



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 11:53 am GMT

PDB ID : 1GWQ  
Title : HUMAN OESTROGEN RECEPTOR ALPHA LIGAND-BINDING DOMAIN IN COMPLEX WITH RALOXIFENE CORE AND TIF2 NRBOX2 PEPTIDE  
Authors : Pike, A.C.W.; Brzozowski, A.M.  
Deposited on : 2002-03-22  
Resolution : 2.45 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 : trunk28620  
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) : recalc28949

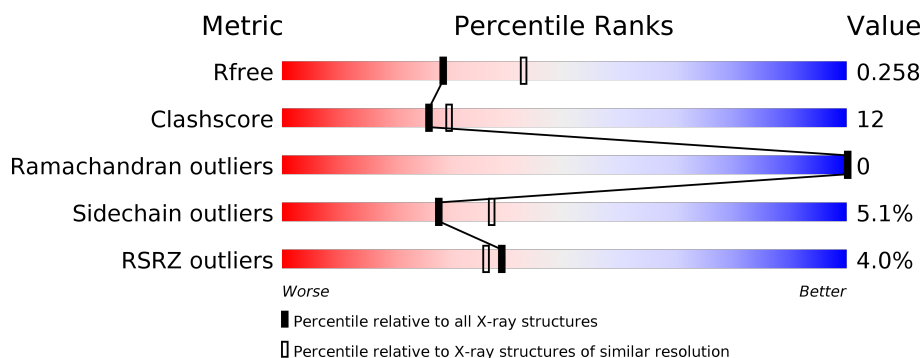
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.45 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	1119 (2.48-2.44)
Clashscore	112137	1193 (2.48-2.44)
Ramachandran outliers	110173	1185 (2.48-2.44)
Sidechain outliers	110143	1185 (2.48-2.44)
RSRZ outliers	101464	1126 (2.48-2.44)

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	248	<div> <div>2%</div> <div>68%</div> <div>26%</div> <div>• •</div> </div>
1	B	248	<div> <div>5%</div> <div>75%</div> <div>19%</div> <div>• •</div> </div>
2	C	9	<div> <div>78%</div> <div>22%</div> </div>
2	D	9	<div> <div>11%</div> <div>67%</div> <div>22%</div> <div>11%</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	ZTW	B	600	-	-	-	X

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4192 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.

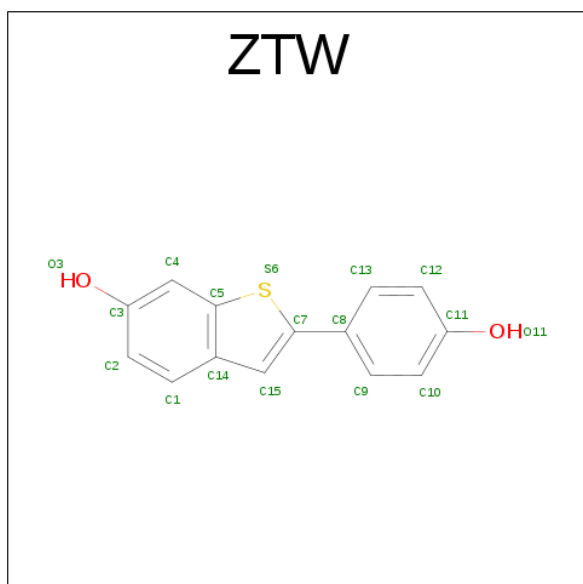
- Molecule 1 is a protein called OESTROGEN RECEPTOR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	243	Total	C	N	O	S	0	6	0
			1952	1246	337	349	20			
1	B	242	Total	C	N	O	S	7	3	0
			1938	1240	334	345	19			

- Molecule 2 is a protein called NUCLEAR RECEPTOR COACTIVATOR 2.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	9	Total	C	N	O	0	0	0
			73	47	15	11			
2	D	9	Total	C	N	O	0	0	0
			76	48	15	13			

- Molecule 3 is RALOXIFENE CORE (three-letter code: ZTW) (formula: C<sub>14</sub>H<sub>10</sub>O<sub>2</sub>S).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	S	0	0
			17	14	2	1		
3	B	1	Total	C	O	S	0	0
			17	14	2	1		

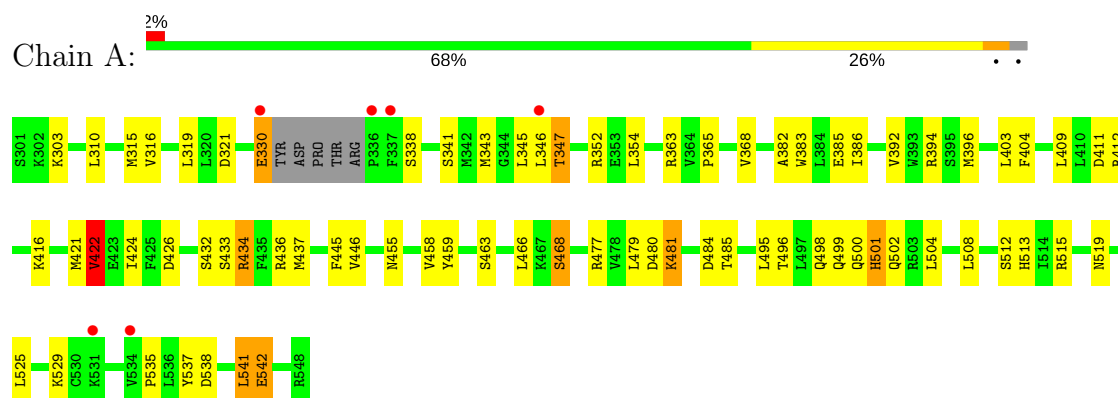
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	59	Total	O	0	0
			59	59		
4	B	60	Total	O	0	0
			60	60		

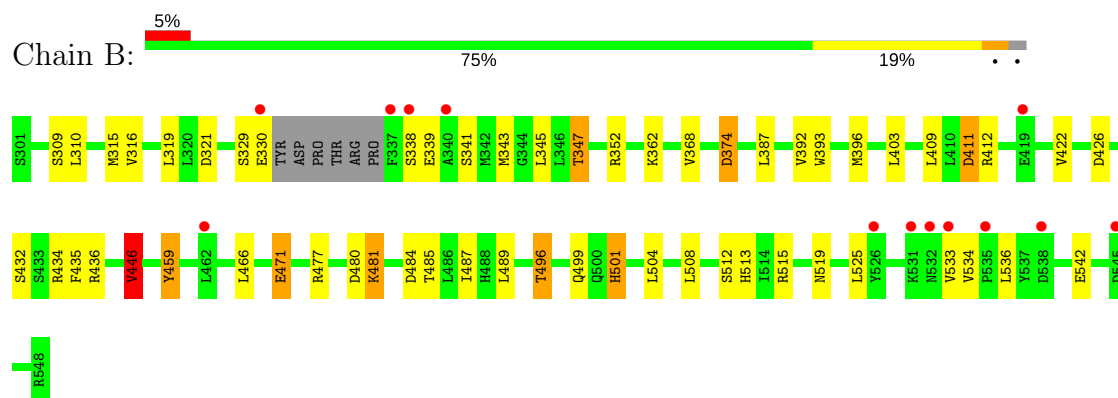
### 3 Residue-property plots [i](#)

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.

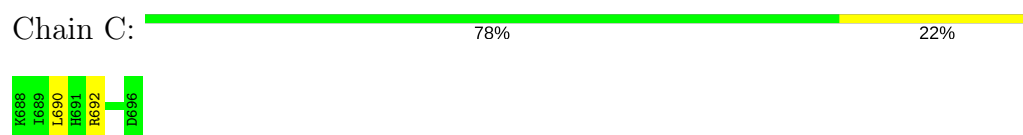
#### • Molecule 1: OESTROGEN RECEPTOR



#### • Molecule 1: OESTROGEN RECEPTOR

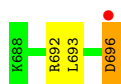


#### • Molecule 2: NUCLEAR RECEPTOR COACTIVATOR 2



#### • Molecule 2: NUCLEAR RECEPTOR COACTIVATOR 2





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	54.82Å 63.62Å 74.72Å 90.00° 98.34° 90.00°	Depositor
Resolution (Å)	20.00 – 2.45 24.66 – 2.43	Depositor EDS
% Data completeness (in resolution range)	99.5 (20.00-2.45) 98.8 (24.66-2.43)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.65 (at 2.44Å)	Xtriage
Refinement program	REFMAC 5.0	Depositor
R, $R_{free}$	0.206 , 0.269 0.203 , 0.258	Depositor DCC
$R_{free}$ test set	973 reflections (5.45%)	DCC
Wilson B-factor (Å <sup>2</sup> )	31.7	Xtriage
Anisotropy	0.527	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 31.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	4192	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.85% 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

### 5.1 Standard geometry

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

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	0/2019	1.59	24/2722 (0.9%)
1	B	0.60	0/1988	1.71	23/2680 (0.9%)
2	C	0.49	0/73	1.52	2/96 (2.1%)
2	D	0.40	0/76	1.19	1/100 (1.0%)
All	All	0.59	0/4156	1.64	50/5598 (0.9%)

There are no bond length outliers.

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	436	ARG	CD-NE-CZ	37.64	176.30	123.60
1	A	436	ARG	CD-NE-CZ	24.95	158.53	123.60
1	B	515	ARG	NE-CZ-NH2	-13.83	113.39	120.30
1	A	394	ARG	NE-CZ-NH2	11.05	125.83	120.30
1	B	484	ASP	CB-CG-OD1	10.69	127.92	118.30
1	B	352	ARG	NE-CZ-NH2	10.43	125.52	120.30
1	A	352	ARG	NE-CZ-NH1	10.19	125.40	120.30
1	A	501[A]	HIS	CA-CB-CG	10.15	130.85	113.60
1	A	501[B]	HIS	CA-CB-CG	10.15	130.85	113.60
1	B	321	ASP	CB-CG-OD2	-10.11	109.20	118.30
1	B	436	ARG	NE-CZ-NH2	9.71	125.15	120.30
2	C	692	ARG	NE-CZ-NH1	9.18	124.89	120.30
1	A	480	ASP	CB-CG-OD1	9.07	126.46	118.30
1	B	459	TYR	CB-CG-CD1	-8.96	115.62	121.00
1	B	515	ARG	NE-CZ-NH1	8.88	124.74	120.30
1	B	412	ARG	NE-CZ-NH1	7.83	124.22	120.30
1	B	446	VAL	CB-CA-C	7.46	125.58	111.40
1	A	321	ASP	CB-CG-OD1	7.43	124.99	118.30
1	B	411	ASP	CB-CG-OD2	7.11	124.69	118.30
1	A	434	ARG	NE-CZ-NH2	-7.10	116.75	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	480	ASP	CB-CG-OD1	7.09	124.68	118.30
1	A	396	MET	CA-CB-CG	7.06	125.30	113.30
1	B	434	ARG	NE-CZ-NH2	-7.05	116.77	120.30
1	A	484	ASP	CB-CG-OD1	6.88	124.49	118.30
1	B	412	ARG	CD-NE-CZ	6.79	133.10	123.60
1	A	477	ARG	NE-CZ-NH2	6.78	123.69	120.30
2	D	692	ARG	NE-CZ-NH1	6.64	123.62	120.30
1	B	501[A]	HIS	CA-CB-CG	6.44	124.55	113.60
1	B	501[B]	HIS	CA-CB-CG	6.44	124.55	113.60
1	B	459	TYR	CA-CB-CG	-6.28	101.46	113.40
1	B	374	ASP	CB-CG-OD1	6.17	123.86	118.30
1	A	412	ARG	NE-CZ-NH2	-5.96	117.32	120.30
1	B	477	ARG	NE-CZ-NH2	5.95	123.28	120.30
1	A	445	PHE	O-C-N	-5.92	113.23	122.70
1	A	352	ARG	NE-CZ-NH2	-5.89	117.36	120.30
1	A	363[A]	ARG	NE-CZ-NH2	-5.82	117.39	120.30
1	A	363[B]	ARG	NE-CZ-NH2	-5.82	117.39	120.30
1	A	477	ARG	NE-CZ-NH1	-5.70	117.45	120.30
1	B	471	GLU	OE1-CD-OE2	-5.68	116.48	123.30
1	A	515	ARG	NE-CZ-NH2	5.64	123.12	120.30
1	B	512	SER	N-CA-CB	5.54	118.81	110.50
1	A	385	GLU	OE1-CD-OE2	-5.53	116.66	123.30
1	A	426	ASP	CB-CG-OD1	5.51	123.26	118.30
1	B	459	TYR	CB-CG-CD2	5.32	124.19	121.00
1	A	352	ARG	CD-NE-CZ	5.21	130.90	123.60
1	A	422	VAL	CB-CA-C	-5.16	101.60	111.40
1	A	412	ARG	CD-NE-CZ	5.06	130.68	123.60
2	C	692	ARG	CD-NE-CZ	5.05	130.67	123.60
1	A	512	SER	N-CA-CB	5.04	118.05	110.50
1	B	426	ASP	CB-CG-OD1	5.03	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	1952	0	1997	62	0
1	B	1938	0	1991	57	0
2	C	73	0	75	0	0
2	D	76	0	77	2	0
3	A	17	0	10	2	0
3	B	17	0	10	3	0
4	A	59	0	0	0	0
4	B	60	0	0	0	0
All	All	4192	0	4160	97	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 12.

All (97) 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:501[B]:HIS:CE1	1:B:501[B]:HIS:CE1	1.83	1.65
1:A:501[B]:HIS:HE1	1:B:501[B]:HIS:CE1	1.06	1.62
1:A:501[B]:HIS:CE1	1:B:501[B]:HIS:HE1	1.23	1.33
1:A:501[B]:HIS:CE1	1:B:501[B]:HIS:ND1	2.24	1.04
1:A:519:ASN:HD22	1:B:519:ASN:ND2	1.61	0.99
1:A:519:ASN:ND2	1:B:519:ASN:HD22	1.61	0.98
1:A:519:ASN:HD22	1:B:519:ASN:HD22	0.91	0.90
1:B:496:THR:HG22	1:B:499:GLN:H	1.36	0.88
1:A:513[A]:HIS:CD2	1:B:459:TYR:HD2	1.92	0.86
1:A:501[B]:HIS:ND1	1:B:501[B]:HIS:ND1	2.25	0.83
1:A:501[B]:HIS:ND1	1:B:501[B]:HIS:CE1	2.45	0.82
1:B:347:THR:HG21	1:B:534:VAL:HG11	1.65	0.78
1:A:513[A]:HIS:NE2	1:B:459:TYR:HD2	1.83	0.77
1:A:347:THR:HG21	1:A:535:PRO:HD2	1.67	0.76
1:B:338:SER:H	1:B:341:SER:HB3	1.49	0.76
1:A:513[A]:HIS:CE1	1:B:459:TYR:CD2	2.76	0.74
1:A:513[A]:HIS:NE2	1:B:459:TYR:CD2	2.57	0.73
1:A:343:MET:O	1:A:347:THR:HB	1.88	0.73
1:B:525:LEU:HD13	3:B:600:ZTW:H4	1.76	0.68
1:A:496:THR:O	1:A:500:GLN:HG3	1.94	0.68
1:B:343:MET:O	1:B:347:THR:HB	1.94	0.68
1:A:392:VAL:HG13	1:A:432:SER:HA	1.75	0.68
1:A:513[A]:HIS:CE1	1:B:459:TYR:HD2	2.12	0.67
1:A:525:LEU:HD13	3:A:600:ZTW:H4	1.79	0.64
1:A:513[A]:HIS:CD2	1:B:459:TYR:CD2	2.82	0.64
1:A:338:SER:H	1:A:341:SER:HB3	1.66	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:319:LEU:CB	1:B:446:VAL:HG13	2.30	0.60
1:A:501[B]:HIS:CE1	1:B:501[B]:HIS:HD1	2.13	0.60
1:B:504:LEU:O	1:B:508:LEU:HG	2.02	0.60
1:B:525:LEU:HD13	3:B:600:ZTW:C4	2.31	0.59
1:A:315:MET:HE1	1:A:365:PRO:HD2	1.85	0.58
1:A:315:MET:CE	1:A:365:PRO:HG2	2.34	0.58
1:B:396:MET:SD	1:B:435:PHE:HB3	2.44	0.58
1:B:319:LEU:HB2	1:B:446:VAL:HG13	1.87	0.56
1:A:495:LEU:HB3	1:A:499:GLN:HB2	1.89	0.55
1:B:315:MET:CE	1:B:319:LEU:HD21	2.37	0.55
1:A:315:MET:HE2	1:A:319:LEU:HD21	1.89	0.55
1:A:310:LEU:O	1:A:481:LYS:HE3	2.07	0.54
1:B:536:LEU:H	1:B:536:LEU:HD12	1.73	0.54
1:A:504:LEU:O	1:A:508:LEU:HG	2.08	0.54
1:B:392:VAL:HG13	1:B:432:SER:HA	1.89	0.53
1:A:538:ASP:O	1:A:542:GLU:HB2	2.08	0.53
1:A:319:LEU:CB	1:A:446:VAL:HG13	2.39	0.53
1:A:382:ALA:O	1:A:386:ILE:HG12	2.08	0.53
1:B:403:LEU:HD13	1:B:409:LEU:HD13	1.90	0.52
1:A:513[A]:HIS:CE1	1:B:459:TYR:CE2	2.97	0.52
1:A:513[A]:HIS:CG	1:B:459:TYR:HD2	2.26	0.51
1:A:529:LYS:HE3	1:A:541:LEU:CD2	2.41	0.51
1:B:339:GLU:HG2	1:B:533:VAL:HG11	1.93	0.51
1:A:319:LEU:HB3	1:A:446:VAL:HG13	1.94	0.50
1:A:315:MET:HE1	1:A:365:PRO:HG2	1.93	0.50
1:B:316:VAL:HG23	1:B:485:THR:CG2	2.40	0.50
1:B:393:TRP:O	1:B:396:MET:HG3	2.11	0.50
1:B:319:LEU:HB3	1:B:446:VAL:HG13	1.92	0.50
1:A:504:LEU:HD23	1:B:504:LEU:HD23	1.94	0.50
1:A:316:VAL:HG23	1:A:485:THR:CG2	2.42	0.49
1:A:403:LEU:HD13	1:A:409:LEU:HD13	1.93	0.49
1:B:496:THR:HB	1:B:499:GLN:OE1	2.13	0.49
1:A:529:LYS:HE3	1:A:541:LEU:HD21	1.95	0.49
1:A:421:MET:HE3	1:A:424:ILE:HB	1.95	0.49
1:B:347:THR:HG21	1:B:534:VAL:CG1	2.37	0.48
1:A:330:GLU:OE2	1:A:345:LEU:HD23	2.14	0.48
1:A:315:MET:HE1	1:A:365:PRO:CG	2.43	0.48
1:B:330:GLU:OE2	1:B:345:LEU:HD23	2.13	0.48
1:A:346:LEU:HD22	1:A:404:PHE:CD1	2.49	0.48
1:B:496:THR:HG22	1:B:499:GLN:N	2.18	0.47
1:A:455:ASN:O	1:A:458:VAL:HG12	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:463:SER:OG	1:A:468:SER:OG	2.33	0.45
1:A:535:PRO:HB2	1:A:537:TYR:CE2	2.51	0.45
1:B:387:LEU:HB3	3:B:600:ZTW:C12	2.47	0.45
1:A:315:MET:HE1	1:A:365:PRO:CD	2.45	0.45
1:A:433:SER:O	1:A:437:MET:HG3	2.16	0.45
1:B:319:LEU:HB2	1:B:446:VAL:CG1	2.47	0.45
1:A:459:TYR:HD2	1:B:513[A]:HIS:CD2	2.35	0.45
1:A:315:MET:CE	1:A:319:LEU:HD21	2.48	0.44
1:B:374:ASP:OD2	1:B:471:GLU:HG3	2.18	0.44
1:B:315:MET:HE3	1:B:319:LEU:HD21	1.98	0.43
1:B:362:LYS:NZ	2:D:696:ASP:OXT	2.47	0.43
1:B:316:VAL:HG21	1:B:489:LEU:HD21	2.00	0.43
1:A:498:GLN:O	1:A:502:GLN:HG3	2.19	0.43
1:A:347:THR:CG2	1:A:537:TYR:HE2	2.32	0.42
1:A:501[A]:HIS:CE1	1:B:487:ILE:HG21	2.53	0.42
1:A:525:LEU:HD13	3:A:600:ZTW:C4	2.46	0.42
1:B:329:SER:OG	1:B:330:GLU:N	2.52	0.42
1:B:534:VAL:O	1:B:536:LEU:HD12	2.20	0.42
1:B:343:MET:HB3	1:B:534:VAL:HG22	2.00	0.42
1:A:315:MET:HE3	1:A:365:PRO:HG2	2.01	0.42
1:A:319:LEU:HB2	1:A:446:VAL:HG13	2.01	0.42
1:A:479:LEU:HD23	1:A:479:LEU:HA	1.93	0.42
1:B:362:LYS:NZ	2:D:693:LEU:O	2.47	0.41
1:B:338:SER:O	1:B:339:GLU:C	2.58	0.41
1:A:354:LEU:HD21	1:A:383:TRP:CE3	2.56	0.41
1:A:347:THR:HG23	1:A:537:TYR:CE2	2.56	0.40
1:B:374:ASP:OD2	1:B:471:GLU:OE2	2.38	0.40
1:A:434:ARG:NE	1:B:459:TYR:OH	2.54	0.40
1:A:416:LYS:HG2	1:A:422:VAL:HG13	2.03	0.40
1:B:310:LEU:O	1:B:481:LYS:HE3	2.21	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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	245/248 (99%)	243 (99%)	2 (1%)	0	100	100
1	B	241/248 (97%)	234 (97%)	7 (3%)	0	100	100
2	C	7/9 (78%)	7 (100%)	0	0	100	100
2	D	7/9 (78%)	7 (100%)	0	0	100	100
All	All	500/514 (97%)	491 (98%)	9 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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	223/224 (100%)	212 (95%)	11 (5%)	29	39
1	B	219/224 (98%)	209 (95%)	10 (5%)	31	42
2	C	7/9 (78%)	6 (86%)	1 (14%)	4	3
2	D	8/9 (89%)	7 (88%)	1 (12%)	5	4
All	All	457/466 (98%)	434 (95%)	23 (5%)	28	38

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

Mol	Chain	Res	Type
1	A	303	LYS
1	A	330	GLU
1	A	347	THR
1	A	368	VAL
1	A	411	ASP
1	A	422	VAL
1	A	466	LEU
1	A	468	SER
1	A	481	LYS
1	A	541	LEU

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Mol	Chain	Res	Type
1	A	542	GLU
1	B	309	SER
1	B	347	THR
1	B	368	VAL
1	B	411	ASP
1	B	422	VAL
1	B	446	VAL
1	B	466	LEU
1	B	481	LYS
1	B	496	THR
1	B	542	GLU
2	C	690	LEU
2	D	696	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	439	ASN
1	A	519	ASN
1	B	439	ASN
1	B	455	ASN
1	B	516	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 carbohydrates in this entry.

### 5.6 Ligand geometry [i](#)

2 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	ZTW	A	600	-	15,19,19	1.24	1 (6%)	18,27,27	1.40	4 (22%)
3	ZTW	B	600	-	15,19,19	0.99	0	18,27,27	0.99	1 (5%)

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	ZTW	A	600	-	-	0/4/4/4	0/3/3/3
3	ZTW	B	600	-	-	0/4/4/4	0/3/3/3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	600	ZTW	C7-S6	-2.72	1.68	1.72

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	600	ZTW	C10-C9-C8	-2.77	117.08	121.10
3	A	600	ZTW	C13-C8-C7	-2.29	116.44	120.21
3	A	600	ZTW	C4-C5-S6	2.13	128.77	124.84
3	B	600	ZTW	C4-C5-S6	2.17	128.83	124.84
3	A	600	ZTW	C9-C10-C11	3.08	123.37	119.88

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 5 short contacts:



Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	600	ZTW	2	0
3	B	600	ZTW	3	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	243/248 (97%)	-0.09	6 (2%) 58 54	14, 29, 79, 100	0
1	B	242/248 (97%)	0.06	13 (5%) 26 24	14, 30, 91, 108	2 (0%)
2	C	9/9 (100%)	0.24	0 100 100	49, 57, 62, 62	0
2	D	9/9 (100%)	0.53	1 (11%) 6 5	53, 66, 70, 79	0
All	All	503/514 (97%)	0.00	20 (3%) 39 36	14, 30, 85, 108	2 (0%)

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	340	ALA	6.1
1	B	526	TYR	4.5
1	B	337	PHE	4.4
1	B	338	SER	4.3
1	A	330	GLU	4.2
1	B	419	GLU	4.0
1	B	533	VAL	3.8
1	B	330	GLU	3.7
1	B	532	ASN	3.7
1	B	531	LYS	3.1
1	A	337	PHE	3.1
1	B	538	ASP	2.6
1	A	531	LYS	2.5
1	B	462	LEU	2.5
2	D	696	ASP	2.5
1	A	336	PRO	2.4
1	B	545	ASP	2.4
1	A	346	LEU	2.3
1	B	535	PRO	2.1
1	A	534	VAL	2.0

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

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

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	ZTW	B	600	17/17	0.81	0.32	3.61	41,50,55,56	0
3	ZTW	A	600	17/17	0.93	0.22	0.69	27,31,37,38	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.