



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 17, 2020 – 09:13 AM BST

PDB ID : 4IW1  
Title : HSA-fructose complex  
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Deposited on : 2013-01-23  
Resolution : 2.56 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

<https://www.wwpdb.org/validation/2017/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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.13.1  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.13.1

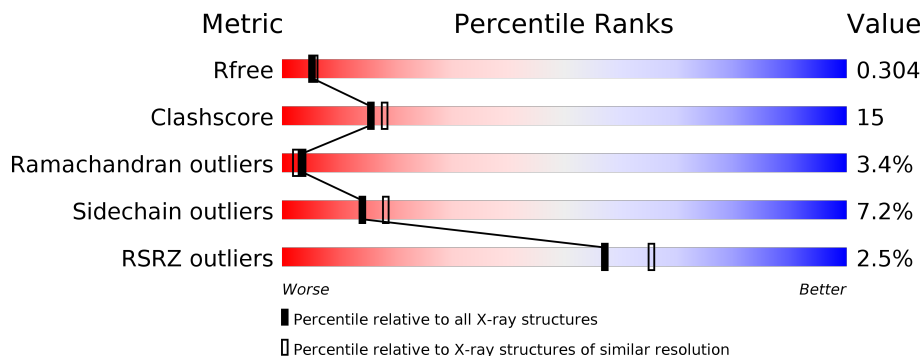
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.56 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))
$R_{free}$	130704	1279 (2.58-2.54)
Clashscore	141614	1327 (2.58-2.54)
Ramachandran outliers	138981	1312 (2.58-2.54)
Sidechain outliers	138945	1312 (2.58-2.54)
RSRZ outliers	127900	1269 (2.58-2.54)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. 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	585	

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
3	PO4	A	602	-	-	X	-
4	FUD	A	603	-	-	X	-

## 2 Entry composition [i](#)

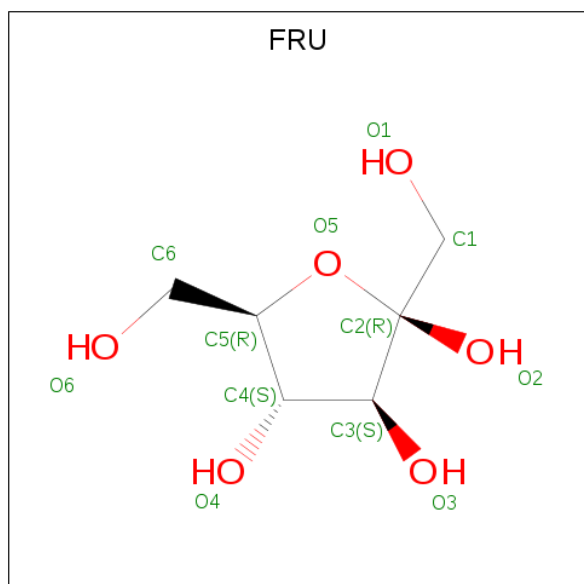
There are 5 unique types of molecules in this entry. The entry contains 4321 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Serum albumin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	560	4289	2716	715	818	40	0	1	0

- Molecule 2 is beta-D-fructofuranose (three-letter code: FRU) (formula: C<sub>6</sub>H<sub>12</sub>O<sub>6</sub>).



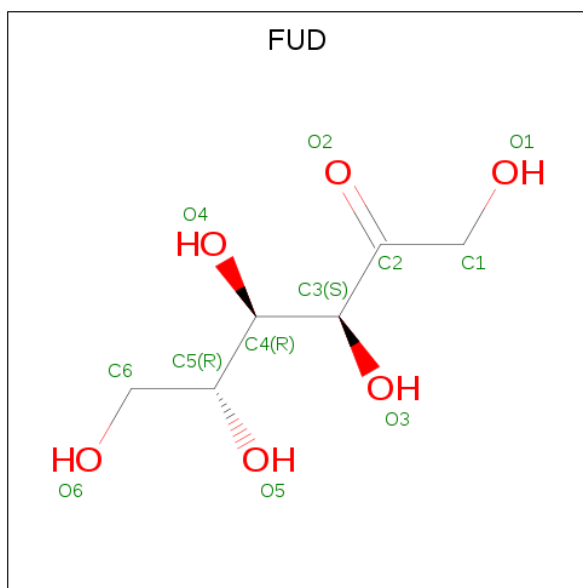
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	12	6	6	0	0

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			5	4	1		

- Molecule 4 is D-fructose (three-letter code: FUD) (formula: C<sub>6</sub>H<sub>12</sub>O<sub>6</sub>).



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

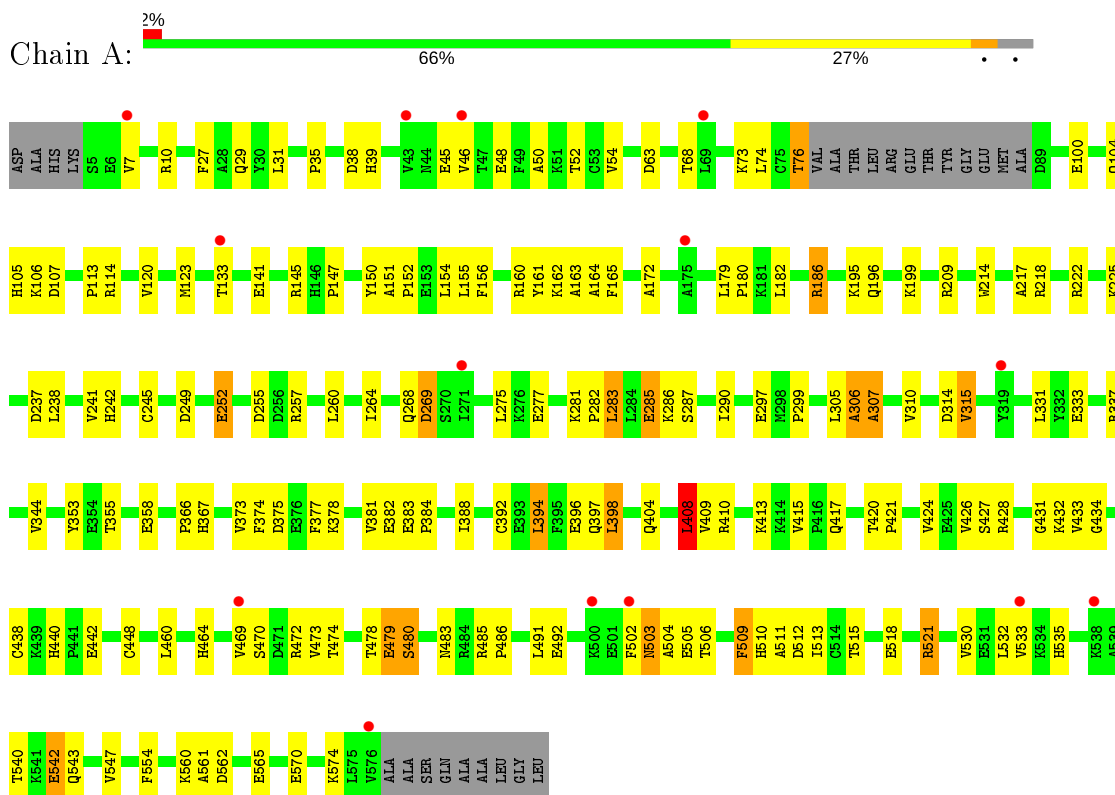
- Molecule 5 is water.

<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
5	A	3	Total	O	0	0
			3	3		

### 3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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: Serum albumin



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	58.36Å 87.98Å 59.35Å 90.00° 102.70° 90.00°	Depositor
Resolution (Å)	48.41 – 2.56 48.37 – 2.56	Depositor EDS
% Data completeness (in resolution range)	97.0 (48.41-2.56) 97.0 (48.37-2.56)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.36 (at 2.54Å)	Xtrriage
Refinement program	CNS 1.3, REFMAC 5.7.0032	Depositor
R, $R_{free}$	0.239 , 0.315 0.235 , 0.304	Depositor DCC
$R_{free}$ test set	954 reflections (5.16%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	54.7	Xtrriage
Anisotropy	0.376	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 46.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	0.017 for l,-k,h	Xtrriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	4321	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	63.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.04% 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, 0.333 respectively for untwinned datasets, and 0.375, 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: PO4, FUD, FRU

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	0.63	0/4376	0.80	2/5943 (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	2

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	408	LEU	CA-CB-CG	6.25	129.66	115.30
1	A	107	ASP	CB-CG-OD1	5.12	122.91	118.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	509	PHE	Peptide
1	A	560	LYS	Peptide

### 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	4289	0	4014	123	0
2	A	12	0	12	5	0
3	A	5	0	0	2	0
4	A	12	0	12	6	0
5	A	3	0	0	0	0
All	All	4321	0	4038	123	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 15.

All (123) 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:50:ALA:O	1:A:54:VAL:HG23	1.65	0.96
1:A:238:LEU:HD21	2:A:601:FRU:H61	1.55	0.89
1:A:510:HIS:CD2	1:A:512:ASP:HB3	2.11	0.85
1:A:238:LEU:HD13	2:A:601:FRU:O2	1.76	0.85
1:A:222:ARG:HH11	4:A:603:FUD:H4	1.42	0.84
1:A:29:GLN:HG2	1:A:147:PRO:HA	1.62	0.81
1:A:373:VAL:HG13	1:A:374:PHE:HD1	1.50	0.76
1:A:543:GLN:O	1:A:547:VAL:HG23	1.86	0.76
1:A:225:LYS:HD3	1:A:299:PRO:HG3	1.68	0.76
1:A:199:LYS:NZ	4:A:603:FUD:O6	2.16	0.75
1:A:141:GLU:OE1	1:A:145:ARG:NH1	2.20	0.70
1:A:156:PHE:CE2	1:A:160:ARG:CZ	2.75	0.69
1:A:424:VAL:O	1:A:428:ARG:HG3	1.91	0.69
1:A:151:ALA:HB3	1:A:152:PRO:HD3	1.73	0.69
1:A:397:GLN:O	1:A:398:LEU:HD23	1.93	0.69
1:A:156:PHE:HE1	1:A:285:GLU:HG3	1.58	0.68
1:A:225:LYS:HE2	1:A:297:GLU:O	1.93	0.68
1:A:509:PHE:HD1	1:A:513:ILE:CD1	2.06	0.68
1:A:485:ARG:HB3	1:A:486:PRO:HD3	1.75	0.68
1:A:409:VAL:O	1:A:413:LYS:HG3	1.94	0.67
1:A:540:THR:HG22	1:A:542:GLU:H	1.60	0.67
1:A:156:PHE:CE2	1:A:160:ARG:NH2	2.64	0.66
1:A:196:GLN:HE22	1:A:242:HIS:CE1	2.13	0.66
1:A:238:LEU:CD2	2:A:601:FRU:H61	2.26	0.65
1:A:570:GLU:O	1:A:574:LYS:N	2.25	0.64
1:A:464:HIS:HE1	1:A:470:SER:H	1.45	0.64
1:A:305:LEU:HD21	1:A:333:GLU:HB2	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:260:LEU:O	1:A:264:ILE:HG13	1.98	0.63
1:A:518:GLU:OE2	1:A:521:ARG:NH1	2.32	0.63
1:A:434:GLY:HA2	1:A:438:CYS:SG	2.39	0.63
1:A:381:VAL:O	1:A:384:PRO:HD2	1.99	0.61
1:A:305:LEU:HD21	1:A:333:GLU:CB	2.30	0.61
1:A:509:PHE:CD1	1:A:513:ILE:HD13	2.36	0.60
1:A:417:GLN:HB2	1:A:470:SER:HB2	1.83	0.59
1:A:510:HIS:O	1:A:512:ASP:N	2.35	0.59
1:A:408:LEU:HD23	1:A:427:SER:HB2	1.84	0.59
1:A:182:LEU:O	1:A:186:ARG:HG2	2.03	0.59
1:A:196:GLN:HE22	1:A:242:HIS:HE1	1.49	0.59
1:A:214:TRP:NE1	3:A:602:PO4:O4	2.36	0.59
1:A:283:LEU:O	1:A:283:LEU:HD12	2.03	0.59
1:A:31:LEU:HD12	1:A:74:LEU:HD22	1.86	0.58
1:A:509:PHE:HD1	1:A:513:ILE:HD13	1.68	0.58
1:A:195:LYS:NZ	4:A:603:FUD:O3	2.36	0.58
1:A:373:VAL:HG13	1:A:374:PHE:CD1	2.37	0.56
1:A:408:LEU:HD11	1:A:530:VAL:HG23	1.86	0.56
1:A:156:PHE:CE1	1:A:285:GLU:HG3	2.40	0.56
1:A:464:HIS:CE1	1:A:469:VAL:H	2.24	0.56
1:A:237:ASP:O	1:A:241:VAL:HG23	2.06	0.55
1:A:472:ARG:HH21	1:A:491:LEU:HD22	1.71	0.55
1:A:474:THR:O	1:A:478:THR:HG23	2.06	0.55
1:A:479:GLU:O	1:A:480:SER:HB2	2.07	0.55
1:A:333:GLU:OE1	1:A:337:ARG:NH2	2.40	0.54
1:A:503:ASN:HD22	1:A:504:ALA:N	2.07	0.53
1:A:502:PHE:HB2	1:A:535:HIS:CE1	2.45	0.52
1:A:222:ARG:NH1	4:A:603:FUD:H4	2.17	0.52
1:A:182:LEU:HD22	1:A:186:ARG:HH21	1.74	0.52
1:A:464:HIS:CE1	1:A:470:SER:H	2.25	0.52
1:A:249:ASP:HB3	1:A:252:GLU:CD	2.30	0.52
1:A:415:VAL:HG11	1:A:473:VAL:HG23	1.91	0.51
1:A:269:ASP:N	1:A:269:ASP:OD1	2.42	0.51
1:A:161:TYR:CZ	1:A:165:PHE:HE2	2.28	0.51
1:A:426:VAL:HG21	1:A:460:LEU:HB2	1.92	0.51
1:A:540:THR:HB	1:A:543:GLN:HG2	1.93	0.51
1:A:46:VAL:HG23	1:A:73:LYS:HG3	1.93	0.50
1:A:73:LYS:O	1:A:76:THR:HG23	2.12	0.50
1:A:388:ILE:O	1:A:392:CYS:HB2	2.11	0.49
1:A:218:ARG:NH1	1:A:222:ARG:HG3	2.27	0.49
1:A:106:LYS:HD3	1:A:147:PRO:HB2	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:392:CYS:O	1:A:396:GLU:HG2	2.12	0.49
1:A:353:TYR:HD1	1:A:373:VAL:HG21	1.77	0.49
1:A:10:ARG:NH1	1:A:255:ASP:OD2	2.45	0.49
1:A:290:ILE:HG22	2:A:601:FRU:H3	1.95	0.49
1:A:404:GLN:O	1:A:408:LEU:HB2	2.13	0.48
1:A:394:LEU:O	1:A:398:LEU:HB2	2.14	0.48
1:A:225:LYS:HD3	1:A:299:PRO:CG	2.41	0.48
1:A:252:GLU:H	1:A:252:GLU:CD	2.11	0.48
1:A:355:THR:O	1:A:358:GLU:HB3	2.14	0.47
1:A:533:VAL:HG12	1:A:533:VAL:O	2.14	0.47
1:A:45:GLU:O	1:A:48[B]:GLU:HB3	2.14	0.47
1:A:27:PHE:HB3	1:A:39:HIS:ND1	2.28	0.47
1:A:333:GLU:HB3	1:A:337:ARG:NH2	2.30	0.47
1:A:161:TYR:CE1	1:A:165:PHE:HE2	2.32	0.46
1:A:472:ARG:NH2	1:A:491:LEU:HD22	2.30	0.46
1:A:509:PHE:HD1	1:A:513:ILE:HD11	1.80	0.46
1:A:283:LEU:C	1:A:283:LEU:HD12	2.35	0.46
1:A:503:ASN:ND2	1:A:505:GLU:H	2.14	0.46
1:A:333:GLU:HB3	1:A:337:ARG:HH21	1.80	0.45
1:A:305:LEU:HD21	1:A:333:GLU:HB3	1.98	0.45
1:A:518:GLU:HA	1:A:518:GLU:OE1	2.16	0.45
1:A:217:ALA:HB2	1:A:331:LEU:HD11	1.98	0.45
1:A:225:LYS:HA	1:A:299:PRO:HG3	1.97	0.45
1:A:162:LYS:O	1:A:164:ALA:N	2.50	0.45
1:A:156:PHE:CD2	1:A:160:ARG:NH2	2.85	0.45
1:A:509:PHE:CD1	1:A:513:ILE:CD1	2.92	0.44
1:A:277:GLU:O	1:A:281:LYS:HG3	2.17	0.44
1:A:120:VAL:O	1:A:123:MET:N	2.49	0.44
1:A:377:PHE:O	1:A:378:LYS:C	2.56	0.43
1:A:373:VAL:HG13	1:A:374:PHE:N	2.33	0.43
1:A:306:ALA:HA	1:A:310:VAL:HB	2.00	0.43
1:A:420:THR:HB	1:A:421:PRO:HD3	1.98	0.43
1:A:225:LYS:NZ	1:A:299:PRO:HA	2.33	0.43
1:A:186:ARG:HG2	1:A:186:ARG:H	1.68	0.42
1:A:100:GLU:O	1:A:104:GLN:HB2	2.19	0.42
1:A:154:LEU:HD12	1:A:154:LEU:HA	1.62	0.42
1:A:214:TRP:CE2	3:A:602:PO4:O4	2.72	0.42
1:A:257:ARG:CZ	1:A:287:SER:HB3	2.50	0.42
1:A:502:PHE:HB2	1:A:535:HIS:NE2	2.34	0.42
1:A:431:GLY:C	1:A:433:VAL:H	2.22	0.41
1:A:268:GLN:HB2	1:A:275:LEU:HD12	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:151:ALA:HB3	1:A:152:PRO:CD	2.46	0.41
1:A:238:LEU:HD21	2:A:601:FRU:C6	2.37	0.41
1:A:483:ASN:O	1:A:486:PRO:HD2	2.19	0.41
1:A:314:ASP:O	1:A:315:VAL:C	2.57	0.41
1:A:154:LEU:O	1:A:155:LEU:C	2.59	0.41
1:A:218:ARG:HH21	4:A:603:FUD:C2	2.32	0.41
1:A:179:LEU:HB2	1:A:180:PRO:HD3	2.03	0.41
1:A:218:ARG:NH2	4:A:603:FUD:H11	2.36	0.41
1:A:366:PRO:O	1:A:367:HIS:C	2.60	0.40
1:A:533:VAL:O	1:A:533:VAL:CG1	2.69	0.40
1:A:306:ALA:O	1:A:307:ALA:C	2.59	0.40
1:A:156:PHE:HE2	1:A:160:ARG:NH2	2.15	0.40
1:A:305:LEU:O	1:A:307:ALA:N	2.54	0.40
1:A:281:LYS:HB3	1:A:282:PRO:HD2	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	557/585 (95%)	492 (88%)	46 (8%)	19 (3%)	<b>3</b> <b>2</b>

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	306	ALA
1	A	442	GLU
1	A	480	SER
1	A	511	ALA
1	A	561	ALA
1	A	163	ALA
1	A	432	LYS

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Mol	Chain	Res	Type
1	A	150	TYR
1	A	172	ALA
1	A	307	ALA
1	A	114	ARG
1	A	410	ARG
1	A	506	THR
1	A	562	ASP
1	A	35	PRO
1	A	448	CYS
1	A	113	PRO
1	A	440	HIS
1	A	315	VAL

### 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	446/511 (87%)	414 (93%)	32 (7%)	<b>14</b> <b>18</b>

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

Mol	Chain	Res	Type
1	A	7	VAL
1	A	38	ASP
1	A	52	THR
1	A	63	ASP
1	A	68	THR
1	A	76	THR
1	A	105	HIS
1	A	133	THR
1	A	186	ARG
1	A	209	ARG
1	A	245	CYS
1	A	252	GLU
1	A	269	ASP
1	A	283	LEU

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Mol	Chain	Res	Type
1	A	285	GLU
1	A	286	LYS
1	A	344	VAL
1	A	375	ASP
1	A	382	GLU
1	A	383	GLU
1	A	394	LEU
1	A	398	LEU
1	A	408	LEU
1	A	479	GLU
1	A	492	GLU
1	A	503	ASN
1	A	515	THR
1	A	521	ARG
1	A	532	LEU
1	A	542	GLU
1	A	554	PHE
1	A	565	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	109	ASN
1	A	196	GLN
1	A	385	GLN
1	A	440	HIS
1	A	459	GLN
1	A	464	HIS
1	A	483	ASN
1	A	503	ASN
1	A	510	HIS

### 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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

3 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
3	PO4	A	602	-	4,4,4	1.03	0	6,6,6	0.41	0
2	FRU	A	601	-	11,12,12	0.70	0	10,18,18	1.83	3 (30%)
4	FUD	A	603	-	10,11,11	0.43	0	9,14,14	0.86	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	FRU	A	601	-	-	0/5/24/24	0/1/1/1
4	FUD	A	603	-	-	6/16/16/16	-

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	FRU	O1-C1-C2	-4.06	103.23	111.86
2	A	601	FRU	O3-C3-C4	-3.31	101.89	113.32
2	A	601	FRU	O4-C4-C3	-2.10	105.87	112.15

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	603	FUD	O1-C1-C2-C3
4	A	603	FUD	O1-C1-C2-O2
4	A	603	FUD	C1-C2-C3-O3
4	A	603	FUD	O2-C2-C3-O3
4	A	603	FUD	O5-C5-C6-O6
4	A	603	FUD	C4-C5-C6-O6

There are no ring outliers.

3 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	602	PO4	2	0
2	A	601	FRU	5	0
4	A	603	FUD	6	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](#)

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	560/585 (95%)	0.13	14 (2%) 57 65	33, 61, 93, 121	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	576	VAL	4.3
1	A	538	LYS	3.4
1	A	500	LYS	3.2
1	A	271	ILE	2.8
1	A	7	VAL	2.7
1	A	502	PHE	2.6
1	A	46	VAL	2.5
1	A	533	VAL	2.4
1	A	319	TYR	2.4
1	A	133	THR	2.3
1	A	469	VAL	2.3
1	A	69	LEU	2.2
1	A	43	VAL	2.2
1	A	175	ALA	2.2

### 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 monosaccharides 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. 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	B-factors(Å <sup>2</sup> )	Q<0.9
2	FRU	A	601	12/12	0.79	0.36	66,83,90,107	0
4	FUD	A	603	12/12	0.89	0.15	70,87,95,101	0
3	PO4	A	602	5/5	0.94	0.33	102,104,110,126	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.