	1001	SFG	C2-N3	4.13	1.39	1.32
2	D	1001	SFG	C2-N3	4.23	1.39	1.32
3	C	1005	849	O26-C21	4.26	1.44	1.37
3	C	1005	849	C10-C21	4.36	1.47	1.40
3	C	1005	849	C1-C2	4.68	1.44	1.37
3	D	1005	849	C1-C2	4.69	1.44	1.37
2	C	1001	SFG	C2-N3	4.82	1.40	1.32
3	B	1005	849	C1-C2	5.19	1.45	1.37
3	A	1005	849	C1-C2	5.44	1.45	1.37

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1001	SFG	N3-C2-N1	-6.29	124.08	128.89
2	A	1001	SFG	N3-C2-N1	-6.09	124.23	128.89
2	D	1001	SFG	N3-C2-N1	-6.06	124.25	128.89
2	C	1001	SFG	N3-C2-N1	-6.02	124.28	128.89
3	B	1005	849	C2-C1-C5	-5.35	118.23	121.93
3	D	1005	849	C2-C1-C5	-5.21	118.33	121.93
3	C	1005	849	C2-C1-C5	-5.16	118.36	121.93
3	A	1005	849	C2-C1-C5	-4.60	118.75	121.93
3	B	1005	849	C3-C4-C6	-2.43	118.24	120.88
3	C	1005	849	C3-C4-C6	-2.22	118.47	120.88
3	D	1005	849	C3-C4-C6	-2.10	118.60	120.88
3	A	1005	849	C3-C4-C6	-2.05	118.65	120.88
3	D	1005	849	C12-C11-C2	2.02	117.19	112.89
3	D	1005	849	C15-N14-C13	2.08	113.39	108.90
2	C	1001	SFG	O2'-C2'-C3'	2.08	118.59	111.83
2	A	1001	SFG	O2'-C2'-C3'	2.09	118.64	111.83
3	B	1005	849	C12-C11-C2	2.21	117.59	112.89
3	C	1005	849	C12-C11-C2	2.23	117.64	112.89
2	D	1001	SFG	C1'-N9-C4	2.25	130.33	126.94
2	B	1001	SFG	C1'-N9-C4	2.25	130.33	126.94
2	C	1001	SFG	C1'-N9-C4	2.35	130.49	126.94
2	D	1001	SFG	O2'-C2'-C3'	2.43	119.73	111.83
3	C	1005	849	C15-N14-C13	2.44	114.19	108.90
3	B	1005	849	C15-N14-C13	2.44	114.19	108.90
3	B	1005	849	C15-C16-C11	3.51	115.45	110.97
3	C	1005	849	C15-C16-C11	3.58	115.55	110.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1005	849	C15-C16-C11	3.67	115.66	110.97
3	D	1005	849	C15-C16-C11	3.77	115.79	110.97
3	C	1005	849	C13-C12-C11	4.62	116.88	110.97
3	D	1005	849	C13-C12-C11	4.74	117.03	110.97
3	A	1005	849	C13-C12-C11	4.77	117.06	110.97
3	B	1005	849	C13-C12-C11	4.77	117.07	110.97
3	B	1005	849	C27-O26-C21	6.09	126.78	117.54
3	D	1005	849	C27-O26-C21	6.36	127.19	117.54
3	C	1005	849	C27-O26-C21	6.73	127.74	117.54
3	A	1005	849	C27-O26-C21	6.98	128.13	117.54

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

8 monomers are involved in 22 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1001	SFG	2	0
3	A	1005	849	7	0
2	B	1001	SFG	1	0
3	B	1005	849	4	0
2	C	1001	SFG	1	0
3	C	1005	849	5	0
2	D	1001	SFG	2	0
3	D	1005	849	5	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	344/348 (98%)	0.17	14 (4%) 41 50	21, 36, 53, 63	0
1	B	344/348 (98%)	0.09	6 (1%) 73 78	24, 37, 49, 62	0
1	C	344/348 (98%)	0.02	8 (2%) 64 70	24, 36, 48, 63	0
1	D	344/348 (98%)	0.06	6 (1%) 73 78	21, 33, 47, 55	0
All	All	1376/1392 (98%)	0.09	34 (2%) 61 67	21, 36, 49, 63	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	478	THR	5.0
1	C	479	GLY	4.5
1	A	478	THR	4.2
1	C	144	GLU	3.8
1	C	478	THR	3.8
1	A	477	TYR	3.7
1	D	478	THR	3.6
1	B	479	GLY	3.6
1	C	477	TYR	3.3
1	A	148	VAL	3.0
1	B	477	TYR	3.0
1	A	400	ILE	2.9
1	D	234	ASP	2.9
1	D	477	TYR	2.8
1	D	256	ILE	2.8
1	D	192	VAL	2.7
1	A	256	ILE	2.6
1	A	479	GLY	2.6
1	A	234	ASP	2.6
1	C	315	TYR	2.5
1	A	257	SER	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	221	GLN	2.4
1	C	248	LEU	2.4
1	A	144	GLU	2.4
1	B	144	GLU	2.4
1	C	378	HIS	2.3
1	C	347	ARG	2.2
1	B	261	GLY	2.2
1	A	270	LEU	2.1
1	A	181	HIS	2.0
1	D	283	SER	2.0
1	A	140	GLU	2.0
1	B	170	THR	2.0
1	A	260	MET	2.0

## 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. 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( $\text{\AA}^2$ )	Q<0.9
3	849	A	1005	28/28	0.88	0.18	0.76	32,40,45,46	0
3	849	C	1005	28/28	0.87	0.17	0.73	37,45,51,52	0
3	849	B	1005	28/28	0.86	0.19	0.71	40,44,51,52	0
2	SFG	A	1001	27/27	0.93	0.15	0.60	29,33,39,40	0
2	SFG	C	1001	27/27	0.92	0.14	0.40	27,30,38,40	0
2	SFG	D	1001	27/27	0.96	0.15	0.33	23,26,35,37	0
2	SFG	B	1001	27/27	0.94	0.14	0.25	27,31,39,41	0
3	849	D	1005	28/28	0.93	0.14	-0.24	30,34,39,40	0

## 6.5 Other polymers

There are no such residues in this entry.