



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:34 PM GMT

PDB ID : 1XU9
Title : Crystal Structure of the Interface Closed Conformation of 11b-hydroxysteroid dehydrogenase isozyme 1
Authors : Hosfield, D.J.; Wu, Y.; Skene, R.J.; Hilger, M.; Jennings, A.; Snell, G.P.; Aertgeerts, K.
Deposited on : 2004-10-25
Resolution : 1.55 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
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

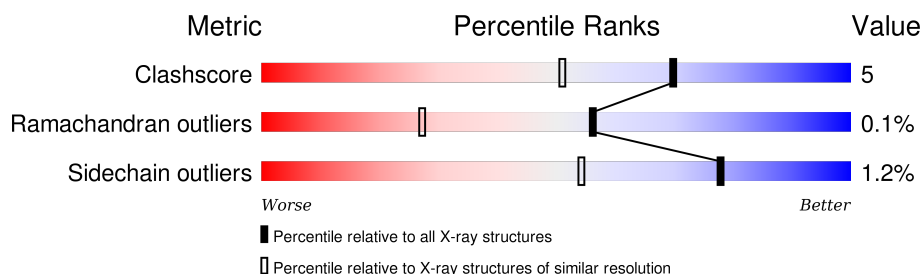
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.55 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))
Clashscore	102246	1014 (1.56-1.56)
Ramachandran outliers	100387	1704 (1.58-1.54)
Sidechain outliers	100360	1702 (1.58-1.54)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	286	<div> <div>86%</div> <div>8% 6%</div> </div>
1	B	286	<div> <div>86%</div> <div>7% • 6%</div> </div>
1	C	286	<div> <div>85%</div> <div>5% 10%</div> </div>
1	D	286	<div> <div>85%</div> <div>6% 10%</div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 9889 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 Corticosteroid 11-beta-dehydrogenase, isozyme 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	269	Total	C	N	O	S	0	11	0
			2109	1346	355	392	16			
1	B	270	Total	C	N	O	S	0	12	0
			2121	1355	357	392	17			
1	C	258	Total	C	N	O	S	0	10	0
			2012	1282	341	375	14			
1	D	258	Total	C	N	O	S	0	12	0
			2016	1283	339	377	17			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	7	MET	-	INitiating methionine	UNP P28845
A	8	LYS	-	cloning artifact	UNP P28845
A	9	HIS	-	cloning artifact	UNP P28845
A	10	GLN	-	cloning artifact	UNP P28845
A	11	HIS	-	cloning artifact	UNP P28845
A	12	GLN	-	cloning artifact	UNP P28845
A	13	HIS	-	cloning artifact	UNP P28845
A	14	GLN	-	cloning artifact	UNP P28845
A	15	HIS	-	cloning artifact	UNP P28845
A	16	GLN	-	cloning artifact	UNP P28845
A	17	HIS	-	cloning artifact	UNP P28845
A	18	GLN	-	cloning artifact	UNP P28845
A	19	HIS	-	cloning artifact	UNP P28845
A	20	GLN	-	cloning artifact	UNP P28845
A	21	GLN	-	cloning artifact	UNP P28845
A	22	PRO	-	cloning artifact	UNP P28845
A	23	LEU	-	cloning artifact	UNP P28845
A	272	SER	CYS	engineered	UNP P28845
B	7	MET	-	INitiating methionine	UNP P28845
B	8	LYS	-	cloning artifact	UNP P28845
B	9	HIS	-	cloning artifact	UNP P28845

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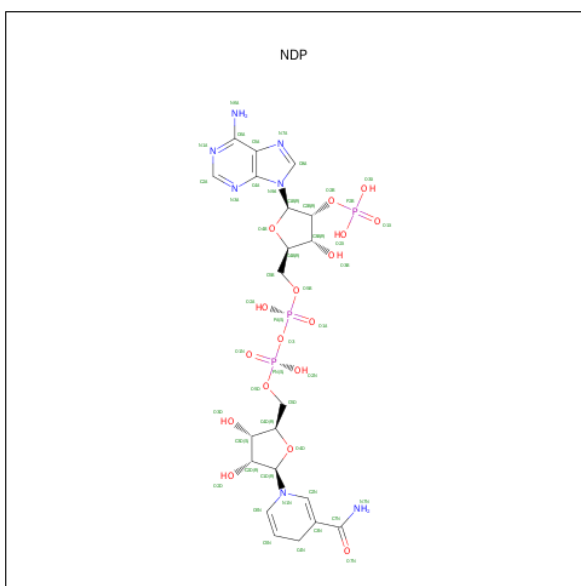
Chain	Residue	Modelled	Actual	Comment	Reference
B	10	GLN	-	cloning artifact	UNP P28845
B	11	HIS	-	cloning artifact	UNP P28845
B	12	GLN	-	cloning artifact	UNP P28845
B	13	HIS	-	cloning artifact	UNP P28845
B	14	GLN	-	cloning artifact	UNP P28845
B	15	HIS	-	cloning artifact	UNP P28845
B	16	GLN	-	cloning artifact	UNP P28845
B	17	HIS	-	cloning artifact	UNP P28845
B	18	GLN	-	cloning artifact	UNP P28845
B	19	HIS	-	cloning artifact	UNP P28845
B	20	GLN	-	cloning artifact	UNP P28845
B	21	GLN	-	cloning artifact	UNP P28845
B	22	PRO	-	cloning artifact	UNP P28845
B	23	LEU	-	cloning artifact	UNP P28845
B	272	SER	CYS	engineered	UNP P28845
C	7	MET	-	INitiating methionine	UNP P28845
C	8	LYS	-	cloning artifact	UNP P28845
C	9	HIS	-	cloning artifact	UNP P28845
C	10	GLN	-	cloning artifact	UNP P28845
C	11	HIS	-	cloning artifact	UNP P28845
C	12	GLN	-	cloning artifact	UNP P28845
C	13	HIS	-	cloning artifact	UNP P28845
C	14	GLN	-	cloning artifact	UNP P28845
C	15	HIS	-	cloning artifact	UNP P28845
C	16	GLN	-	cloning artifact	UNP P28845
C	17	HIS	-	cloning artifact	UNP P28845
C	18	GLN	-	cloning artifact	UNP P28845
C	19	HIS	-	cloning artifact	UNP P28845
C	20	GLN	-	cloning artifact	UNP P28845
C	21	GLN	-	cloning artifact	UNP P28845
C	22	PRO	-	cloning artifact	UNP P28845
C	23	LEU	-	cloning artifact	UNP P28845
C	272	SER	CYS	engineered	UNP P28845
D	7	MET	-	INitiating methionine	UNP P28845
D	8	LYS	-	cloning artifact	UNP P28845
D	9	HIS	-	cloning artifact	UNP P28845
D	10	GLN	-	cloning artifact	UNP P28845
D	11	HIS	-	cloning artifact	UNP P28845
D	12	GLN	-	cloning artifact	UNP P28845
D	13	HIS	-	cloning artifact	UNP P28845
D	14	GLN	-	cloning artifact	UNP P28845
D	15	HIS	-	cloning artifact	UNP P28845

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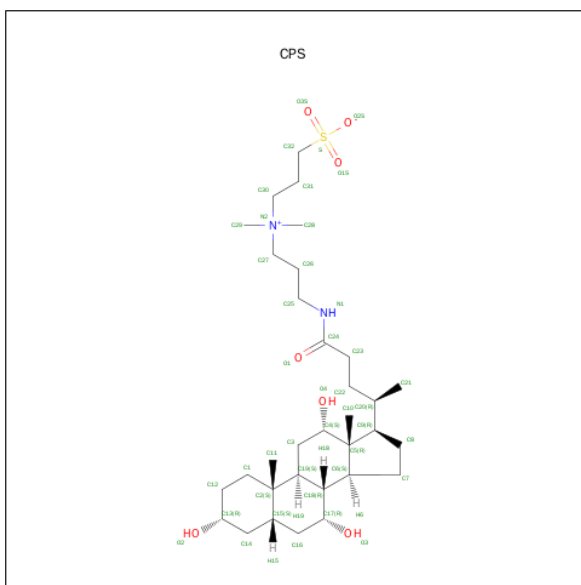
Chain	Residue	Modelled	Actual	Comment	Reference
D	16	GLN	-	cloning artifact	UNP P28845
D	17	HIS	-	cloning artifact	UNP P28845
D	18	GLN	-	cloning artifact	UNP P28845
D	19	HIS	-	cloning artifact	UNP P28845
D	20	GLN	-	cloning artifact	UNP P28845
D	21	GLN	-	cloning artifact	UNP P28845
D	22	PRO	-	cloning artifact	UNP P28845
D	23	LEU	-	cloning artifact	UNP P28845
D	272	SER	CYS	engineered	UNP P28845

- Molecule 2 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NDP) (formula: $C_{21}H_{30}N_7O_{17}P_3$).



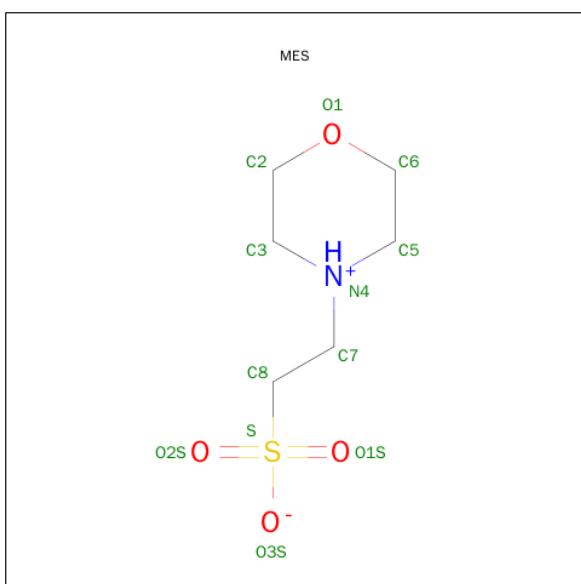
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	B	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	C	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	D	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 3 is 3-[(3-CHOLAMIDOPROPYL)DIMETHYLAMMONIO]-1-PROPANESULFONATE (three-letter code: CPS) (formula: $C_{32}H_{58}N_2O_7S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 42	C 32	N 2	O 7	S 1	0	0
3	B	1	Total 42	C 32	N 2	O 7	S 1	0	0
3	C	1	Total 42	C 32	N 2	O 7	S 1	0	0
3	D	1	Total 42	C 32	N 2	O 7	S 1	0	0

- Molecule 4 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: $\text{C}_6\text{H}_{13}\text{NO}_4\text{S}$).



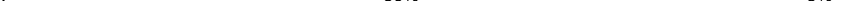

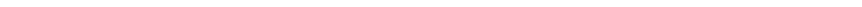
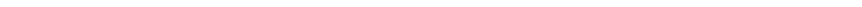











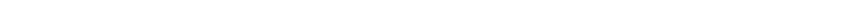


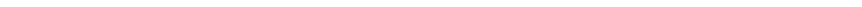

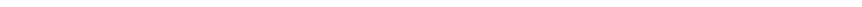
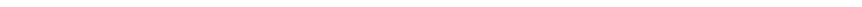

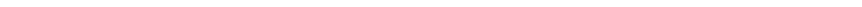


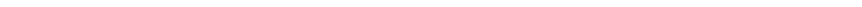
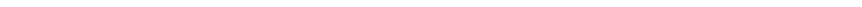

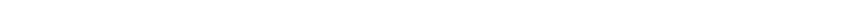
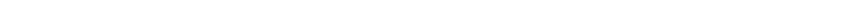

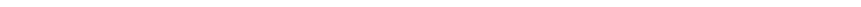
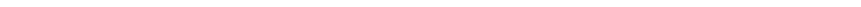




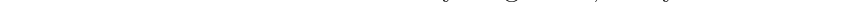



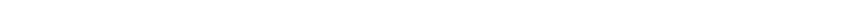





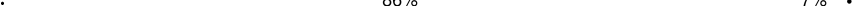

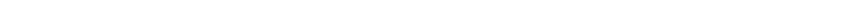
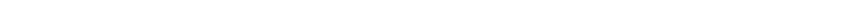

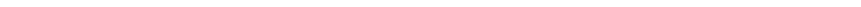












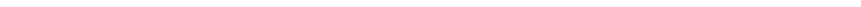





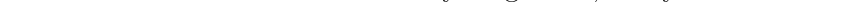

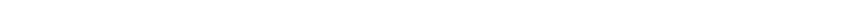

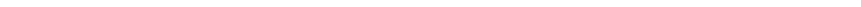
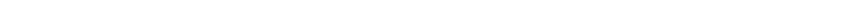







Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 5 is water.


Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	350	Total	O	0	0
			350	350		
5	B	293	Total	O	0	0
			293	293		
5	C	339	Total	O	0	0
			339	339		
5	D	277	Total	O	0	0
			277	277		

Note EDS was not executed.

- Chain A: 
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- Chain B:  86% 7% 6%
- | Residue | State |
|---------|--------|
| MET | Green |
| LVS | Green |
| HIS | Green |
| GLN | Green |
| HIS | Green |
| GLN | Green |
| HIS | Green |
| GLN | Green |
| HIS | Green |
| GLN | Green |
| HIS | Green |
| GLN | Green |
| HIS | Green |
| Q20 | Green |
| T40 | Yellow |
| S67 | Yellow |
| H87 | Yellow |
| D114 | Yellow |
| H120 | Yellow |
| T124 | Orange |
| I133 | Orange |
| M140 | Orange |
| S146 | Orange |
| S169 | Orange |
| M179 | Orange |
| F194 | Orange |
| S195 | Green |
| S196 | Green |
| I197 | Green |
| D219 | Green |
| Q253 | Green |
| D259 | Green |
| R269 | Green |
| N270 | Green |
| P271 | Green |
| S283 | Orange |
| R288 | Orange |
| P289 | Red |
| ILE | Red |
| ASU | Grey |
| LVS | Grey |

- Chain C:**
-
- | Amino Acid | Percentage |
|------------|------------|
| MET | 85% |
| LNS | 85% |
| HIS | 85% |
| GLN | 85% |
| HIS | 85% |
| GLN | 85% |
| HIS | 85% |
| GLN | 85% |
| GLN | 85% |
| GLN | 85% |
| GLN | 85% |
| HIS | 85% |
| Q20 | 5% |
| K56 | 5% |
| E69 | 5% |
| H87 | 5% |
| H130 | 5% |
| D131 | 5% |
| H135 | 5% |
| I136 | 5% |
| R137 | 5% |
| S169 | 5% |
| Y183 | 5% |
| F193 | 5% |
| I197 | 5% |
| S202 | 5% |
| I230 | 5% |
| VAL | 5% |
| HIS | 5% |
| MET | 5% |
| GLN | 5% |
| GLN | 5% |
| A235 | 5% |
| K274 | 5% |
| S281 | 5% |
| THR | 5% |
| SER | 5% |
| TVA | 5% |
| ASN | 5% |
| ASP | 5% |
| ARG | 5% |
| PHE | 5% |
| ILE | 5% |
| ASN | 5% |

- Chain D: 

4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	57.43 Å 152.67 Å 73.92 Å 90.00° 93.77° 90.00°	Depositor
Resolution (Å)	20.00 – 1.55	Depositor
% Data completeness (in resolution range)	92.1 (20.00-1.55)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.04	Depositor
Refinement program	REFMAC 5.1.19	Depositor
R, R_{free}	0.157 , 0.181	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	9889	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NDP, CPS, MES

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	0.45	0/2192	0.67	1/2960 (0.0%)
1	B	0.78	4/2205 (0.2%)	1.14	7/2977 (0.2%)
1	C	0.43	0/2087	0.65	0/2817
1	D	0.41	0/2101	0.63	1/2833 (0.0%)
All	All	0.54	4/8585 (0.0%)	0.80	9/11587 (0.1%)

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	C	0	1
All	All	0	3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	289	PHE	CD1-CE1	19.26	1.77	1.39
1	B	289	PHE	CB-CG	-15.72	1.24	1.51
1	B	289	PHE	CG-CD2	13.16	1.58	1.38
1	B	289	PHE	C-O	8.09	1.38	1.23

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	289	PHE	CB-CG-CD1	-35.98	95.61	120.80
1	B	289	PHE	CB-CG-CD2	32.85	143.80	120.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	289	PHE	CA-CB-CG	9.35	136.34	113.90
1	A	219	ASP	CB-CG-OD2	7.26	124.83	118.30
1	D	219	ASP	CB-CG-OD2	6.33	123.99	118.30
1	B	289	PHE	CG-CD1-CE1	6.09	127.50	120.80
1	B	219	ASP	CB-CG-OD2	5.99	123.69	118.30
1	B	114	ASP	CB-CG-OD2	5.63	123.36	118.30
1	B	259	ASP	CB-CG-OD2	5.34	123.11	118.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	169	SER	Peptide
1	B	169	SER	Peptide
1	C	169	SER	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	2109	0	2145	29	0
1	B	2121	0	2161	29	0
1	C	2012	0	2054	12	0
1	D	2016	0	2056	11	0
2	A	48	0	26	0	0
2	B	48	0	26	0	0
2	C	48	0	26	0	0
2	D	48	0	26	1	0
3	A	42	0	58	0	0
3	B	42	0	58	1	0
3	C	42	0	58	3	0
3	D	42	0	58	3	0
4	A	12	0	12	0	0
5	A	350	0	0	19	0
5	B	293	0	0	7	0
5	C	339	0	0	11	0
5	D	277	0	0	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	9889	0	8764	79	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 5.

All (79) 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:289:PHE:CE1	1:B:289:PHE:CD1	1.77	1.67
1:A:126:LEU:HB2	1:B:289:PHE:CE1	1.42	1.52
1:A:126:LEU:CB	1:B:289:PHE:CE1	2.02	1.30
5:A:2739:HOH:O	1:B:289:PHE:CE1	1.98	1.16
5:A:2739:HOH:O	1:B:289:PHE:CD2	2.06	1.09
1:A:196[B]:SER:OG	5:A:1946:HOH:O	1.83	0.95
1:A:67[B]:SER:OG	5:A:1974:HOH:O	2.00	0.79
1:B:67[B]:SER:OG	5:B:1892:HOH:O	2.00	0.78
1:A:126:LEU:N	1:B:289:PHE:CE1	2.56	0.70
1:A:224:MET:O	1:A:232:HIS:HE1	1.75	0.69
1:B:253:GLN:OE1	5:B:2819:HOH:O	2.10	0.69
1:B:288:ARG:CG	1:B:289:PHE:H	2.05	0.69
1:B:196[B]:SER:OG	5:B:1712:HOH:O	1.95	0.69
1:A:40:THR:OG1	1:A:120:HIS:HD2	1.79	0.66
1:D:40:THR:OG1	1:D:120:HIS:HD2	1.79	0.66
1:A:126:LEU:O	1:B:289:PHE:CD1	2.50	0.64
1:A:124[B]:THR:CG2	5:A:2365:HOH:O	2.44	0.63
1:B:40:THR:OG1	1:B:120:HIS:HD2	1.82	0.61
1:A:124[B]:THR:HG21	5:A:1548:HOH:O	2.02	0.60
1:A:94:GLU:CD	5:A:1820:HOH:O	2.40	0.59
1:A:124[B]:THR:HG21	5:A:2365:HOH:O	2.02	0.59
1:C:202[A]:SER:OG	5:C:532:HOH:O	2.16	0.58
1:D:197:ILE:C	5:D:533:HOH:O	2.41	0.58
1:B:124:THR:O	5:B:2529:HOH:O	2.17	0.58
1:B:288:ARG:CG	1:B:289:PHE:N	2.67	0.57
1:C:56:LYS:NZ	5:C:404:HOH:O	2.37	0.56
1:A:126:LEU:HB3	1:B:289:PHE:CE1	2.26	0.56
1:A:269:ARG:NE	5:A:2252:HOH:O	2.38	0.55
1:A:224:MET:O	1:A:232:HIS:CE1	2.58	0.55
5:A:2739:HOH:O	1:B:289:PHE:CD1	2.58	0.55
1:B:288:ARG:HG2	1:B:289:PHE:N	2.21	0.55
1:B:288:ARG:HG2	1:B:289:PHE:H	1.71	0.54
1:B:283:SER:HB2	5:B:2162:HOH:O	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:229:GLY:O	1:A:232:HIS:HD2	1.90	0.54
1:A:124[B]:THR:CG2	5:A:1548:HOH:O	2.57	0.53
1:D:120:HIS:HE1	1:D:146:SER:OG	1.91	0.53
1:A:140:MET:HB3	5:A:1708:HOH:O	2.09	0.52
1:A:120:HIS:HE1	1:A:146:SER:OG	1.93	0.52
1:B:120:HIS:HE1	1:B:146:SER:OG	1.93	0.52
1:B:87:HIS:HD2	5:B:1603:HOH:O	1.92	0.51
1:C:87:HIS:HD2	5:C:306:HOH:O	1.93	0.51
1:B:269:ARG:HG3	1:B:271:PRO:HD3	1.92	0.51
3:B:6:CPS:H4	3:B:6:CPS:H21A	1.93	0.51
1:A:286:MET:SD	1:B:179[B]:MET:HE3	2.51	0.50
3:C:293:CPS:H32	5:C:376:HOH:O	2.12	0.49
3:C:293:CPS:H21A	3:C:293:CPS:H4	1.96	0.47
1:C:135:HIS:HD2	5:C:372:HOH:O	1.96	0.47
1:D:140[A]:MET:HG2	5:D:314:HOH:O	2.14	0.46
1:A:124[B]:THR:HG23	5:A:2365:HOH:O	2.10	0.46
3:D:293:CPS:H21A	3:D:293:CPS:H4	1.98	0.46
1:C:69[A]:GLU:HG3	5:C:334:HOH:O	2.16	0.45
1:B:140:MET:HA	1:B:140:MET:HE3	1.97	0.45
1:A:253:GLN:NE2	5:A:2092:HOH:O	2.50	0.45
1:D:183:TYR:CZ	3:D:293:CPS:H16	2.51	0.45
1:C:137:ARG:NH1	5:C:373:HOH:O	2.48	0.45
1:C:274:LYS:HD2	5:C:440:HOH:O	2.17	0.44
1:A:122:THR:O	1:A:124[B]:THR:HG23	2.17	0.44
3:D:293:CPS:H31	3:D:293:CPS:H29A	1.82	0.44
1:D:179[B]:MET:HE3	5:D:535:HOH:O	2.18	0.44
1:A:138:LYS:NZ	5:A:2815:HOH:O	2.41	0.44
1:A:286:MET:SD	1:B:179[B]:MET:CE	3.06	0.43
1:D:56:LYS:HE2	1:D:81:LEU:HD22	2.00	0.43
1:B:194:PHE:HA	1:B:197[A]:ILE:HG12	2.01	0.43
1:A:87:HIS:HE1	5:A:2028:HOH:O	2.01	0.43
1:C:20:GLN:HA	5:C:532:HOH:O	2.17	0.42
1:D:201:TYR:CE2	5:D:533:HOH:O	2.71	0.42
1:C:130:HIS:N	5:C:370:HOH:O	2.53	0.42
1:C:183:TYR:CZ	3:C:293:CPS:H16	2.55	0.42
1:D:20:GLN:HA	5:D:434:HOH:O	2.20	0.42
1:D:202[A]:SER:OG	5:D:434:HOH:O	2.22	0.41
1:B:288:ARG:CD	1:B:289:PHE:H	2.32	0.41
1:C:193:PHE:CZ	1:C:197[B]:ILE:HD13	2.55	0.41
1:A:253:GLN:NE2	5:A:2822:HOH:O	2.51	0.41
1:D:119:ASN:ND2	2:D:4:NDP:H4D	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:87:HIS:HD2	5:A:1625:HOH:O	2.04	0.41
1:A:149:VAL:HG22	1:B:133:ILE:HD13	2.02	0.40
1:C:274:LYS:HE2	5:C:469:HOH:O	2.21	0.40
1:B:140:MET:HB3	5:B:2368:HOH:O	2.20	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	278/286 (97%)	268 (96%)	10 (4%)	0	100	100
1	B	280/286 (98%)	269 (96%)	10 (4%)	1 (0%)	39	14
1	C	264/286 (92%)	257 (97%)	7 (3%)	0	100	100
1	D	266/286 (93%)	257 (97%)	9 (3%)	0	100	100
All	All	1088/1144 (95%)	1051 (97%)	36 (3%)	1 (0%)	56	27

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	288	ARG

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	237/243 (98%)	233 (98%)	4 (2%)	68	38
1	B	239/243 (98%)	235 (98%)	4 (2%)	68	38
1	C	225/243 (93%)	222 (99%)	3 (1%)	76	51
1	D	228/243 (94%)	226 (99%)	2 (1%)	84	66
All	All	929/972 (96%)	916 (99%)	13 (1%)	78	48

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	GLU
1	A	124[A]	THR
1	A	124[B]	THR
1	A	285	ASN
1	B	124	THR
1	B	270	ASN
1	B	283	SER
1	B	289	PHE
1	C	69[A]	GLU
1	C	69[B]	GLU
1	C	131	ASP
1	D	109	LEU
1	D	270	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	77	HIS
1	A	87	HIS
1	A	105	GLN
1	A	119	ASN
1	A	120	HIS
1	A	232	HIS
1	A	234	GLN
1	A	253	GLN
1	A	270	ASN
1	A	285	ASN
1	B	87	HIS
1	B	119	ASN
1	B	120	HIS
1	B	234	GLN

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Mol	Chain	Res	Type
1	B	270	ASN
1	C	87	HIS
1	C	119	ASN
1	C	135	HIS
1	C	270	ASN
1	D	119	ASN
1	D	120	HIS
1	D	135	HIS
1	D	270	ASN

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	NDP	A	1	-	42,52,52	1.37	3 (7%)	55,80,80	2.04	14 (25%)
4	MES	A	293	-	11,12,12	0.54	0	14,16,16	1.33	1 (7%)
3	CPS	A	5	-	44,45,45	0.75	0	67,70,70	1.06	3 (4%)
2	NDP	B	2	-	42,52,52	1.41	5 (11%)	55,80,80	1.84	13 (23%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	CPS	B	6	-	44,45,45	0.67	0	67,70,70	1.31	5 (7%)
3	CPS	C	293	-	44,45,45	0.67	0	67,70,70	1.38	4 (5%)
2	NDP	C	3	-	42,52,52	1.39	5 (11%)	55,80,80	1.98	14 (25%)
3	CPS	D	293	-	44,45,45	0.57	0	67,70,70	1.36	3 (4%)
2	NDP	D	4	-	42,52,52	1.48	3 (7%)	55,80,80	2.12	15 (27%)

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	NDP	A	1	-	-	0/30/77/77	0/5/5/5
4	MES	A	293	-	-	0/6/14/14	0/1/1/1
3	CPS	A	5	-	-	0/25/90/90	0/4/4/4
2	NDP	B	2	-	-	0/30/77/77	0/5/5/5
3	CPS	B	6	-	-	0/25/90/90	0/4/4/4
3	CPS	C	293	-	-	0/25/90/90	0/4/4/4
2	NDP	C	3	-	-	0/30/77/77	0/5/5/5
3	CPS	D	293	-	-	0/25/90/90	0/4/4/4
2	NDP	D	4	-	-	0/30/77/77	0/5/5/5

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2	NDP	C4N-C5N	-4.91	1.38	1.49
2	A	1	NDP	C4N-C5N	-4.73	1.38	1.49
2	D	4	NDP	C4N-C5N	-4.72	1.38	1.49
2	C	3	NDP	C4N-C5N	-4.49	1.39	1.49
2	C	3	NDP	C3B-C2B	-2.23	1.48	1.53
2	C	3	NDP	C2N-C3N	2.03	1.39	1.34
2	B	2	NDP	C2N-C3N	2.07	1.39	1.34
2	B	2	NDP	C2A-N3A	2.27	1.36	1.32
2	B	2	NDP	C6N-C5N	2.56	1.38	1.33
2	D	4	NDP	C6N-C5N	2.70	1.38	1.33
2	A	1	NDP	C6N-C5N	2.85	1.38	1.33
2	C	3	NDP	C6N-C5N	2.88	1.38	1.33
2	A	1	NDP	C4A-N3A	4.60	1.42	1.35
2	C	3	NDP	C4A-N3A	4.67	1.42	1.35
2	B	2	NDP	C4A-N3A	4.68	1.42	1.35
2	D	4	NDP	C4A-N3A	5.51	1.43	1.35

All (72) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	4	NDP	N3A-C2A-N1A	-9.03	121.98	128.89
2	A	1	NDP	N3A-C2A-N1A	-7.94	122.81	128.89
2	C	3	NDP	N3A-C2A-N1A	-7.00	123.54	128.89
2	B	2	NDP	N3A-C2A-N1A	-5.96	124.33	128.89
2	C	3	NDP	C4A-C5A-N7A	-4.29	105.53	109.48
2	B	2	NDP	C4A-C5A-N7A	-4.28	105.54	109.48
2	D	4	NDP	C4A-C5A-N7A	-4.02	105.78	109.48
2	A	1	NDP	C3N-C2N-N1N	-3.42	118.24	123.14
2	C	3	NDP	O2B-P2B-O1X	-3.41	98.60	107.11
2	C	3	NDP	C3N-C2N-N1N	-3.32	118.39	123.14
2	D	4	NDP	O3B-C3B-C2B	-3.20	101.93	111.16
2	A	1	NDP	O3B-C3B-C2B	-3.15	102.06	111.16
2	D	4	NDP	O2B-P2B-O1X	-3.10	99.37	107.11
3	A	5	CPS	C19-C18-C17	-3.08	108.28	111.92
2	A	1	NDP	O3D-C3D-C2D	-3.05	101.92	111.83
2	A	1	NDP	C4A-C5A-N7A	-2.96	106.75	109.48
3	B	6	CPS	C19-C18-C17	-2.94	108.45	111.92
2	A	1	NDP	C4N-C5N-C6N	-2.87	117.85	122.58
2	B	2	NDP	C3N-C2N-N1N	-2.82	119.11	123.14
3	B	6	CPS	C5-C9-C20	-2.78	116.12	119.50
2	C	3	NDP	O3B-C3B-C2B	-2.75	103.20	111.16
3	D	293	CPS	C19-C18-C17	-2.65	108.80	111.92
2	B	2	NDP	O3D-C3D-C2D	-2.58	103.44	111.83
2	B	2	NDP	O3B-C3B-C2B	-2.57	103.73	111.16
2	B	2	NDP	O3B-C3B-C4B	-2.53	103.46	111.05
2	C	3	NDP	O3D-C3D-C2D	-2.35	104.20	111.83
2	D	4	NDP	O3D-C3D-C2D	-2.33	104.24	111.83
3	B	6	CPS	C29-N2-C30	-2.31	103.56	109.43
3	C	293	CPS	C19-C18-C17	-2.29	109.22	111.92
2	A	1	NDP	O3D-C3D-C4D	-2.18	104.52	111.05
2	D	4	NDP	O3B-C3B-C4B	-2.14	104.65	111.05
2	A	1	NDP	O3B-C3B-C4B	-2.10	104.75	111.05
2	C	3	NDP	C1D-N1N-C6N	-2.10	116.11	120.81
2	C	3	NDP	C4N-C5N-C6N	-2.06	119.19	122.58
2	B	2	NDP	O2B-P2B-O1X	-2.05	101.98	107.11
2	D	4	NDP	C5B-C4B-C3B	-2.03	107.14	115.21
3	C	293	CPS	C3-C19-C18	2.01	113.59	110.73
3	C	293	CPS	C19-C18-C6	2.17	112.61	109.62
2	C	3	NDP	C2A-N1A-C6A	2.19	122.69	118.77
2	B	2	NDP	C2A-N1A-C6A	2.31	122.89	118.77
3	A	5	CPS	C19-C18-C6	2.34	112.85	109.62
2	D	4	NDP	C2B-C3B-C4B	2.36	107.43	101.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	4	NDP	O3X-P2B-O1X	2.39	118.27	110.58
2	D	4	NDP	O3X-P2B-O2X	2.39	116.48	107.38
2	D	4	NDP	C5N-C4N-C3N	2.39	119.12	112.52
2	C	3	NDP	C5N-C4N-C3N	2.40	119.13	112.52
2	B	2	NDP	O2D-C2D-C1D	2.41	118.35	109.94
2	B	2	NDP	C5N-C4N-C3N	2.52	119.45	112.52
2	A	1	NDP	C2A-N1A-C6A	2.52	123.27	118.77
2	D	4	NDP	C2A-N1A-C6A	2.59	123.39	118.77
2	A	1	NDP	O2B-C2B-C1B	2.59	120.12	110.02
2	A	1	NDP	O2D-C2D-C1D	2.60	119.04	109.94
2	C	3	NDP	O2D-C2D-C1D	2.62	119.11	109.94
2	C	3	NDP	O2B-C2B-C1B	2.72	120.64	110.02
2	B	2	NDP	O2B-C2B-C1B	2.77	120.82	110.02
2	D	4	NDP	O2D-C2D-C1D	2.81	119.77	109.94
2	A	1	NDP	C5N-C4N-C3N	2.82	120.28	112.52
3	A	5	CPS	O3S-S-C32	3.14	109.59	106.91
3	D	293	CPS	O1S-S-C32	3.16	109.60	106.91
2	D	4	NDP	O4D-C1D-N1N	3.29	115.00	108.07
2	C	3	NDP	O4D-C1D-N1N	3.41	115.27	108.07
2	B	2	NDP	O4B-C1B-N9A	3.43	115.28	108.10
2	C	3	NDP	O4B-C1B-N9A	3.54	115.50	108.10
2	A	1	NDP	O4B-C1B-N9A	3.63	115.69	108.10
2	B	2	NDP	O4D-C1D-N1N	3.70	115.88	108.07
4	A	293	MES	C5-N4-C3	3.86	117.26	108.90
2	A	1	NDP	O4D-C1D-N1N	4.06	116.64	108.07
3	B	6	CPS	O1S-S-C32	4.35	110.61	106.91
3	B	6	CPS	O3S-S-C32	4.40	110.66	106.91
2	D	4	NDP	O4B-C1B-N9A	4.46	117.44	108.10
3	D	293	CPS	O3S-S-C32	7.40	113.22	106.91
3	C	293	CPS	O1S-S-C32	8.08	113.80	106.91

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	6	CPS	1	0
3	C	293	CPS	3	0
3	D	293	CPS	3	0
2	D	4	NDP	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

6.5 Other polymers [i](#)

EDS was not executed - this section will therefore be empty.