



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:31 PM GMT

PDB ID : 3ZQI  
Title : Structure of Tetracycline repressor in complex with inducer peptide- TIP2  
Authors : Sevvana, M.; Goeke, D.; Stoeckle, C.; Kaspar, D.; Grubmueller, S.; Goetz, C.;  
Wimmer, C.; Berens, C.; Klotzsche, M.; Muller, Y.A.; Hillen, W.  
Deposited on : 2011-06-09  
Resolution : 1.50 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

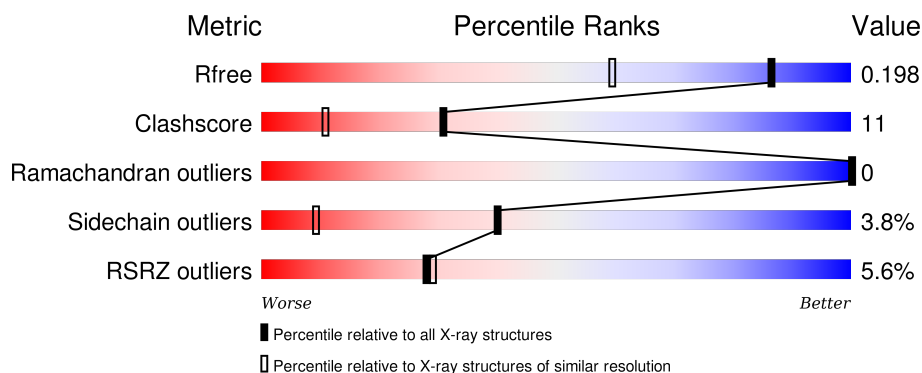
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.50 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	2072 (1.50-1.50)
Clashscore	102246	2274 (1.50-1.50)
Ramachandran outliers	100387	2218 (1.50-1.50)
Sidechain outliers	100360	2216 (1.50-1.50)
RSRZ outliers	91569	2075 (1.50-1.50)

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	208	<div> <div>4%</div> <div>79%</div> <div>16%</div> <div>• •</div> </div>
1	B	208	<div> <div>5%</div> <div>72%</div> <div>19%</div> <div>5%</div> <div>•</div> </div>
2	C	16	<div> <div>6%</div> <div>69%</div> <div>13%</div> <div>6%</div> <div>13%</div> </div>
2	D	16	<div> <div>19%</div> <div>50%</div> <div>38%</div> <div>13%</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
3	EDO	A	1206	-	-	-	X
3	EDO	B	1206	-	-	-	X
3	EDO	B	1207	-	-	-	X
5	PG4	B	1208	-	-	X	X

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 4011 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 TETRACYCLINE REPRESSOR PROTEIN CLASS B FROM TRANSPOSON TN10, TETRACYCLINE REPRESSOR PROTEIN CLASS D.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	201	Total	C	N	O	S	0	7	0
			1633	1036	286	309	2			
1	B	201	Total	C	N	O	S	0	8	0
			1640	1042	287	309	2			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	68	SER	CYS	ENGINEERED MUTATION	UNP P04483
A	88	ASN	CYS	ENGINEERED MUTATION	UNP P04483
A	121	THR	CYS	ENGINEERED MUTATION	UNP P04483
A	144	SER	CYS	ENGINEERED MUTATION	UNP P04483
B	68	SER	CYS	ENGINEERED MUTATION	UNP P04483
B	88	ASN	CYS	ENGINEERED MUTATION	UNP P04483
B	121	THR	CYS	ENGINEERED MUTATION	UNP P04483
B	144	SER	CYS	ENGINEERED MUTATION	UNP P04483

- Molecule 2 is a protein called INDUCER PEPTIDE TIP2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	14	Total	C	N	O	S	0	0	0
			127	84	25	16	2			
2	D	14	Total	C	N	O	S	0	1	0
			132	88	25	16	3			

- Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).

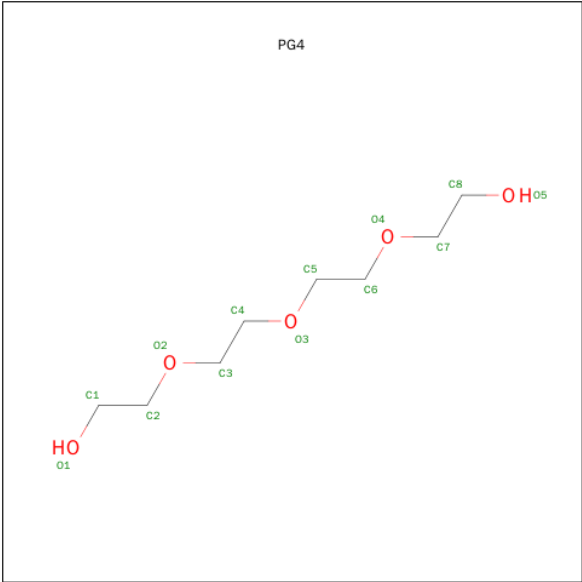


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mg	0	0
			1	1		

- Molecule 5 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: C<sub>8</sub>H<sub>18</sub>O<sub>5</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			13	8	5		

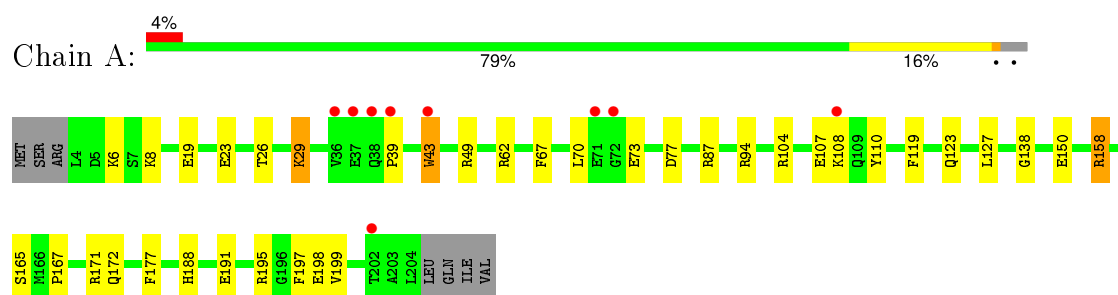
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	225	Total	O	0	0
			225	225		
6	B	208	Total	O	0	0
			208	208		
6	C	6	Total	O	0	0
			6	6		
6	D	6	Total	O	0	0
			6	6		

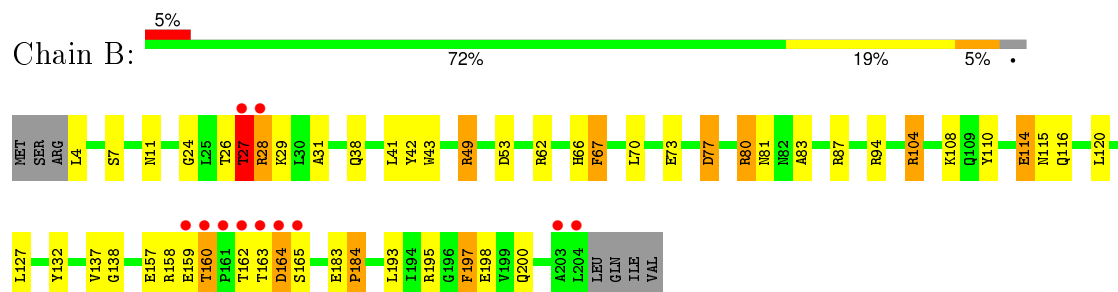
### 3 Residue-property plots [i](#)

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.

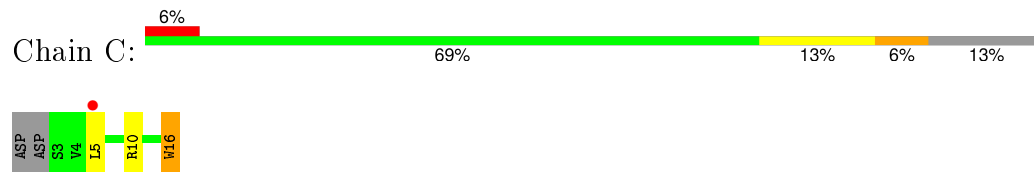
- Molecule 1: TETRACYCLINE REPRESSOR PROTEIN CLASS B FROM TRANSPOSON TN10, TETRACYCLINE REPRESSOR PROTEIN CLASS D



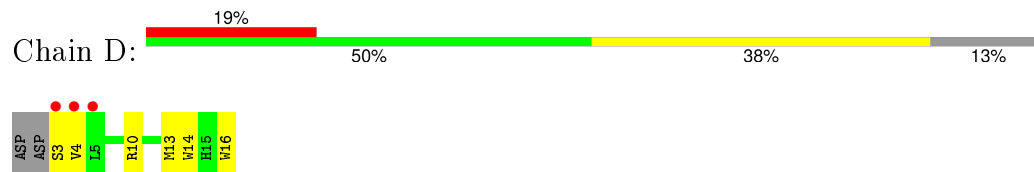
- Molecule 1: TETRACYCLINE REPRESSOR PROTEIN CLASS B FROM TRANSPOSON TN10, TETRACYCLINE REPRESSOR PROTEIN CLASS D



- Molecule 2: INDUCER PEPTIDE TIP2



- Molecule 2: INDUCER PEPTIDE TIP2



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	56.70 Å 78.56 Å 103.88 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.88 – 1.50 24.88 – 1.50	Depositor EDS
% Data completeness (in resolution range)	100.0 (24.88-1.50) 96.0 (24.88-1.50)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	5.05 (at 1.50 Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.160 , 0.195 0.162 , 0.198	Depositor DCC
$R_{free}$ test set	3597 reflections (5.26%)	DCC
Wilson B-factor (Å <sup>2</sup> )	13.6	Xtriage
Anisotropy	0.032	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 47.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 71932 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	4011	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.63% 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: PG4, MG, EDO

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	1.38	13/1686 (0.8%)	1.37	15/2282 (0.7%)
1	B	1.42	11/1696 (0.6%)	1.47	29/2295 (1.3%)
2	C	1.35	1/133 (0.8%)	1.20	1/179 (0.6%)
2	D	1.33	0/141	1.29	0/189
All	All	1.40	25/3656 (0.7%)	1.41	45/4945 (0.9%)

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	157	GLU	CD-OE1	8.72	1.35	1.25
1	B	104	ARG	CG-CD	-7.70	1.32	1.51
1	B	157	GLU	CG-CD	7.70	1.63	1.51
1	B	43	TRP	CB-CG	-6.90	1.37	1.50
2	C	16	TRP	CZ3-CH2	6.66	1.50	1.40
1	B	198	GLU	CB-CG	-6.55	1.39	1.52
1	A	198	GLU	CB-CG	-6.26	1.40	1.52
1	B	27	THR	CB-CG2	-6.25	1.31	1.52
1	A	43	TRP	CB-CG	6.10	1.61	1.50
1	B	73	GLU	CD-OE2	6.00	1.32	1.25
1	B	132	TYR	CD1-CE1	5.99	1.48	1.39
1	A	199	VAL	CB-CG2	-5.98	1.40	1.52
1	A	177	PHE	CD2-CE2	5.98	1.51	1.39
1	A	73	GLU	CG-CD	5.89	1.60	1.51
1	A	150	GLU	CG-CD	5.87	1.60	1.51
1	A	8	LYS	CE-NZ	5.79	1.63	1.49
1	A	19	GLU	CG-CD	5.78	1.60	1.51
1	A	104	ARG	CZ-NH1	5.77	1.40	1.33
1	A	165	SER	CB-OG	5.75	1.49	1.42
1	A	107	GLU	CD-OE1	5.64	1.31	1.25
1	A	104	ARG	CZ-NH2	5.60	1.40	1.33
1	A	150	GLU	CD-OE1	5.48	1.31	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	114	GLU	CG-CD	5.41	1.60	1.51
1	B	42	TYR	CD1-CE1	5.39	1.47	1.39
1	B	73	GLU	CB-CG	-5.35	1.42	1.52

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	62	ARG	NE-CZ-NH2	-13.91	113.35	120.30
1	A	49	ARG	NE-CZ-NH2	-12.56	114.02	120.30
1	A	158	ARG	NE-CZ-NH2	-11.32	114.64	120.30
1	A	94	ARG	NE-CZ-NH2	-10.90	114.85	120.30
1	B	49[A]	ARG	NE-CZ-NH2	-10.64	114.98	120.30
1	B	49[B]	ARG	NE-CZ-NH2	-10.64	114.98	120.30
1	A	87	ARG	NE-CZ-NH2	-10.48	115.06	120.30
1	B	158	ARG	NE-CZ-NH1	10.21	125.40	120.30
1	B	49[A]	ARG	NE-CZ-NH1	10.15	125.38	120.30
1	B	49[B]	ARG	NE-CZ-NH1	10.15	125.38	120.30
1	B	87	ARG	NE-CZ-NH1	10.08	125.34	120.30
1	B	87	ARG	NE-CZ-NH2	-10.03	115.28	120.30
1	A	49	ARG	NE-CZ-NH1	9.71	125.16	120.30
1	B	94	ARG	NE-CZ-NH2	-9.18	115.71	120.30
1	B	80	ARG	NE-CZ-NH2	-8.74	115.93	120.30
1	A	87	ARG	NE-CZ-NH1	8.18	124.39	120.30
1	A	197	PHE	CB-CG-CD2	-7.93	115.25	120.80
1	B	62	ARG	NE-CZ-NH2	-7.62	116.49	120.30
1	B	62	ARG	NE-CZ-NH1	6.95	123.78	120.30
1	B	49[A]	ARG	CD-NE-CZ	6.79	133.11	123.60
1	B	49[B]	ARG	CD-NE-CZ	6.79	133.11	123.60
1	B	158	ARG	NE-CZ-NH2	-6.60	117.00	120.30
1	A	77	ASP	CB-CG-OD1	6.52	124.17	118.30
1	A	70	LEU	CB-CG-CD2	6.47	121.99	111.00
1	B	27	THR	CB-CA-C	-6.32	94.54	111.60
1	B	67	PHE	CB-CG-CD2	-6.17	116.48	120.80
1	A	158	ARG	CB-CA-C	6.14	122.68	110.40
1	B	110	TYR	CB-CG-CD2	-6.14	117.32	121.00
1	B	80	ARG	NE-CZ-NH1	5.94	123.27	120.30
1	A	110	TYR	CB-CG-CD2	-5.91	117.45	121.00
1	B	193	LEU	CB-CG-CD1	-5.83	101.09	111.00
1	B	132	TYR	CB-CG-CD2	-5.59	117.64	121.00
1	B	157	GLU	CG-CD-OE1	5.54	129.38	118.30
1	B	77[A]	ASP	CB-CG-OD2	-5.49	113.36	118.30
1	B	77[B]	ASP	CB-CG-OD2	-5.49	113.36	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	70	LEU	CA-CB-CG	-5.48	102.70	115.30
1	A	119	PHE	CB-CG-CD2	-5.46	116.98	120.80
1	B	197	PHE	CB-CG-CD1	5.33	124.53	120.80
1	A	158	ARG	N-CA-CB	-5.29	101.08	110.60
1	B	120	LEU	CB-CG-CD1	-5.23	102.10	111.00
1	A	62	ARG	NE-CZ-NH1	5.23	122.92	120.30
1	B	110	TYR	CB-CG-CD1	5.10	124.06	121.00
1	B	70	LEU	CB-CG-CD2	5.06	119.59	111.00
1	B	67	PHE	CB-CG-CD1	5.02	124.32	120.80
2	C	10	ARG	NE-CZ-NH1	5.00	122.80	120.30

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	1633	0	1640	17	0
1	B	1640	0	1652	56	0
2	C	127	0	120	1	0
2	D	132	0	129	7	0
3	A	8	0	12	0	0
3	B	12	0	18	0	0
4	A	1	0	0	0	0
5	B	13	0	18	9	0
6	A	225	0	0	9	0
6	B	208	0	0	19	0
6	C	6	0	0	1	0
6	D	6	0	0	1	0
All	All	4011	0	3589	78	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 11.

All (78) 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:4:LEU:N	6:B:2001:HOH:O	1.65	1.24
1:B:115:ASN:HB3	6:B:2130:HOH:O	1.34	1.19
1:A:26:THR:OG1	1:A:29:LYS:HD2	1.40	1.16
1:B:28:ARG:HD3	1:B:28:ARG:N	1.54	1.16
1:A:171:ARG:HG2	6:A:2188:HOH:O	0.96	1.11
1:B:77[A]:ASP:OD1	6:B:2094:HOH:O	1.68	1.10
5:B:1208:PG4:H52	6:B:2062:HOH:O	1.56	1.06
1:B:28:ARG:CD	1:B:28:ARG:H	1.72	1.01
1:B:195:ARG:HD3	6:B:2196:HOH:O	1.66	0.95
1:B:28:ARG:H	1:B:28:ARG:HD3	0.78	0.93
6:A:2137:HOH:O	1:B:164:ASP:HB2	1.67	0.92
1:A:123:GLN:NE2	6:A:2100:HOH:O	1.81	0.91
1:B:26:THR:HG23	1:B:29:LYS:H	1.34	0.91
1:B:116:GLN:HE22	2:D:14:TRP:HE1	1.20	0.84
5:B:1208:PG4:C5	6:B:2206:HOH:O	2.29	0.80
1:A:23:GLU:OE2	6:A:2034:HOH:O	2.03	0.77
1:A:108:LYS:HD2	6:A:2135:HOH:O	1.85	0.76
1:B:164:ASP:OD1	1:B:165:SER:N	2.19	0.76
1:B:49[A]:ARG:HD2	1:B:53:ASP:OD1	1.86	0.75
1:B:26:THR:OG1	1:B:28:ARG:NE	2.23	0.71
1:A:26:THR:OG1	1:A:29:LYS:CD	2.31	0.70
1:B:115:ASN:HB3	6:B:2131:HOH:O	1.92	0.69
5:B:1208:PG4:C5	6:B:2205:HOH:O	2.41	0.68
1:B:28:ARG:CD	1:B:28:ARG:N	2.44	0.67
1:B:27:THR:HG23	1:B:28:ARG:CD	2.26	0.66
1:A:191:GLU:OE2	1:A:195[B]:ARG:NH2	2.30	0.65
1:B:115:ASN:CG	6:B:2131:HOH:O	2.36	0.64
1:B:115:ASN:CB	6:B:2131:HOH:O	2.45	0.62
1:B:83:ALA:HB2	1:B:137[A]:VAL:CG1	2.32	0.59
1:B:26:THR:HG23	1:B:29:LYS:N	2.13	0.59
1:B:27:THR:OG1	1:B:28:ARG:N	2.31	0.57
1:B:116:GLN:NE2	2:D:14:TRP:HE1	1.97	0.57
1:B:108:LYS:NZ	6:B:2123:HOH:O	1.79	0.56
1:A:138:GLY:HA3	2:C:16:TRP:CE2	2.41	0.56
1:B:66:HIS:HD2	6:B:2080:HOH:O	1.87	0.56
1:A:127[A]:LEU:HD23	6:B:2174:HOH:O	2.05	0.55
1:B:49[A]:ARG:HD2	1:B:53:ASP:CG	2.27	0.55
1:A:39:PRO:O	1:A:43:TRP:CD1	2.60	0.54
1:B:4:LEU:CA	6:B:2001:HOH:O	2.35	0.54
1:B:26:THR:CG2	1:B:29:LYS:H	2.11	0.54
1:A:171:ARG:CG	6:A:2188:HOH:O	1.83	0.54
1:B:26:THR:HG22	1:B:29:LYS:HD3	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:138:GLY:HA3	2:D:16:TRP:CE2	2.43	0.54
1:A:171:ARG:CD	6:A:2188:HOH:O	2.42	0.52
1:B:11:ASN:OD1	5:B:1208:PG4:H21	2.09	0.52
1:B:104:ARG:CZ	2:D:13[B]:MET:HE1	2.40	0.52
1:A:188:HIS:CE1	1:B:200[A]:GLN:OE1	2.65	0.50
1:B:7[A]:SER:OG	5:B:1208:PG4:H22	2.12	0.50
1:B:77[A]:ASP:OD1	1:B:80:ARG:NH1	2.44	0.49
5:B:1208:PG4:C5	6:B:2062:HOH:O	2.34	0.49
1:A:158:ARG:HG3	6:B:2057:HOH:O	2.13	0.48
1:A:172:GLN:CD	1:B:127[A]:LEU:CD2	2.81	0.48
1:B:24:GLY:O	1:B:29:LYS:HE2	2.15	0.46
1:A:171:ARG:NE	6:A:2188:HOH:O	2.49	0.45
1:B:27:THR:HG23	1:B:28:ARG:HD3	1.97	0.45
5:B:1208:PG4:H51	6:B:2206:HOH:O	2.07	0.44
1:B:27:THR:HG23	1:B:28:ARG:NE	2.32	0.44
5:B:1208:PG4:H52	6:B:2206:HOH:O	2.05	0.44
1:B:104:ARG:CZ	2:D:13[B]:MET:CE	2.96	0.43
1:B:28:ARG:HB3	1:B:38:GLN:CD	2.39	0.43
1:B:183:GLU:HB3	1:B:184:PRO:HD3	1.99	0.43
1:A:167:PRO:HB2	1:B:114:GLU:CD	2.39	0.43
1:B:28:ARG:HB3	1:B:38:GLN:OE1	2.19	0.43
1:B:197:PHE:HA	1:B:200[B]:GLN:HE21	1.85	0.42
1:B:160:THR:HB	6:C:2003:HOH:O	2.19	0.42
1:B:49[A]:ARG:NH2	6:B:2053:HOH:O	2.52	0.41
1:B:26:THR:CG2	1:B:29:LYS:HD3	2.51	0.41
6:A:2167:HOH:O	2:D:10:ARG:HG2	2.20	0.41
1:B:11:ASN:OD1	5:B:1208:PG4:C2	2.69	0.41
1:B:66:HIS:CG	1:B:81:ASN:HB3	2.56	0.40
1:B:31:ALA:HB2	1:B:41[B]:LEU:HD12	2.03	0.40
2:D:4:VAL:HG13	6:D:2001:HOH:O	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	206/208 (99%)	205 (100%)	1 (0%)	0	100	100
1	B	207/208 (100%)	205 (99%)	2 (1%)	0	100	100
2	C	12/16 (75%)	11 (92%)	1 (8%)	0	100	100
2	D	13/16 (81%)	12 (92%)	1 (8%)	0	100	100
All	All	438/448 (98%)	433 (99%)	5 (1%)	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	179/179 (100%)	176 (98%)	3 (2%)	68	37
1	B	180/179 (101%)	171 (95%)	9 (5%)	30	5
2	C	11/13 (85%)	10 (91%)	1 (9%)	12	1
2	D	12/13 (92%)	11 (92%)	1 (8%)	14	1
All	All	382/384 (100%)	368 (96%)	14 (4%)	40	10

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	LYS
1	A	29	LYS
1	A	67	PHE
1	B	27	THR
1	B	28	ARG
1	B	67	PHE
1	B	159	GLU
1	B	160	THR
1	B	162	THR
1	B	163	THR

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Mol	Chain	Res	Type
1	B	164	ASP
1	B	184	PRO
2	C	5	LEU
2	D	3	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	38	GLN
1	A	44	HIS
1	A	115	ASN
1	B	66	HIS
1	B	116	GLN
1	B	152	GLN
1	B	179	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](#)

Of 7 ligands modelled in this entry, 1 is monoatomic - leaving 6 for Mogul analysis.

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	EDO	A	1205	-	3,3,3	1.18	0	2,2,2	1.06	0
3	EDO	A	1206	-	3,3,3	0.43	0	2,2,2	1.35	0
3	EDO	B	1205	-	3,3,3	0.44	0	2,2,2	0.85	0
3	EDO	B	1206	-	3,3,3	0.54	0	2,2,2	0.74	0
3	EDO	B	1207	-	3,3,3	1.05	0	2,2,2	0.76	0
5	PG4	B	1208	-	12,12,12	1.07	1 (8%)	11,11,11	1.23	1 (9%)

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	EDO	A	1205	-	-	0/1/1/1	0/0/0/0
3	EDO	A	1206	-	-	0/1/1/1	0/0/0/0
3	EDO	B	1205	-	-	0/1/1/1	0/0/0/0
3	EDO	B	1206	-	-	0/1/1/1	0/0/0/0
3	EDO	B	1207	-	-	0/1/1/1	0/0/0/0
5	PG4	B	1208	-	-	0/10/10/10	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	1208	PG4	C6-C5	2.11	1.59	1.48

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	1208	PG4	O2-C2-C1	2.02	119.74	110.43

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	1208	PG4	9	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	201/208 (96%)	0.01	9 (4%) 37 38	7, 15, 33, 41	0
1	B	201/208 (96%)	-0.02	11 (5%) 29 30	7, 14, 34, 52	0
2	C	14/16 (87%)	0.55	1 (7%) 19 19	11, 24, 43, 43	0
2	D	14/16 (87%)	1.13	3 (21%) 1 1	9, 20, 34, 42	0
All	All	430/448 (95%)	0.05	24 (5%) 28 29	7, 15, 35, 52	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	162	THR	8.0
1	B	163	THR	7.9
1	A	43	TRP	5.0
1	B	164	ASP	4.5
2	D	3	SER	4.0
1	B	160	THR	3.6
1	B	28	ARG	3.6
2	C	5	LEU	3.5
2	D	5	LEU	3.5
1	A	71	GLU	3.3
1	B	165	SER	3.3
2	D	4	VAL	3.0
1	B	204	LEU	2.8
1	A	202	THR	2.7
1	A	72	GLY	2.7
1	A	108	LYS	2.5
1	B	159	GLU	2.4
1	A	39	PRO	2.4
1	B	161	PRO	2.4
1	A	36	VAL	2.2
1	A	37	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	38	GLN	2.1
1	B	27	THR	2.1
1	B	203	ALA	2.1

## 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(Å <sup>2</sup> )	Q<0.9
5	PG4	B	1208	13/13	0.76	0.18	7.41	22,33,39,39	0
3	EDO	B	1207	4/4	0.82	0.15	4.36	20,22,28,32	0
3	EDO	B	1206	4/4	0.92	0.17	3.72	21,30,35,38	0
3	EDO	A	1206	4/4	0.92	0.12	3.06	27,31,36,37	0
3	EDO	A	1205	4/4	0.96	0.06	-0.67	15,17,20,24	0
3	EDO	B	1205	4/4	0.96	0.07	-	17,22,26,27	0
4	MG	A	1207	1/1	0.99	0.20	-	23,23,23,23	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.