



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 9, 2018 – 12:13 PM EST

PDB ID : 5WZX  
Title : Structural basis for a pentacyclic oleanane-type triterpenoid as a ligand of FXR  
Authors : Lu, Y.; Li, Y.  
Deposited on : 2017-01-19  
Resolution : 2.95 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20030736  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20030736

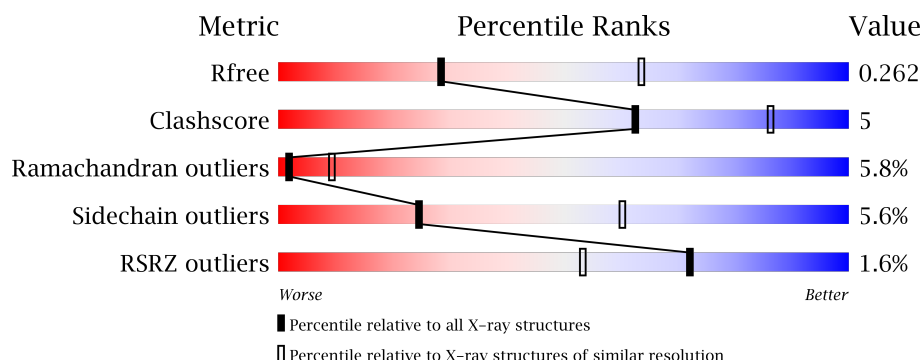
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.95 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	100719	2395 (3.00-2.92)
Clashscore	112137	2773 (3.00-2.92)
Ramachandran outliers	110173	2680 (3.00-2.92)
Sidechain outliers	110143	2683 (3.00-2.92)
RSRZ outliers	101464	2421 (3.00-2.92)

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	228	<div> <div>0.1%</div> <div>79% 12% 8%</div> </div>
1	B	228	<div> <div>2%</div> <div>71% 18% 7%</div> </div>
2	C	12	<div> <div>75% 17% 8%</div> </div>
2	D	12	<div> <div>8%</div> <div>67% 25% 8%</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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	7VX	A	501	-	-	-	X
3	7VX	B	501	-	-	-	X

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 3784 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.

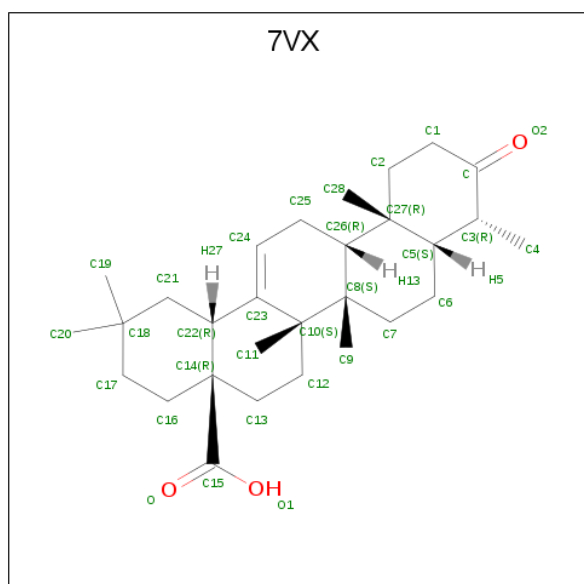
- Molecule 1 is a protein called Bile acid receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	210	Total	C	N	O	S	0	0	0
			1725	1104	292	318	11			
1	B	212	Total	C	N	O	S	0	0	0
			1747	1120	295	321	11			

- Molecule 2 is a protein called SRC2-3 peptide from Nuclear receptor coactivator 2.

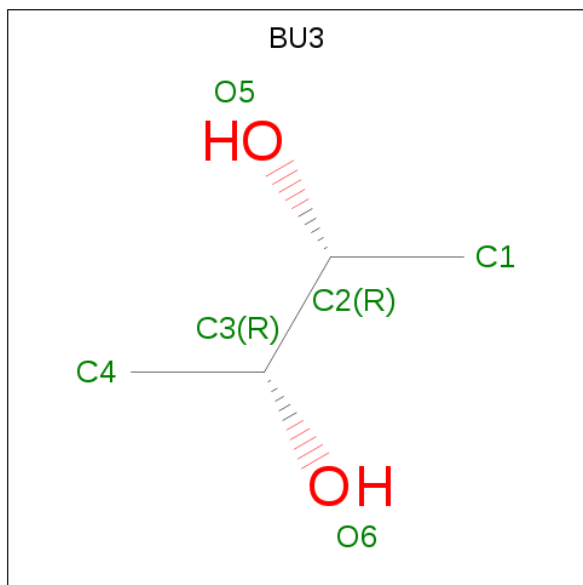
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	12	Total	C	N	O	0	0	0
			102	65	17	20			
2	D	11	Total	C	N	O	0	0	0
			90	59	15	16			

- Molecule 3 is (4aR,6aR,6aS,6bS,8aS,9R,12aR,14bR)-2,2,6a,6b,9,12a-hexamethyl-10-oxidanylidene-1,3,4,5,6,6a,7,8,8a,9,11,12,13,14b-tetradecahydronicene-4a-carboxylic acid (three-letter code: 7VX) (formula: C<sub>29</sub>H<sub>44</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			32	29	3		
3	B	1	Total	C	O	0	0
			32	29	3		

- Molecule 4 is (R,R)-2,3-BUTANEDIOL (three-letter code: BU3) (formula: C<sub>4</sub>H<sub>10</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	4	2		
4	A	1	Total	C	O	0	0
			6	4	2		
4	A	1	Total	C	O	0	0
			6	4	2		

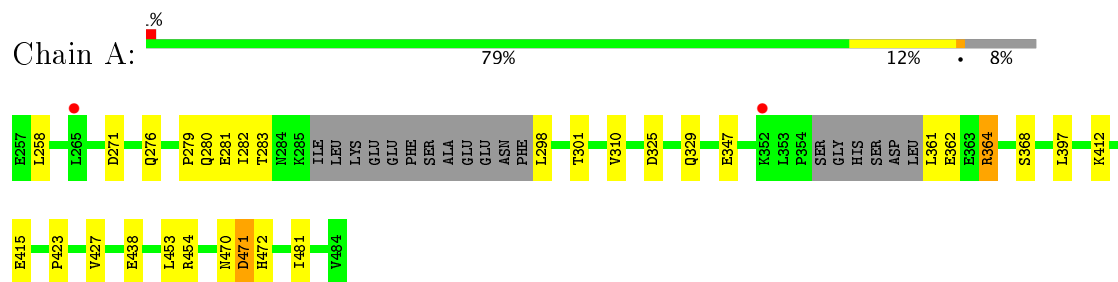
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	17	Total	O	0	0
			17	17		
5	B	20	Total	O	0	0
			20	20		
5	D	1	Total	O	0	0
			1	1		

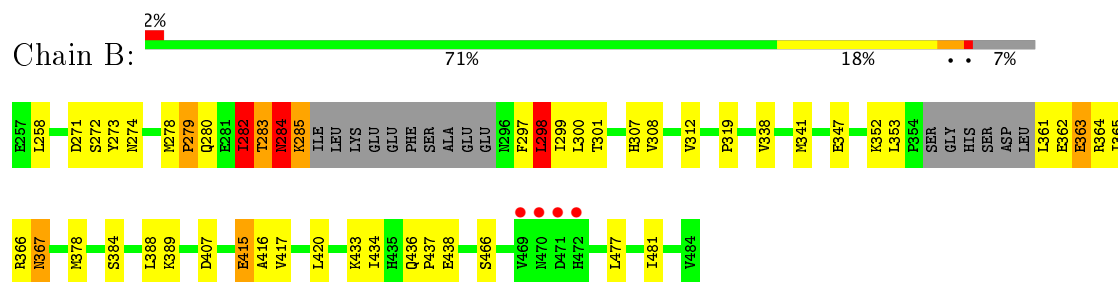
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

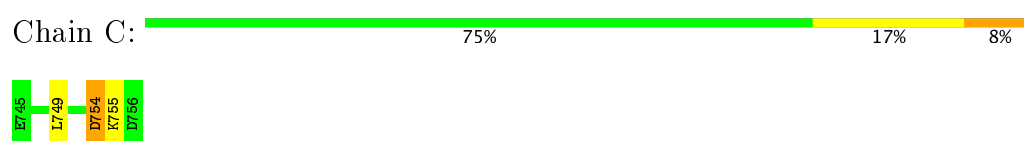
- Molecule 1: Bile acid receptor



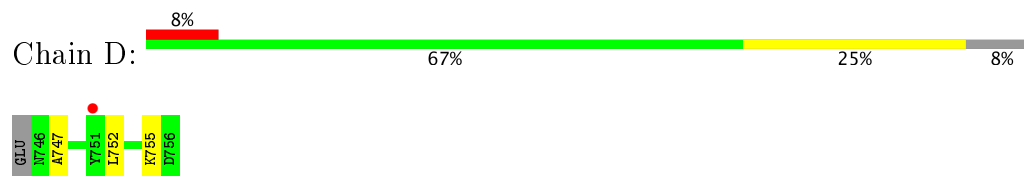
- Molecule 1: Bile acid receptor



- Molecule 2: SRC2-3 peptide from Nuclear receptor coactivator 2



- Molecule 2: SRC2-3 peptide from Nuclear receptor coactivator 2



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	74.56 Å 86.08 Å 177.30 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	88.65 – 2.95 47.56 – 2.95	Depositor EDS
% Data completeness (in resolution range)	86.5 (88.65-2.95) 86.5 (47.56-2.95)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.52 (at 2.96 Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, $R_{free}$	0.208 , 0.263 0.211 , 0.262	Depositor DCC
$R_{free}$ test set	514 reflections (5.04%)	DCC
Wilson B-factor (Å <sup>2</sup> )	83.0	Xtriage
Anisotropy	0.064	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 51.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	3784	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	86.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.17% 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: BU3, 7VX

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	0.59	0/1760	0.81	0/2376
1	B	0.60	0/1783	0.86	2/2407 (0.1%)
2	C	0.76	0/102	0.98	0/136
2	D	0.77	0/90	0.85	0/120
All	All	0.61	0/3735	0.84	2/5039 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	274	ASN	N-CA-C	6.49	128.51	111.00
1	B	274	ASN	C-N-CA	5.42	135.24	121.70

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	1725	0	1729	14	0
1	B	1747	0	1753	21	0
2	C	102	0	103	1	0
2	D	90	0	93	1	0
3	A	32	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	32	0	0	0	0
4	A	18	0	30	1	0
5	A	17	0	0	0	0
5	B	20	0	0	2	0
5	D	1	0	0	0	0
All	All	3784	0	3708	37	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 (37) 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:388:LEU:N	5:B:601:HOH:O	2.13	0.82
1:B:365:ILE:HA	1:B:367:ASN:HD22	1.54	0.72
1:B:384:SER:O	5:B:601:HOH:O	2.12	0.67
1:B:415:GLU:O	1:B:417:VAL:N	2.27	0.67
1:B:298:LEU:O	1:B:300:LEU:N	2.32	0.62
1:B:363:GLU:OE1	1:B:366:ARG:NH1	2.32	0.62
1:B:365:ILE:HA	1:B:367:ASN:ND2	2.18	0.58
1:A:258:LEU:HD11	1:A:427:VAL:HG23	1.86	0.58
1:B:282:ILE:HD12	1:B:352:LYS:HE2	1.89	0.54
1:A:276:GLN:CB	1:A:310:VAL:HG13	2.38	0.54
1:B:273:TYR:OH	1:B:347:GLU:OE1	2.20	0.54
1:A:276:GLN:HB3	1:A:310:VAL:HG13	1.90	0.54
1:A:298:LEU:HD22	1:A:364:ARG:HH21	1.74	0.52
1:B:282:ILE:O	1:B:283:THR:HG23	2.09	0.52
1:A:423:PRO:O	1:A:427:VAL:HG12	2.11	0.51
1:A:298:LEU:HD22	1:A:364:ARG:NH2	2.25	0.51
1:A:471:ASP:O	1:A:471:ASP:CG	2.49	0.51
1:B:284:ASN:OD1	1:B:284:ASN:N	2.36	0.50
1:B:308:VAL:HG11	1:B:477:LEU:HD11	1.95	0.48
1:A:470:ASN:O	1:A:472:HIS:N	2.47	0.47
2:C:754:ASP:OD1	2:C:754:ASP:N	2.45	0.46
1:A:325:ASP:O	1:A:329:GLN:HG3	2.16	0.45
1:A:361:LEU:HD11	4:A:502:BU3:H2	1.99	0.45
1:A:368:SER:HB2	3:A:501:7VX:C4	2.47	0.45
1:A:279:PRO:O	1:A:281:GLU:N	2.49	0.45
3:A:501:7VX:C11	3:A:501:7VX:C15	2.96	0.44
1:B:319:PRO:O	1:B:420:LEU:HD22	2.19	0.43
1:B:278:MET:HG2	1:B:279:PRO:CD	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:361:LEU:O	1:B:361:LEU:HD12	2.19	0.42
1:B:312:VAL:HG11	2:D:752:LEU:HB2	2.01	0.42
1:A:415:GLU:CD	1:A:415:GLU:N	2.73	0.42
1:A:453:LEU:O	1:A:454:ARG:C	2.59	0.41
1:B:436:GLN:N	1:B:437:PRO:HD3	2.36	0.41
1:B:362:GLU:O	1:B:364:ARG:N	2.53	0.41
1:B:361:LEU:HA	1:B:362:GLU:HA	1.86	0.41
1:B:283:THR:C	1:B:285:LYS:H	2.23	0.41
1:B:307:HIS:HB3	1:B:341:MET:HG2	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	204/228 (90%)	186 (91%)	13 (6%)	5 (2%)	6	29
1	B	206/228 (90%)	169 (82%)	20 (10%)	17 (8%)	1	4
2	C	10/12 (83%)	8 (80%)	1 (10%)	1 (10%)	1	2
2	D	9/12 (75%)	7 (78%)	0	2 (22%)	0	0
All	All	429/480 (89%)	370 (86%)	34 (8%)	25 (6%)	2	9

All (25) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	280	GLN
1	A	471	ASP
1	B	279	PRO
1	B	280	GLN
1	B	299	ILE
1	B	363	GLU

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Mol	Chain	Res	Type
1	B	367	ASN
1	B	416	ALA
1	A	438	GLU
1	B	283	THR
1	B	297	PHE
1	B	298	LEU
1	B	353	LEU
1	B	415	GLU
1	B	438	GLU
2	D	755	LYS
1	B	284	ASN
2	D	747	ALA
1	A	283	THR
1	B	407	ASP
1	B	272	SER
2	C	755	LYS
1	B	282	ILE
1	A	282	ILE
1	B	434	ILE

### 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	194/211 (92%)	186 (96%)	8 (4%)	35	71
1	B	197/211 (93%)	184 (93%)	13 (7%)	19	53
2	C	11/11 (100%)	9 (82%)	2 (18%)	2	8
2	D	9/11 (82%)	9 (100%)	0	100	100
All	All	411/444 (93%)	388 (94%)	23 (6%)	25	60

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	271	ASP
1	A	301	THR

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Mol	Chain	Res	Type
1	A	347	GLU
1	A	362	GLU
1	A	364	ARG
1	A	397	LEU
1	A	412	LYS
1	A	481	ILE
1	B	258	LEU
1	B	271	ASP
1	B	282	ILE
1	B	284	ASN
1	B	285	LYS
1	B	298	LEU
1	B	301	THR
1	B	338	VAL
1	B	378	MET
1	B	389	LYS
1	B	433	LYS
1	B	466	SER
1	B	481	ILE
2	C	749	LEU
2	C	754	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	367	ASN
1	B	421	GLN

### 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

5 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	7VX	A	501	-	33,36,36	0.93	2 (6%)	56,62,62	2.81	26 (46%)
4	BU3	A	502	-	4,5,5	0.59	0	6,6,6	0.43	0
4	BU3	A	503	-	4,5,5	0.50	0	6,6,6	0.61	0
4	BU3	A	504	-	4,5,5	0.44	0	6,6,6	0.93	0
3	7VX	B	501	-	33,36,36	0.88	1 (3%)	56,62,62	2.68	26 (46%)

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
3	7VX	A	501	-	-	0/0/93/93	0/5/5/5
4	BU3	A	502	-	-	0/4/4/4	0/0/0/0
4	BU3	A	503	-	-	0/4/4/4	0/0/0/0
4	BU3	A	504	-	-	0/4/4/4	0/0/0/0
3	7VX	B	501	-	-	0/0/93/93	0/5/5/5

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	501	7VX	C8-C10	-2.40	1.54	1.58
3	B	501	7VX	C17-C18	-2.22	1.50	1.53
3	A	501	7VX	C6-C5	2.00	1.57	1.53

All (52) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	501	7VX	C28-C27-C5	-5.40	103.44	111.16
3	B	501	7VX	C28-C27-C26	-4.46	104.69	112.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	501	7VX	C12-C10-C23	-4.36	106.16	111.64
3	B	501	7VX	C19-C18-C17	-3.99	103.87	110.03
3	A	501	7VX	C9-C8-C10	-3.69	105.70	110.01
3	B	501	7VX	C4-C3-C	-3.60	106.13	112.15
3	B	501	7VX	C28-C27-C5	-3.57	106.06	111.16
3	B	501	7VX	C11-C10-C8	-3.47	109.00	112.34
3	A	501	7VX	C14-C22-C23	-3.45	109.92	112.17
3	A	501	7VX	C11-C10-C8	-3.43	109.04	112.34
3	B	501	7VX	C16-C17-C18	-3.25	109.88	113.18
3	A	501	7VX	C12-C10-C23	-3.10	107.75	111.64
3	B	501	7VX	C9-C8-C7	-3.09	102.37	107.90
3	B	501	7VX	O2-C-C3	-3.05	118.40	122.06
3	B	501	7VX	C10-C8-C26	-3.01	103.48	108.02
3	A	501	7VX	C7-C8-C26	-2.94	105.59	109.06
3	B	501	7VX	C16-C14-C15	-2.88	103.01	110.95
3	A	501	7VX	C28-C27-C26	-2.81	107.74	112.94
3	A	501	7VX	C10-C8-C26	-2.80	103.79	108.02
3	A	501	7VX	C9-C8-C7	-2.56	103.31	107.90
3	A	501	7VX	C16-C14-C22	-2.47	107.07	109.80
3	B	501	7VX	C8-C26-C27	-2.43	113.73	116.74
3	A	501	7VX	C4-C3-C	-2.35	108.22	112.15
3	A	501	7VX	C2-C1-C	-2.02	107.88	112.70
3	B	501	7VX	C20-C18-C17	2.05	113.19	110.03
3	B	501	7VX	C12-C10-C8	2.09	113.09	110.50
3	B	501	7VX	C21-C22-C14	2.13	114.25	112.02
3	A	501	7VX	C25-C26-C27	2.41	116.97	113.36
3	B	501	7VX	C13-C12-C10	2.59	117.73	113.81
3	A	501	7VX	C18-C21-C22	2.61	118.16	114.39
3	A	501	7VX	C16-C17-C18	2.73	115.95	113.18
3	A	501	7VX	C26-C27-C5	2.79	113.43	107.65
3	A	501	7VX	C13-C12-C10	2.85	118.12	113.81
3	B	501	7VX	C26-C27-C5	2.98	113.82	107.65
3	A	501	7VX	C1-C-C3	3.03	120.69	116.08
3	B	501	7VX	C7-C6-C5	3.07	118.47	113.12
3	B	501	7VX	C6-C5-C27	3.10	115.64	111.57
3	B	501	7VX	C4-C3-C5	3.36	120.00	113.97
3	B	501	7VX	C8-C10-C23	3.49	113.43	108.79
3	A	501	7VX	C7-C6-C5	3.65	119.48	113.12
3	A	501	7VX	C12-C10-C8	3.75	115.14	110.50
3	B	501	7VX	C16-C14-C22	3.88	114.09	109.80
3	A	501	7VX	C6-C5-C3	3.90	120.62	113.89
3	A	501	7VX	C5-C3-C	3.98	117.32	109.98

*Continued on next page...*

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	501	7VX	C9-C8-C26	4.11	118.20	111.85
3	B	501	7VX	C1-C-C3	4.29	122.62	116.08
3	A	501	7VX	C9-C8-C26	4.51	118.83	111.85
3	A	501	7VX	C2-C27-C5	4.58	114.50	107.79
3	B	501	7VX	C2-C27-C5	5.08	115.24	107.79
3	A	501	7VX	C17-C18-C21	5.42	115.73	108.69
3	B	501	7VX	C7-C8-C10	7.62	118.36	110.25
3	A	501	7VX	C7-C8-C10	9.74	120.62	110.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	501	7VX	2	0
4	A	502	BU3	1	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	210/228 (92%)	-0.09	2 (0%) 82 66	49, 80, 126, 153	0
1	B	212/228 (92%)	-0.07	4 (1%) 67 47	52, 83, 134, 174	0
2	C	12/12 (100%)	-0.12	0 100 100	59, 70, 97, 104	0
2	D	11/12 (91%)	0.16	1 (9%) 10 5	66, 86, 97, 103	0
All	All	445/480 (92%)	-0.07	7 (1%) 72 53	49, 81, 131, 174	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	352	LYS	3.3
1	B	471	ASP	3.0
1	B	469	VAL	2.9
1	B	470	ASN	2.6
2	D	751	TYR	2.1
1	A	265	LEU	2.0
1	B	472	HIS	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.



## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	7VX	B	501	32/32	0.91	0.37	3.12	93,112,128,137	0
3	7VX	A	501	32/32	0.94	0.31	2.30	85,93,111,123	0
4	BU3	A	503	6/6	0.73	0.27	-	89,105,106,107	0
4	BU3	A	504	6/6	0.76	0.35	-	99,100,109,110	0
4	BU3	A	502	6/6	0.80	0.19	-	125,132,136,137	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.