



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:01 PM GMT

PDB ID : 1W73  
Title : BINARY STRUCTURE OF HUMAN DECR SOLVED BY SEMET SAD.  
Authors : Alphey, M.S.; Byres, E.; Hunter, W.N.  
Deposited on : 2004-08-27  
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

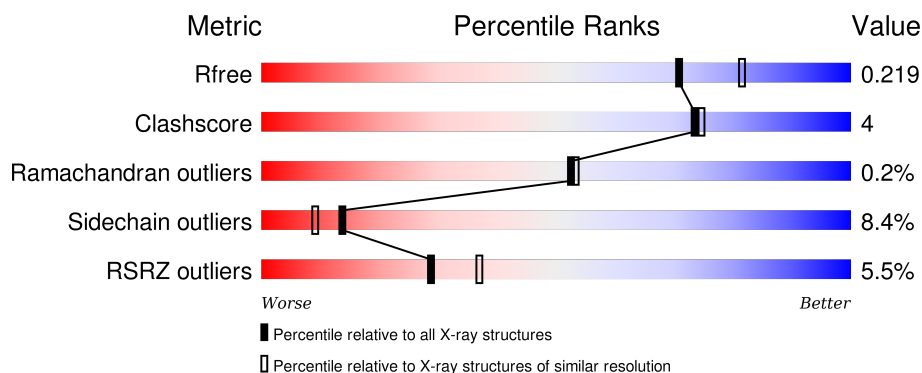
# 1 Overall quality at a glance

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	302	<div> <div>6%</div> <div>82% 12% • 5%</div> </div>
1	B	302	<div> <div>3%</div> <div>80% 12% • 7%</div> </div>
1	C	302	<div> <div>6%</div> <div>79% 13% • 6%</div> </div>
1	D	302	<div> <div>5%</div> <div>79% 11% • 7%</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	GOL	A	1329	-	-	-	X
2	GOL	D	1328	-	-	-	X
3	NAP	B	1330	-	-	-	X

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 9296 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 2,4-DIENOYL-COA REDUCTASE.

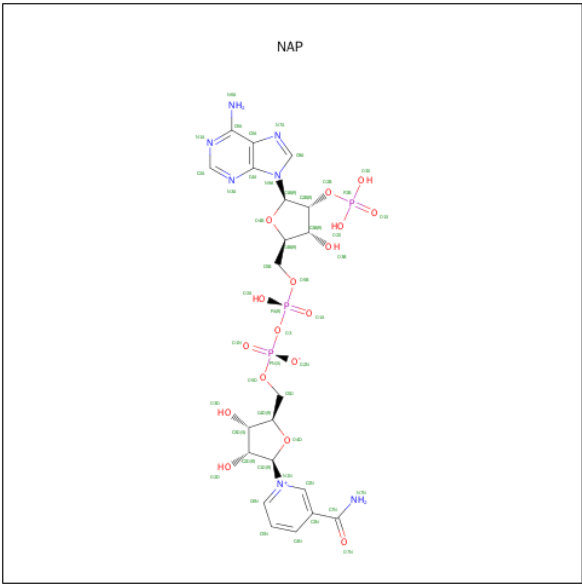
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	288	Total	C	N	O	S	Se	0	5	0
			2178	1380	370	417	4	7			
1	B	282	Total	C	N	O	S	Se	0	5	0
			2137	1354	363	408	4	8			
1	C	284	Total	C	N	O	S	Se	0	5	0
			2148	1361	363	413	4	7			
1	D	280	Total	C	N	O	S	Se	0	1	0
			2102	1332	357	401	4	8			

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: C<sub>21</sub>H<sub>28</sub>N<sub>7</sub>O<sub>17</sub>P<sub>3</sub>).



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

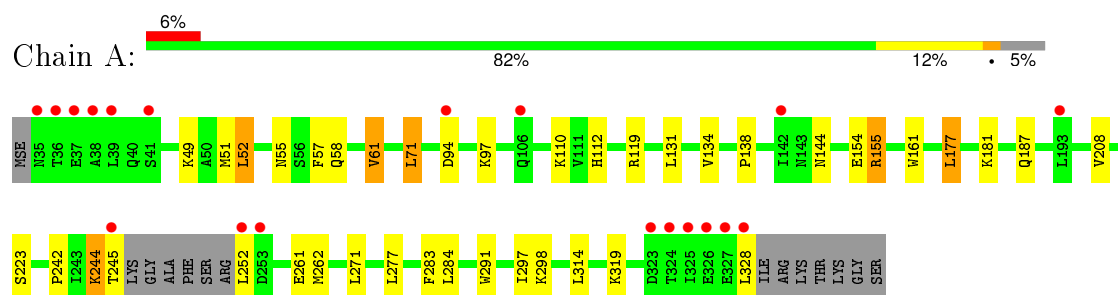
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	132	Total	O	0	0
			132	132		
4	B	137	Total	O	0	0
			137	137		
4	C	119	Total	O	0	0
			119	119		
4	D	139	Total	O	0	0
			139	139		

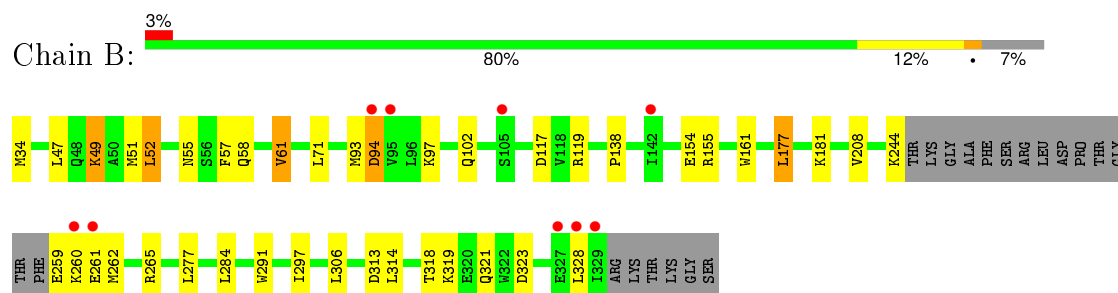
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). 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.

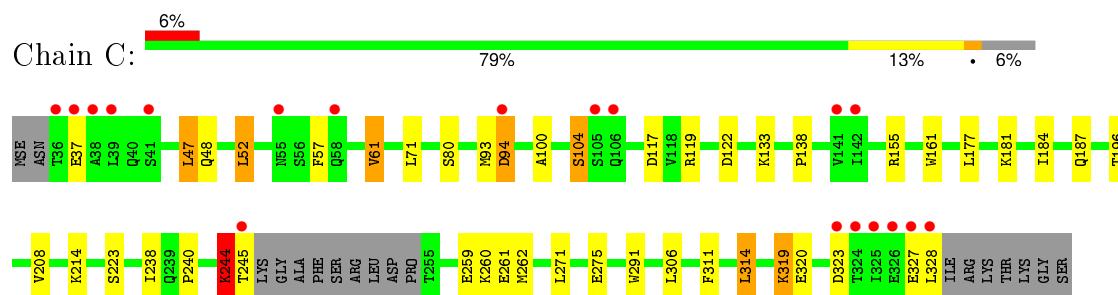
#### • Molecule 1: 2,4-DIENOYL-COA REDUCTASE



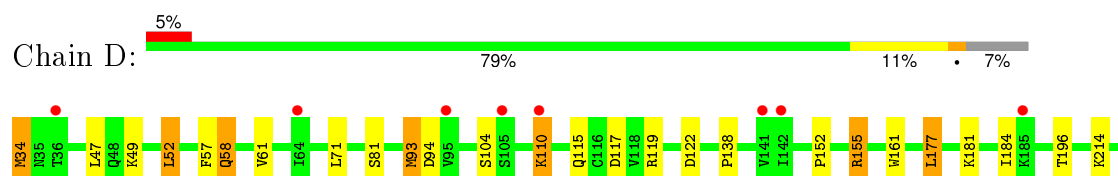
#### • Molecule 1: 2,4-DIENOYL-COA REDUCTASE

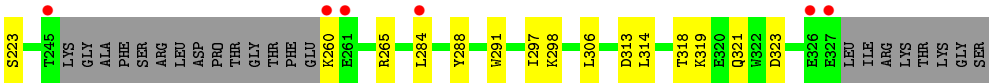


#### • Molecule 1: 2,4-DIENOYL-COA REDUCTASE



#### • Molecule 1: 2,4-DIENOYL-COA REDUCTASE





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	62.95Å 131.74Å 71.30Å 90.00° 92.49° 90.00°	Depositor
Resolution (Å)	70.71 – 2.10 29.89 – 2.10	Depositor EDS
% Data completeness (in resolution range)	96.5 (70.71-2.10) 96.5 (29.89-2.10)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	5.26 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, $R_{free}$	0.180 , 0.221 0.180 , 0.219	Depositor DCC
$R_{free}$ test set	3302 reflections (5.34%)	DCC
Wilson B-factor (Å <sup>2</sup> )	23.1	Xtriage
Anisotropy	0.765	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 50.5	EDS
Estimated twinning fraction	0.070 for h,-k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 65177 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	9296	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.46% of the height of the origin peak. No significant pseudotranslation is detected.*

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

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



## 5 Model quality [i](#)

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

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

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	0.49	0/2229	0.72	2/3004 (0.1%)
1	B	0.47	0/2188	0.72	5/2944 (0.2%)
1	C	0.47	0/2198	0.70	5/2961 (0.2%)
1	D	0.47	0/2134	0.69	5/2873 (0.2%)
All	All	0.48	0/8749	0.71	17/11782 (0.1%)

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	94[A]	ASP	CB-CG-OD2	6.71	124.34	118.30
1	C	94[B]	ASP	CB-CG-OD2	6.71	124.34	118.30
1	B	117	ASP	CB-CG-OD2	5.91	123.62	118.30
1	C	122	ASP	CB-CG-OD2	5.91	123.61	118.30
1	B	323	ASP	CB-CG-OD2	5.86	123.57	118.30
1	B	313	ASP	CB-CG-OD2	5.55	123.29	118.30
1	D	323	ASP	CB-CG-OD2	5.53	123.27	118.30
1	B	94[A]	ASP	CB-CG-OD2	5.37	123.14	118.30
1	B	94[B]	ASP	CB-CG-OD2	5.37	123.14	118.30
1	C	323	ASP	CB-CG-OD2	5.35	123.12	118.30
1	D	313	ASP	CB-CG-OD2	5.33	123.10	118.30
1	D	117	ASP	CB-CG-OD2	5.27	123.05	118.30
1	C	117	ASP	CB-CG-OD2	5.23	123.01	118.30
1	D	94	ASP	CB-CG-OD2	5.21	122.99	118.30
1	A	94[A]	ASP	CB-CG-OD2	5.05	122.85	118.30
1	A	94[B]	ASP	CB-CG-OD2	5.05	122.85	118.30
1	D	122	ASP	CB-CG-OD2	5.04	122.83	118.30

There are no chirality outliers.

There are no planarity outliers.

## 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	2178	0	2200	23	0
1	B	2137	0	2165	20	0
1	C	2148	0	2169	19	0
1	D	2102	0	2133	17	0
2	A	6	0	8	0	0
2	D	6	0	8	1	0
3	A	48	0	25	3	0
3	B	48	0	25	2	0
3	C	48	0	25	1	0
3	D	48	0	25	2	0
4	A	132	0	0	4	1
4	B	137	0	0	3	1
4	C	119	0	0	0	0
4	D	139	0	0	4	0
All	All	9296	0	8783	72	1

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 4.

All (72) 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:D:61:VAL:HG22	1:D:138:PRO:HA	1.54	0.87
4:B:2063:HOH:O	1:C:319:LYS:HD2	1.83	0.78
1:A:61:VAL:HG22	1:A:138:PRO:HA	1.65	0.77
1:A:71:LEU:HB2	3:A:1330:NAP:H51N	1.73	0.69
2:D:1328:GOL:O3	4:D:2133:HOH:O	2.11	0.68
1:A:154:GLU:HG2	1:C:184:ILE:HD13	1.79	0.64
1:B:61:VAL:HG22	1:B:138:PRO:HA	1.80	0.63
1:C:187:GLN:OE1	1:D:34:MSE:N	2.31	0.62
1:B:71:LEU:HB2	3:B:1330:NAP:H51N	1.80	0.61
1:C:177:LEU:HD22	1:C:181:LYS:HE3	1.82	0.60
1:D:318:THR:H	1:D:321:GLN:HE21	1.50	0.59
1:A:71:LEU:HG	1:A:277:LEU:CD2	2.32	0.59
1:B:47:LEU:HD22	1:B:49:LYS:HD2	1.84	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:71:LEU:HG	1:A:277:LEU:HD22	1.83	0.58
1:A:284:LEU:HD11	1:A:297:ILE:HD12	1.87	0.57
1:B:284:LEU:HD11	1:B:297:ILE:HD13	1.88	0.55
1:D:298:LYS:HE2	4:D:2116:HOH:O	2.06	0.54
1:A:208:VAL:HG22	1:C:223:SER:HB3	1.91	0.53
1:A:144:ASN:OD1	3:A:1330:NAP:H4D	2.08	0.53
1:A:52:LEU:HD23	1:A:57:PHE:HE1	1.74	0.53
3:A:1330:NAP:H5N	4:A:2073:HOH:O	2.09	0.52
1:C:161:TRP:C	1:C:161:TRP:CD1	2.83	0.52
1:D:52:LEU:HD23	1:D:57:PHE:HE1	1.75	0.51
1:A:55:ASN:ND2	1:A:58[A]:GLN:OE1	2.44	0.51
1:A:187:GLN:NE2	1:B:34:MSE:N	2.58	0.51
1:D:177:LEU:HD22	1:D:181:LYS:HE3	1.92	0.50
3:D:1330:NAP:H5N	4:D:2078:HOH:O	2.11	0.50
1:D:58:GLN:NE2	4:D:2016:HOH:O	2.43	0.50
1:D:71:LEU:HB2	3:D:1330:NAP:H51N	1.92	0.50
1:C:61:VAL:HG22	1:C:138:PRO:HA	1.93	0.50
1:B:318:THR:H	1:B:321:GLN:HE21	1.60	0.49
1:B:177:LEU:HD22	1:B:181:LYS:HE3	1.94	0.49
1:A:154:GLU:CG	1:C:184:ILE:HD13	2.44	0.48
1:A:187:GLN:HE22	1:B:34:MSE:N	2.11	0.48
1:A:283:PHE:CD1	1:B:51:MSE:SE	3.17	0.47
1:A:244:LYS:HG2	1:A:271:LEU:HD12	1.97	0.47
1:A:161:TRP:CD1	1:A:161:TRP:C	2.87	0.47
1:D:284:LEU:HD11	1:D:297:ILE:HD12	1.96	0.47
1:A:155:ARG:HD3	4:A:2052:HOH:O	2.14	0.47
1:C:47:LEU:HD13	1:D:288:TYR:CE1	2.50	0.46
1:C:259:GLU:HA	1:C:262:MSE:HE3	1.98	0.46
1:D:196:THR:HA	1:D:214:LYS:HD2	1.97	0.46
1:B:261:GLU:CD	1:B:265[B]:ARG:HH12	2.19	0.46
1:A:112[B]:HIS:CD2	1:A:134:VAL:HG21	2.51	0.45
1:B:318:THR:HG23	1:B:321:GLN:NE2	2.31	0.45
1:C:71:LEU:HB2	3:C:1330:NAP:H51N	1.97	0.45
3:B:1330:NAP:H5N	4:B:2082:HOH:O	2.17	0.45
1:C:100:ALA:O	1:C:104:SER:HB2	2.18	0.44
1:B:154:GLU:HG2	1:D:184:ILE:HD13	1.99	0.44
1:D:93:MSE:HG2	1:D:115:GLN:HB2	1.99	0.44
1:C:196:THR:HA	1:C:214:LYS:HD2	2.00	0.44
1:A:223:SER:HB3	1:C:208:VAL:HG22	2.01	0.43
1:A:51:MSE:HE3	4:A:2010:HOH:O	2.18	0.43
1:B:55:ASN:HB2	4:B:2018:HOH:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:71:LEU:HG	1:B:277:LEU:HD22	2.00	0.43
1:D:152:PRO:HB2	1:D:155:ARG:HG3	2.01	0.42
1:B:208:VAL:HG22	1:D:223:SER:HB3	2.02	0.42
1:C:244:LYS:HG2	1:C:271:LEU:HD12	2.01	0.42
1:C:48:GLN:HG2	1:C:275:GLU:HB2	2.01	0.42
1:B:318:THR:HG23	1:B:321:GLN:HE21	1.84	0.42
1:C:238:ILE:O	1:C:240:PRO:HD3	2.21	0.41
1:B:261:GLU:C	1:B:262:MSE:CA	2.89	0.41
1:D:161:TRP:CD1	1:D:161:TRP:C	2.93	0.41
1:B:52:LEU:HD23	1:B:57:PHE:HE1	1.84	0.41
1:A:242:PRO:HB2	1:A:271:LEU:HD22	2.02	0.41
1:C:311:PHE:O	1:C:314:LEU:HB2	2.20	0.41
1:D:110:LYS:HD2	1:D:110:LYS:HA	1.79	0.41
1:A:177:LEU:HD22	1:A:181:LYS:HE3	2.03	0.41
1:B:161:TRP:C	1:B:161:TRP:CD1	2.94	0.41
1:A:298:LYS:HE2	4:A:2114:HOH:O	2.20	0.40
1:C:52:LEU:HD23	1:C:57:PHE:HE1	1.86	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:2032:HOH:O	4:B:2044:HOH:O[1_556]	2.11	0.09

## 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	289/302 (96%)	281 (97%)	7 (2%)	1 (0%)	46	45
1	B	281/302 (93%)	273 (97%)	8 (3%)	0	100	100
1	C	285/302 (94%)	276 (97%)	8 (3%)	1 (0%)	39	37

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	277/302 (92%)	270 (98%)	7 (2%)	0	100	100
All	All	1132/1208 (94%)	1100 (97%)	30 (3%)	2 (0%)	52	53

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	244	LYS
1	A	244	LYS

### 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	236/235 (100%)	218 (92%)	18 (8%)	16	12
1	B	231/235 (98%)	212 (92%)	19 (8%)	14	10
1	C	233/235 (99%)	210 (90%)	23 (10%)	10	6
1	D	226/235 (96%)	208 (92%)	18 (8%)	15	11
All	All	926/940 (98%)	848 (92%)	78 (8%)	14	9

All (78) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	LYS
1	A	52	LEU
1	A	61	VAL
1	A	71	LEU
1	A	97	LYS
1	A	110	LYS
1	A	119	ARG
1	A	131	LEU
1	A	155	ARG
1	A	177	LEU
1	A	245	THR
1	A	252	LEU

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Mol	Chain	Res	Type
1	A	261	GLU
1	A	262	MSE
1	A	291	TRP
1	A	314	LEU
1	A	319	LYS
1	A	328	LEU
1	B	49	LYS
1	B	52	LEU
1	B	61	VAL
1	B	93	MSE
1	B	94[A]	ASP
1	B	94[B]	ASP
1	B	97	LYS
1	B	102	GLN
1	B	119	ARG
1	B	155	ARG
1	B	177	LEU
1	B	244	LYS
1	B	259	GLU
1	B	260	LYS
1	B	291	TRP
1	B	306	LEU
1	B	314	LEU
1	B	319	LYS
1	B	328	LEU
1	C	37	GLU
1	C	47	LEU
1	C	52	LEU
1	C	61	VAL
1	C	80	SER
1	C	93	MSE
1	C	94[A]	ASP
1	C	94[B]	ASP
1	C	104	SER
1	C	119	ARG
1	C	133	LYS
1	C	155	ARG
1	C	244	LYS
1	C	245	THR
1	C	260	LYS
1	C	261	GLU
1	C	291	TRP

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Mol	Chain	Res	Type
1	C	306	LEU
1	C	314	LEU
1	C	319	LYS
1	C	320	GLU
1	C	327	GLU
1	C	328	LEU
1	D	34	MSE
1	D	47	LEU
1	D	49	LYS
1	D	52	LEU
1	D	58	GLN
1	D	81	SER
1	D	93	MSE
1	D	104	SER
1	D	110	LYS
1	D	119	ARG
1	D	155	ARG
1	D	177	LEU
1	D	260	LYS
1	D	265	ARG
1	D	291	TRP
1	D	306	LEU
1	D	314	LEU
1	D	319	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	55	ASN
1	A	85	GLN
1	A	126	ASN
1	A	182	GLN
1	A	321	GLN
1	B	55	ASN
1	B	126	ASN
1	B	182	GLN
1	B	321	GLN
1	C	85	GLN
1	C	126	ASN
1	C	187	GLN
1	D	55	ASN
1	D	85	GLN

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Mol	Chain	Res	Type
1	D	126	ASN
1	D	143	ASN
1	D	182	GLN
1	D	321	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 [i](#)

6 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	GOL	A	1329	-	5,5,5	0.44	0	5,5,5	0.27	0
3	NAP	A	1330	-	42,52,52	1.59	2 (4%)	54,80,80	1.82	5 (9%)
3	NAP	B	1330	-	42,52,52	1.62	3 (7%)	54,80,80	1.70	3 (5%)
3	NAP	C	1330	-	42,52,52	1.59	3 (7%)	54,80,80	1.80	2 (3%)
2	GOL	D	1328	-	5,5,5	0.39	0	5,5,5	0.42	0
3	NAP	D	1330	-	42,52,52	1.63	3 (7%)	54,80,80	1.85	4 (7%)

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	GOL	A	1329	-	-	0/4/4/4	0/0/0/0
3	NAP	A	1330	-	-	0/27/67/67	0/5/5/5
3	NAP	B	1330	-	-	0/27/67/67	0/5/5/5
3	NAP	C	1330	-	-	0/27/67/67	0/5/5/5
2	GOL	D	1328	-	-	0/4/4/4	0/0/0/0
3	NAP	D	1330	-	-	0/27/67/67	0/5/5/5

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	1330	NAP	C2A-N1A	2.11	1.37	1.33
3	C	1330	NAP	C2A-N1A	2.12	1.37	1.33
3	B	1330	NAP	C2A-N1A	2.14	1.38	1.33
3	A	1330	NAP	C2A-N3A	2.86	1.37	1.32
3	C	1330	NAP	C2A-N3A	3.14	1.37	1.32
3	B	1330	NAP	C2A-N3A	3.47	1.38	1.32
3	D	1330	NAP	C2A-N3A	3.54	1.38	1.32
3	A	1330	NAP	O7N-C7N	8.28	1.41	1.24
3	C	1330	NAP	O7N-C7N	8.33	1.41	1.24
3	D	1330	NAP	O7N-C7N	8.43	1.42	1.24
3	B	1330	NAP	O7N-C7N	8.49	1.42	1.24

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	1330	NAP	N3A-C2A-N1A	-11.71	119.93	128.89
3	C	1330	NAP	N3A-C2A-N1A	-11.26	120.27	128.89
3	A	1330	NAP	N3A-C2A-N1A	-11.13	120.38	128.89
3	B	1330	NAP	N3A-C2A-N1A	-10.55	120.82	128.89
3	B	1330	NAP	PN-O3-PA	-2.94	124.46	132.73
3	D	1330	NAP	C4A-C5A-N7A	-2.18	107.47	109.48
3	D	1330	NAP	PN-O3-PA	-2.17	126.65	132.73
3	A	1330	NAP	C1B-N9A-C4A	-2.16	123.68	126.94
3	C	1330	NAP	PN-O3-PA	-2.14	126.71	132.73
3	A	1330	NAP	PN-O3-PA	-2.07	126.91	132.73
3	B	1330	NAP	C4D-O4D-C1D	2.29	112.24	109.72
3	D	1330	NAP	O4D-C1D-N1N	2.33	110.69	108.13
3	A	1330	NAP	C4D-O4D-C1D	2.48	112.44	109.72
3	A	1330	NAP	O4D-C1D-N1N	2.95	111.38	108.13

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1330	NAP	3	0
3	B	1330	NAP	2	0
3	C	1330	NAP	1	0
2	D	1328	GOL	1	0
3	D	1330	NAP	2	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	281/302 (93%)	0.22	19 (6%)	20 28	13, 23, 57, 82	0
1	B	274/302 (90%)	0.11	9 (3%)	50 59	13, 24, 41, 88	0
1	C	277/302 (91%)	0.22	19 (6%)	20 27	14, 26, 52, 83	0
1	D	272/302 (90%)	0.20	14 (5%)	32 40	15, 27, 43, 64	0
All	All	1104/1208 (91%)	0.19	61 (5%)	29 37	13, 25, 51, 88	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	328	LEU	8.7
1	A	325	ILE	6.9
1	A	328	LEU	6.4
1	B	328	LEU	6.2
1	A	327	GLU	6.1
1	A	326	GLU	6.1
1	C	36	THR	5.5
1	A	324	THR	5.4
1	B	329	ILE	5.4
1	A	35	ASN	4.6
1	C	326	GLU	4.5
1	A	39	LEU	4.4
1	A	245	THR	4.4
1	C	325	ILE	4.2
1	B	260	LYS	4.1
1	A	252	LEU	4.1
1	A	38	ALA	3.9
1	C	327	GLU	3.8
1	D	245	THR	3.7
1	B	261	GLU	3.6
1	D	327	GLU	3.5

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Mol	Chain	Res	Type	RSRZ
1	B	95	VAL	3.5
1	C	41	SER	3.4
1	C	245	THR	3.4
1	A	41	SER	3.4
1	C	323	ASP	3.4
1	C	38	ALA	3.4
1	A	36	THR	3.2
1	B	327	GLU	3.1
1	C	324	THR	3.0
1	C	37	GLU	2.9
1	D	261	GLU	2.9
1	B	142	ILE	2.8
1	A	142	ILE	2.7
1	A	106	GLN	2.7
1	A	94[A]	ASP	2.6
1	B	105	SER	2.6
1	C	58[A]	GLN	2.5
1	C	39	LEU	2.5
1	A	323	ASP	2.4
1	B	94[A]	ASP	2.4
1	C	106	GLN	2.3
1	C	141	VAL	2.3
1	C	105	SER	2.3
1	C	94[A]	ASP	2.3
1	D	95	VAL	2.3
1	C	142	ILE	2.2
1	D	260	LYS	2.2
1	A	193	LEU	2.2
1	A	37	GLU	2.2
1	D	110	LYS	2.2
1	D	36	THR	2.2
1	D	64	ILE	2.2
1	A	253	ASP	2.1
1	D	185	LYS	2.1
1	D	141	VAL	2.1
1	C	55[A]	ASN	2.1
1	D	326	GLU	2.1
1	D	105	SER	2.1
1	D	142	ILE	2.0
1	D	284	LEU	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
2	GOL	A	1329	6/6	0.81	0.37	8.94	57,59,60,61	0
2	GOL	D	1328	6/6	0.85	0.34	3.94	41,43,43,46	0
3	NAP	B	1330	48/48	0.89	0.19	2.13	22,46,54,55	0
3	NAP	C	1330	48/48	0.89	0.17	1.34	20,54,57,59	0
3	NAP	D	1330	48/48	0.88	0.19	1.24	21,52,57,58	0
3	NAP	A	1330	48/48	0.92	0.16	0.72	15,44,51,52	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.