



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 02:38 pm GMT

PDB ID : 3KDZ  
Title : X-ray crystal structure of a tyrosine aminomutase mutant construct with bound ligand  
Authors : Cooke, H.A.; Bruner, S.D.  
Deposited on : 2009-10-23  
Resolution : 2.20 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

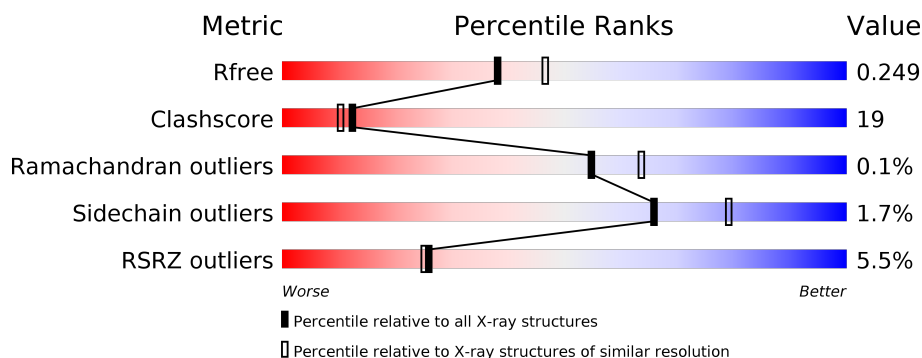
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.20 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	4002 (2.20-2.20)
Clashscore	112137	4730 (2.20-2.20)
Ramachandran outliers	110173	4656 (2.20-2.20)
Sidechain outliers	110143	4657 (2.20-2.20)
RSRZ outliers	101464	4033 (2.20-2.20)

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	537	<div> <div>7%</div> <div>68%</div> <div>29%</div> <div>..</div> </div>
1	B	537	<div> <div>4%</div> <div>72%</div> <div>24%</div> <div>...</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	TYR	A	600	-	-	X	X
2	TYR	B	600	-	-	-	X

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8266 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.

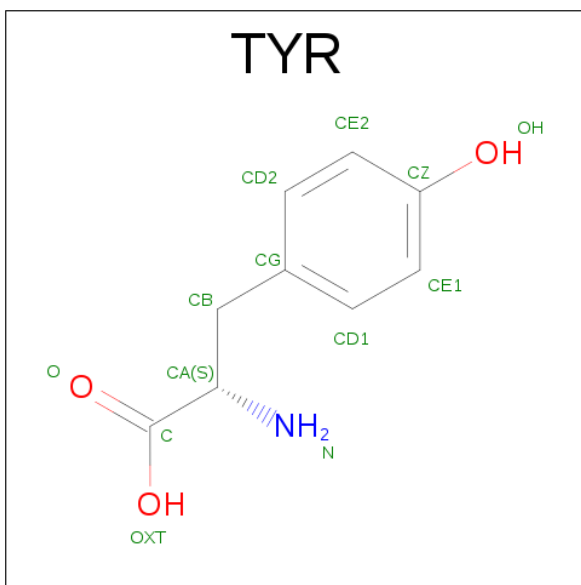
- Molecule 1 is a protein called Histidine ammonia-lyase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	527	Total	C	N	O	S	0	0	0
			4012	2508	728	768	8			
1	B	527	Total	C	N	O	S	0	0	0
			4012	2508	728	768	8			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	63	PHE	TYR	ENGINEERED	UNP Q8GMG0
A	152	MDO	ALA	CHROMOPHORE	UNP Q8GMG0
A	152	MDO	SER	CHROMOPHORE	UNP Q8GMG0
A	152	MDO	GLY	CHROMOPHORE	UNP Q8GMG0
B	63	PHE	TYR	ENGINEERED	UNP Q8GMG0
B	152	MDO	ALA	CHROMOPHORE	UNP Q8GMG0
B	152	MDO	SER	CHROMOPHORE	UNP Q8GMG0
B	152	MDO	GLY	CHROMOPHORE	UNP Q8GMG0

- Molecule 2 is TYROSINE (three-letter code: TYR) (formula: C<sub>9</sub>H<sub>11</sub>NO<sub>3</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			13	9	1	3		
2	B	1	Total	C	N	O	0	0
			13	9	1	3		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	94	Total	O	0	0
			94	94		
3	B	122	Total	O	0	0
			122	122		





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	92.72Å 145.94Å 74.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.40 – 2.20 39.41 – 2.20	Depositor EDS
% Data completeness (in resolution range)	95.9 (39.40-2.20) 95.8 (39.41-2.20)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.69 (at 2.20Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.205 , 0.249 0.203 , 0.249	Depositor DCC
$R_{free}$ test set	5052 reflections (10.07%)	DCC
Wilson B-factor (Å <sup>2</sup> )	34.9	Xtriage
Anisotropy	0.659	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 40.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	8266	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.50% 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: MDO

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	0.35	1/4058 (0.0%)	0.61	1/5499 (0.0%)
1	B	0.51	8/4058 (0.2%)	0.69	6/5499 (0.1%)
All	All	0.43	9/8116 (0.1%)	0.65	7/10998 (0.1%)

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	1
1	B	0	3
All	All	0	4

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	294	GLY	C-O	-8.15	1.10	1.23
1	B	297	VAL	CB-CG1	-7.86	1.36	1.52
1	B	300	SER	C-O	-7.09	1.09	1.23
1	B	300	SER	CB-OG	-6.58	1.33	1.42
1	B	293	ALA	CA-CB	-6.49	1.38	1.52
1	B	300	SER	CA-CB	-6.13	1.43	1.52
1	B	297	VAL	CB-CG2	-6.09	1.40	1.52
1	A	342	PRO	N-CD	5.65	1.55	1.47
1	B	295	LYS	CA-C	-5.61	1.38	1.52

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	306	LYS	CB-CA-C	-11.39	87.62	110.40
1	B	306	LYS	N-CA-C	9.16	135.73	111.00
1	B	304	LEU	N-CA-C	-7.44	90.91	111.00
1	B	298	GLN	N-CA-C	-6.39	93.76	111.00
1	B	310	LEU	CB-CA-C	-6.02	98.76	110.20
1	B	300	SER	CB-CA-C	-5.80	99.08	110.10
1	A	341	ASN	CB-CA-C	5.43	121.26	110.40

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	152	MDO	Mainchain
1	B	152	MDO	Mainchain,Peptide
1	B	294	GLY	Mainchain

## 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	4012	0	4043	192	0
1	B	4012	0	4041	165	1
2	A	13	0	8	7	0
2	B	13	0	8	4	0
3	A	94	0	0	5	0
3	B	122	0	0	5	0
All	All	8266	0	8100	304	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 19.

All (304) 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:152:MDO:CB2	2:A:600:TYR:N	1.87	1.36
1:A:295:LYS:HB2	3:A:679:HOH:O	1.37	1.18

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:295:LYS:HE2	1:B:297:VAL:N	1.62	1.14
1:A:152:MDO:HB22	2:A:600:TYR:N	1.50	1.14
1:A:73:ILE:CD1	1:B:304:LEU:HB3	1.82	1.09
1:B:299:ARG:N	1:B:299:ARG:HD2	1.66	1.06
1:A:350:ILE:HD12	1:A:350:ILE:H	1.22	1.03
1:A:73:ILE:HD12	1:B:306:LYS:HZ3	1.21	1.02
1:B:299:ARG:HD2	1:B:299:ARG:H	1.18	1.02
1:A:152:MDO:HB21	1:B:308:TYR:OH	1.60	1.02
1:A:73:ILE:HD11	1:B:304:LEU:HB3	1.42	1.00
1:B:295:LYS:HE2	1:B:297:VAL:H	0.82	0.97
1:B:35:VAL:H	1:B:137:ASN:HD21	1.07	0.96
1:B:306:LYS:O	1:B:311:ARG:NH2	1.98	0.95
1:B:235:LEU:HD11	1:B:386:VAL:HG11	1.47	0.94
1:B:299:ARG:HH11	1:B:299:ARG:HG2	1.33	0.93
1:A:63:PHE:HZ	2:A:600:TYR:HB3	1.34	0.92
1:B:299:ARG:H	1:B:299:ARG:CD	1.81	0.90
1:A:35:VAL:H	1:A:137:ASN:HD21	1.17	0.90
1:B:295:LYS:CE	1:B:297:VAL:H	1.79	0.89
1:B:447:MET:HG3	1:B:450:ILE:HD12	1.52	0.89
1:A:53:GLU:O	1:A:57:GLU:HG2	1.73	0.88
1:A:251:HIS:HD2	1:A:260:GLN:HG2	1.36	0.88
1:A:284:ARG:HD2	1:B:60:ILE:HD12	1.56	0.86
1:A:73:ILE:HD13	1:B:304:LEU:HB3	1.56	0.86
1:A:73:ILE:HD12	1:B:306:LYS:NZ	1.91	0.84
1:A:63:PHE:CZ	2:A:600:TYR:HB3	2.13	0.83
1:A:73:ILE:CD1	1:B:304:LEU:CB	2.55	0.82
1:A:75:MET:HE1	1:B:299:ARG:HA	1.62	0.80
1:B:295:LYS:HE3	1:B:296:ASP:H	1.45	0.80
1:A:251:HIS:CD2	1:A:260:GLN:HG2	2.18	0.78
1:A:64:GLY:HA3	1:A:201:LEU:HD22	1.65	0.78
1:B:247:LEU:HB3	1:B:249:GLU:OE2	1.83	0.77
1:A:344:PHE:HE1	1:A:350:ILE:HG13	1.50	0.76
1:A:73:ILE:HD11	1:B:304:LEU:CB	2.13	0.75
1:A:96:GLY:H	1:A:161:HIS:HE1	1.32	0.75
1:A:350:ILE:H	1:A:350:ILE:CD1	1.98	0.74
1:A:498:ARG:HE	1:A:498:ARG:HA	1.52	0.73
1:A:350:ILE:HD12	1:A:350:ILE:N	2.03	0.72
1:A:75:MET:HE1	1:B:299:ARG:CA	2.18	0.72
1:A:60:ILE:HG22	1:B:288:GLN:NE2	2.06	0.71
1:A:407:HIS:HD2	1:A:507:ASP:H	1.39	0.70
1:B:235:LEU:HD11	1:B:386:VAL:CG1	2.19	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:198:LYS:HG2	1:A:198:LYS:O	1.91	0.70
1:A:73:ILE:CD1	1:B:306:LYS:HZ3	2.00	0.69
1:B:392:SER:HB3	1:B:395:LEU:HB2	1.76	0.67
1:A:61:PRO:HG2	1:B:287:LEU:HB3	1.76	0.67
1:A:135:TYR:HA	1:A:140:ILE:HG23	1.75	0.67
1:A:350:ILE:HD13	1:B:280:HIS:HE1	1.60	0.67
1:A:75:MET:HE1	1:B:299:ARG:CB	2.24	0.67
1:A:207:THR:H	1:A:339:ASN:HD22	1.41	0.67
1:B:295:LYS:HE3	1:B:296:ASP:N	2.11	0.66
1:B:15:VAL:HG13	1:B:115:ASN:HD22	1.60	0.66
1:A:402:GLY:N	3:A:662:HOH:O	2.14	0.66
1:A:172:VAL:HG21	1:A:184:VAL:HG21	1.78	0.65
1:B:295:LYS:CD	1:B:297:VAL:O	2.44	0.65
1:A:75:MET:CE	1:B:299:ARG:HB3	2.25	0.65
1:A:400:VAL:HG21	1:A:404:PRO:HB3	1.77	0.65
1:A:508:ARG:NH1	1:A:513:ASP:OD2	2.30	0.65
2:B:600:TYR:O	3:B:692:HOH:O	2.14	0.65
1:B:299:ARG:NH1	1:B:299:ARG:HG2	2.10	0.64
1:B:82:GLU:HG3	1:B:197:PHE:CD1	2.33	0.64
1:B:184:VAL:O	1:B:188:ARG:HG2	1.98	0.63
1:A:123:ALA:HB2	1:A:344:PHE:CD2	2.34	0.63
1:A:135:TYR:HA	1:A:140:ILE:CG2	2.28	0.63
1:A:330:LYS:HE2	1:A:330:LYS:HA	1.79	0.63
1:B:249:GLU:H	1:B:249:GLU:CD	2.00	0.63
1:B:301:GLU:HG3	1:B:302:ILE:HD12	1.80	0.62
1:B:298:GLN:C	1:B:299:ARG:HD2	2.19	0.62
1:A:35:VAL:HG12	1:A:137:ASN:ND2	2.14	0.62
1:B:309:SER:O	1:B:313:ILE:HD11	1.99	0.62
1:A:15:VAL:HG23	1:A:35:VAL:HG23	1.82	0.61
1:B:330:LYS:HE2	1:B:330:LYS:HA	1.80	0.61
1:B:298:GLN:CA	1:B:299:ARG:HD2	2.30	0.61
1:B:392:SER:HB2	3:B:691:HOH:O	2.01	0.61
1:A:69:TYR:HB3	2:A:600:TYR:CE2	2.35	0.61
1:A:232:THR:HG21	1:A:313:ILE:HG12	1.83	0.61
1:A:172:VAL:CG2	1:A:184:VAL:HG21	2.30	0.60
1:A:75:MET:HE3	1:B:299:ARG:HB3	1.82	0.60
1:A:73:ILE:HD13	1:B:304:LEU:CB	2.24	0.60
1:B:35:VAL:N	1:B:137:ASN:HD21	1.90	0.60
1:B:231:VAL:HB	1:B:470:TYR:CE1	2.37	0.60
1:A:475:GLN:NE2	1:A:504:LEU:HB3	2.17	0.60
1:A:344:PHE:CE1	1:A:350:ILE:HG13	2.35	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:311:ARG:NH2	2:B:600:TYR:O	2.35	0.60
1:B:407:HIS:CD2	1:B:507:ASP:H	2.20	0.59
1:B:401:SER:HB2	3:B:609:HOH:O	2.03	0.59
1:B:69:TYR:HB3	2:B:600:TYR:CE2	2.38	0.59
1:B:251:HIS:HD2	1:B:260:GLN:HE21	1.49	0.59
1:A:355:ASN:ND2	1:B:315:GLN:HE21	2.00	0.59
1:A:73:ILE:HB	1:B:306:LYS:HD2	1.84	0.59
1:B:407:HIS:HD2	1:B:507:ASP:H	1.49	0.59
1:B:207:THR:H	1:B:339:ASN:HD22	1.51	0.59
1:A:123:ALA:HB2	1:A:344:PHE:HD2	1.67	0.59
1:A:305:GLN:HE22	1:B:341:ASN:HD21	1.50	0.59
1:A:308:TYR:CZ	1:B:442:GLN:HG2	2.37	0.58
1:B:284:ARG:HH22	1:B:285:ARG:NH2	2.02	0.58
1:A:251:HIS:CD2	1:A:260:GLN:HE21	2.21	0.58
1:B:17:GLY:H	1:B:115:ASN:ND2	2.01	0.58
1:B:299:ARG:N	1:B:299:ARG:CD	2.30	0.58
1:A:82:GLU:OE2	1:A:196:ARG:HB3	2.04	0.58
1:A:70:GLY:O	1:B:306:LYS:HE2	2.03	0.57
1:A:62:ILE:HB	1:A:65:VAL:HG13	1.87	0.57
1:B:71:GLU:H	1:B:438:ASN:HD21	1.50	0.57
1:A:402:GLY:CA	3:A:662:HOH:O	2.53	0.57
1:A:230:ILE:HD13	1:A:530:ILE:HD13	1.86	0.57
1:A:55:ILE:HG22	1:A:60:ILE:HD11	1.85	0.57
1:A:355:ASN:HD22	1:B:315:GLN:HE21	1.52	0.56
1:A:59:ASN:OD1	1:A:79:LYS:HG3	2.05	0.56
1:A:65:VAL:HG23	1:A:197:PHE:CD1	2.39	0.56
1:A:251:HIS:HD2	1:A:260:GLN:CG	2.14	0.56
1:A:60:ILE:C	1:A:60:ILE:HD12	2.26	0.56
1:A:15:VAL:CG2	1:A:35:VAL:HG23	2.36	0.56
1:B:295:LYS:HD2	1:B:297:VAL:O	2.04	0.56
1:A:310:LEU:HA	1:A:313:ILE:HD11	1.88	0.55
1:A:38:PRO:HB2	1:A:40:GLU:HG2	1.88	0.55
1:A:346:GLU:HG3	1:A:346:GLU:O	2.05	0.55
1:A:311:ARG:HH21	1:B:356:PHE:HE2	1.55	0.55
1:A:402:GLY:HA3	3:A:662:HOH:O	2.07	0.55
1:B:298:GLN:HA	1:B:299:ARG:HD2	1.89	0.55
1:A:74:TYR:HD2	1:A:75:MET:HE2	1.72	0.55
1:B:520:ALA:O	1:B:525:GLU:HG2	2.07	0.55
1:A:523:ARG:NE	1:A:525:GLU:OE1	2.37	0.54
1:A:239:VAL:HG22	1:A:239:VAL:O	2.06	0.54
1:A:359:GLN:N	1:A:360:PRO:HD2	2.22	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:307:ALA:HB2	1:B:385:ARG:NH2	2.23	0.54
1:B:280:HIS:CE1	1:B:284:ARG:HD3	2.43	0.54
1:A:135:TYR:CA	1:A:140:ILE:HG23	2.38	0.53
1:A:46:GLN:O	1:A:49:ARG:HB3	2.08	0.53
1:A:152:MDO:CB2	2:A:600:TYR:CA	2.82	0.53
1:A:389:ARG:HG2	3:A:674:HOH:O	2.08	0.53
1:A:315:GLN:HE21	1:B:355:ASN:ND2	2.07	0.53
1:A:313:ILE:H	1:A:313:ILE:HD12	1.73	0.53
1:A:392:SER:O	1:A:393:TYR:HB2	2.09	0.52
1:A:447:MET:HE2	1:A:450:ILE:CD1	2.39	0.52
1:A:284:ARG:HG3	1:B:61:PRO:HG2	1.91	0.52
1:A:308:TYR:OH	1:B:152:MDO:CB2	2.57	0.52
1:B:392:SER:O	1:B:393:TYR:HB2	2.10	0.52
1:A:73:ILE:CG2	1:B:306:LYS:HD2	2.39	0.52
1:B:359:GLN:N	1:B:360:PRO:HD2	2.24	0.52
1:A:207:THR:H	1:A:339:ASN:ND2	2.06	0.52
1:B:113:ARG:NH1	1:B:206:GLY:HA3	2.25	0.52
1:B:15:VAL:CG1	1:B:115:ASN:HD22	2.21	0.52
1:B:313:ILE:HB	1:B:314:PRO:HD3	1.91	0.52
1:A:300:SER:O	1:B:74:TYR:HB2	2.10	0.51
1:A:351:PHE:CZ	1:B:250:GLY:HA3	2.45	0.51
1:A:73:ILE:HD13	1:B:304:LEU:O	2.11	0.51
1:A:364:ALA:O	1:A:368:VAL:HG23	2.11	0.51
1:B:11:VAL:CG1	1:B:12:PRO:HD3	2.41	0.51
1:A:60:ILE:HD12	1:A:60:ILE:O	2.10	0.51
1:A:350:ILE:HD13	1:B:280:HIS:CE1	2.43	0.51
1:B:251:HIS:CD2	1:B:260:GLN:HE21	2.27	0.51
1:A:509:TYR:CE2	1:A:511:ALA:HB3	2.45	0.51
1:B:124:VAL:CG1	1:B:128:ILE:HB	2.41	0.51
1:A:82:GLU:HG3	1:A:197:PHE:CG	2.46	0.50
1:A:71:GLU:OE2	1:B:303:TYR:CE1	2.64	0.50
1:B:302:ILE:HD12	1:B:302:ILE:N	2.27	0.50
1:A:71:GLU:OE2	1:B:303:TYR:HE1	1.93	0.50
1:B:82:GLU:OE2	1:B:196:ARG:HB3	2.11	0.50
1:A:235:LEU:O	1:A:235:LEU:HD23	2.11	0.50
1:A:296:ASP:HA	1:B:78:ASP:HB2	1.93	0.50
1:A:69:TYR:CE1	1:A:89:LEU:HD22	2.46	0.49
1:A:96:GLY:H	1:A:161:HIS:CE1	2.21	0.49
1:B:159:LEU:HB3	1:B:204:ILE:HA	1.94	0.49
1:B:96:GLY:H	1:B:161:HIS:HE1	1.58	0.49
1:A:244:SER:HB3	1:A:245:PRO:HD3	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:342:PRO:HB2	1:B:350:ILE:HG22	1.95	0.49
1:A:346:GLU:O	1:A:348:LYS:HG2	2.12	0.49
1:A:63:PHE:CE1	1:B:306:LYS:NZ	2.72	0.49
1:B:11:VAL:N	1:B:12:PRO:CD	2.75	0.49
1:A:205:ASN:O	1:A:340:ASP:HA	2.13	0.49
1:B:159:LEU:HD12	1:B:204:ILE:O	2.12	0.49
1:A:152:MDO:HB22	2:A:600:TYR:CA	2.39	0.49
1:A:75:MET:HA	1:A:75:MET:HE2	1.95	0.49
1:A:338:ALA:HB3	1:B:254:ALA:HB1	1.94	0.49
1:A:82:GLU:HG3	1:A:197:PHE:CD1	2.48	0.49
1:B:101:PHE:CE1	1:B:146:GLU:HA	2.47	0.48
1:B:308:TYR:HE2	1:B:311:ARG:HH22	1.61	0.48
1:B:295:LYS:HB3	1:B:298:GLN:HG2	1.95	0.48
1:A:315:GLN:HG2	1:B:354:ALA:O	2.13	0.48
1:B:486:LEU:HB3	1:B:490:ALA:HB3	1.95	0.48
1:A:17:GLY:O	1:A:119:LYS:HG3	2.12	0.48
1:B:283:LEU:O	1:B:287:LEU:HB2	2.13	0.48
1:A:79:LYS:HA	1:A:197:PHE:HZ	1.79	0.48
1:A:76:GLN:HB3	1:B:291:LYS:NZ	2.29	0.48
1:B:509:TYR:CE2	1:B:511:ALA:HB3	2.49	0.47
1:A:225:ALA:O	1:A:229:GLU:HG3	2.14	0.47
1:A:257:HIS:HE1	1:B:334:GLU:OE2	1.98	0.47
1:A:231:VAL:HB	1:A:470:TYR:CE1	2.50	0.47
1:A:113:ARG:HD2	1:A:208:SER:OG	2.15	0.47
1:B:284:ARG:HH22	1:B:285:ARG:HH21	1.60	0.47
1:A:63:PHE:HA	1:A:67:THR:OG1	2.14	0.47
1:A:503:THR:HG22	1:A:504:LEU:N	2.30	0.47
1:A:407:HIS:CD2	1:A:507:ASP:H	2.26	0.47
1:B:306:LYS:C	1:B:311:ARG:HH21	2.06	0.47
1:A:287:LEU:HD13	1:A:302:ILE:HB	1.97	0.47
1:B:71:GLU:H	1:B:438:ASN:ND2	2.11	0.47
1:B:72:MET:HE3	3:B:683:HOH:O	2.15	0.47
1:B:276:LEU:HD12	1:B:486:LEU:HD23	1.97	0.46
1:B:523:ARG:NH2	1:B:525:GLU:HG3	2.30	0.46
1:A:60:ILE:HD11	1:A:62:ILE:HD11	1.97	0.46
1:A:284:ARG:HG2	1:A:284:ARG:HH11	1.79	0.46
1:B:17:GLY:H	1:B:115:ASN:HD21	1.62	0.46
1:B:152:MDO:HB2	1:B:445:VAL:HG12	1.96	0.46
1:B:525:GLU:OE1	1:B:528:ARG:NH2	2.49	0.46
1:A:129:LEU:HD12	1:A:129:LEU:N	2.31	0.46
1:B:113:ARG:HH12	1:B:206:GLY:HA3	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:408:SER:HB3	1:B:411:ALA:HB3	1.97	0.46
1:B:198:LYS:HG2	1:B:198:LYS:O	2.16	0.46
1:B:329:HIS:HD2	1:B:332:ARG:NH1	2.14	0.46
1:A:184:VAL:O	1:A:187:GLU:HB3	2.15	0.46
1:A:447:MET:HE2	1:A:447:MET:HB3	1.90	0.46
1:A:330:LYS:HZ1	1:B:257:HIS:HD2	1.62	0.46
1:A:75:MET:CE	1:B:299:ARG:HA	2.39	0.46
1:B:494:TYR:CZ	1:B:498:ARG:HG3	2.50	0.46
1:A:113:ARG:NH1	1:A:206:GLY:HA3	2.30	0.46
1:A:313:ILE:HB	1:A:314:PRO:HD3	1.97	0.46
1:B:150:LEU:O	1:B:445:VAL:HB	2.16	0.46
1:A:344:PHE:CE1	1:A:350:ILE:HG23	2.51	0.46
1:A:65:VAL:HG22	1:A:66:THR:HG23	1.98	0.46
1:A:276:LEU:HD11	1:A:482:ARG:HB3	1.98	0.45
1:A:494:TYR:CE2	1:A:498:ARG:HG3	2.51	0.45
1:B:280:HIS:HB2	3:B:606:HOH:O	2.15	0.45
1:A:63:PHE:O	1:A:65:VAL:HG12	2.16	0.45
1:B:205:ASN:O	1:B:340:ASP:HA	2.17	0.45
1:B:307:ALA:CB	1:B:385:ARG:CZ	2.94	0.45
1:A:101:PHE:CE2	1:A:146:GLU:HG2	2.52	0.45
1:A:235:LEU:C	1:A:235:LEU:HD23	2.37	0.45
1:A:53:GLU:HA	1:A:56:ALA:HB3	1.98	0.45
1:A:147:ILE:HG22	1:A:148:GLY:N	2.32	0.45
1:A:399:LEU:O	1:A:479:ILE:HD11	2.17	0.45
1:A:81:LYS:HB3	1:A:84:GLU:HB3	1.97	0.45
1:B:304:LEU:HA	1:B:304:LEU:HD13	1.76	0.45
1:B:310:LEU:HD23	1:B:310:LEU:N	2.30	0.45
1:B:69:TYR:HB3	2:B:600:TYR:CD2	2.51	0.45
1:A:124:VAL:CG1	1:A:128:ILE:HB	2.46	0.44
1:B:124:VAL:HG12	1:B:125:ARG:N	2.32	0.44
1:B:486:LEU:O	1:B:491:LYS:NZ	2.50	0.44
1:A:73:ILE:CB	1:B:306:LYS:HD2	2.47	0.44
1:A:354:ALA:O	1:B:315:GLN:HG2	2.17	0.44
1:B:11:VAL:N	1:B:12:PRO:HD3	2.32	0.44
1:A:32:ARG:HH12	1:A:103:GLU:HG2	1.83	0.44
1:A:341:ASN:OD1	1:A:342:PRO:HA	2.18	0.44
1:A:37:VAL:HG13	1:A:37:VAL:O	2.18	0.44
1:B:523:ARG:HH21	1:B:525:GLU:HG3	1.82	0.44
1:A:330:LYS:NZ	1:B:257:HIS:HD2	2.15	0.44
1:A:75:MET:HA	1:A:75:MET:CE	2.47	0.44
1:B:132:LEU:HD12	1:B:132:LEU:HA	1.89	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:52:PHE:HZ	1:A:197:PHE:HB3	1.83	0.43
1:A:52:PHE:CZ	1:A:197:PHE:HB3	2.53	0.43
1:A:498:ARG:HA	1:A:501:VAL:O	2.18	0.43
1:B:64:GLY:HA3	1:B:201:LEU:HD22	2.00	0.43
1:A:35:VAL:N	1:A:137:ASN:HD21	1.99	0.43
1:A:97:VAL:HG23	1:A:147:ILE:HB	2.00	0.43
1:B:69:TYR:CE2	1:B:89:LEU:HD22	2.53	0.43
1:A:146:GLU:OE1	1:A:453:ARG:HD2	2.19	0.43
1:B:70:GLY:HA3	1:B:438:ASN:HD21	1.83	0.43
1:A:260:GLN:NE2	1:B:355:ASN:HD21	2.17	0.43
1:B:26:ARG:HD2	1:B:219:GLY:HA3	2.00	0.43
1:B:263:THR:O	1:B:267:MET:HG2	2.17	0.43
1:B:295:LYS:CE	1:B:296:ASP:N	2.79	0.43
1:A:260:GLN:HE22	1:B:355:ASN:HD21	1.66	0.43
1:A:303:TYR:HB3	1:A:305:GLN:O	2.19	0.43
1:A:230:ILE:HA	1:A:270:LEU:HD13	2.01	0.43
1:A:416:PRO:O	1:A:420:LEU:HG	2.18	0.43
1:A:231:VAL:HB	1:A:470:TYR:CZ	2.53	0.42
1:B:381:ARG:NE	1:B:381:ARG:HA	2.33	0.42
1:A:62:ILE:HB	1:A:65:VAL:CG1	2.49	0.42
1:B:244:SER:N	1:B:245:PRO:HD2	2.34	0.42
1:B:399:LEU:O	1:B:479:ILE:HD11	2.18	0.42
1:A:346:GLU:O	1:A:347:GLY:C	2.58	0.42
1:A:255:ARG:HD3	1:A:260:GLN:OE1	2.20	0.42
1:A:407:HIS:NE2	1:A:506:VAL:HB	2.35	0.42
1:A:251:HIS:HD2	1:A:260:GLN:HE21	1.67	0.42
1:B:299:ARG:CG	1:B:299:ARG:NH1	2.78	0.42
1:B:308:TYR:CE2	1:B:311:ARG:NH2	2.87	0.42
1:B:475:GLN:HE21	1:B:479:ILE:HD11	1.85	0.42
1:A:159:LEU:HB3	1:A:204:ILE:HA	2.02	0.42
1:A:313:ILE:N	1:A:313:ILE:HD12	2.35	0.41
1:A:305:GLN:HE21	1:B:352:HIS:HB3	1.85	0.41
1:A:378:LEU:O	1:A:382:GLN:HG3	2.19	0.41
1:A:443:ASP:OD2	1:B:309:SER:OG	2.36	0.41
1:B:113:ARG:HD2	1:B:208:SER:OG	2.20	0.41
1:B:307:ALA:HB1	1:B:385:ARG:CZ	2.50	0.41
1:A:63:PHE:CD1	1:B:306:LYS:NZ	2.88	0.41
1:A:387:LEU:HD23	1:A:399:LEU:HG	2.03	0.41
1:B:113:ARG:HA	1:B:113:ARG:HD2	1.64	0.41
1:A:260:GLN:HE22	1:B:355:ASN:ND2	2.18	0.41
1:A:42:ILE:O	1:A:45:ALA:HB3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:35:VAL:HG12	1:A:137:ASN:HD21	1.82	0.41
1:A:76:GLN:OE1	1:B:287:LEU:HD22	2.21	0.41
1:A:150:LEU:O	1:A:445:VAL:HB	2.21	0.41
1:A:237:GLU:OE2	1:A:490:ALA:HB2	2.20	0.41
1:A:69:TYR:CZ	1:A:89:LEU:HD22	2.56	0.41
1:B:295:LYS:CE	1:B:297:VAL:N	2.56	0.41
1:A:515:GLU:HA	1:A:515:GLU:OE2	2.20	0.41
1:A:470:TYR:CE1	1:A:521:LEU:HD21	2.57	0.40
1:B:306:LYS:O	1:B:307:ALA:C	2.59	0.40
1:A:251:HIS:NE2	1:A:314:PRO:CB	2.84	0.40
1:A:60:ILE:HA	1:A:61:PRO:HD3	1.83	0.40
1:A:263:THR:O	1:A:267:MET:HG2	2.22	0.40
1:A:431:SER:HA	1:A:446:SER:O	2.21	0.40
1:B:127:ILE:HG23	1:B:128:ILE:N	2.36	0.40
1:B:307:ALA:HB2	1:B:385:ARG:CZ	2.51	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)
1:B:423:GLU:OE1	1:B:453:ARG:NH2[2_555]	2.09	0.11

## 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	524/537 (98%)	498 (95%)	25 (5%)	1 (0%)	51	58
1	B	524/537 (98%)	505 (96%)	19 (4%)	0	100	100
All	All	1048/1074 (98%)	1003 (96%)	44 (4%)	1 (0%)	55	63

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	505	GLY

### 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	416/425 (98%)	412 (99%)	4 (1%)	80	89
1	B	416/425 (98%)	406 (98%)	10 (2%)	54	67
All	All	832/850 (98%)	818 (98%)	14 (2%)	66	79

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

Mol	Chain	Res	Type
1	A	65	VAL
1	A	103	GLU
1	A	140	ILE
1	A	498	ARG
1	B	132	LEU
1	B	175	ASP
1	B	252	ASP
1	B	290	ASP
1	B	292	GLU
1	B	295	LYS
1	B	297	VAL
1	B	299	ARG
1	B	304	LEU
1	B	447	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	121	HIS
1	A	137	ASN
1	A	161	HIS
1	A	205	ASN

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Mol	Chain	Res	Type
1	A	227	GLN
1	A	251	HIS
1	A	257	HIS
1	A	305	GLN
1	A	339	ASN
1	A	355	ASN
1	A	390	HIS
1	A	407	HIS
1	A	442	GLN
1	B	115	ASN
1	B	121	HIS
1	B	137	ASN
1	B	161	HIS
1	B	205	ASN
1	B	227	GLN
1	B	251	HIS
1	B	257	HIS
1	B	280	HIS
1	B	288	GLN
1	B	326	HIS
1	B	329	HIS
1	B	339	ASN
1	B	355	ASN
1	B	407	HIS
1	B	438	ASN
1	B	442	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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
1	MDO	A	152	1	12,13,14	2.36	5 (41%)	12,18,20	3.19	2 (16%)
1	MDO	B	152	1	12,13,14	2.46	6 (50%)	12,18,20	3.00	2 (16%)

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
1	MDO	A	152	1	-	0/4/23/24	0/1/1/1
1	MDO	B	152	1	-	0/4/23/24	0/1/1/1

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	152	MDO	CA2-N2	-2.75	1.33	1.39
1	B	152	MDO	C2-N3	-2.72	1.33	1.39
1	A	152	MDO	C2-N3	-2.71	1.33	1.39
1	B	152	MDO	CA2-N2	-2.67	1.33	1.39
1	B	152	MDO	CA2-C2	-2.18	1.39	1.43
1	B	152	MDO	C1-N3	-2.15	1.33	1.37
1	A	152	MDO	CA2-C2	-2.14	1.39	1.43
1	A	152	MDO	C1-N3	-2.13	1.33	1.37
1	B	152	MDO	O-C	2.19	1.33	1.19
1	A	152	MDO	O2-C2	6.13	1.36	1.23
1	B	152	MDO	O2-C2	6.14	1.36	1.23

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	152	MDO	O2-C2-CA2	-7.97	126.67	130.97
1	B	152	MDO	O2-C2-CA2	-7.39	126.98	130.97
1	B	152	MDO	CA2-C2-N3	6.70	106.29	103.30
1	A	152	MDO	CA2-C2-N3	7.28	106.55	103.30

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	152	MDO	5	0
1	B	152	MDO	2	0

## 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	$\# Z  > 2$
2	TYR	A	600	-	9,13,13	0.44	0	12,17,17	0.71	0
2	TYR	B	600	-	9,13,13	0.29	0	12,17,17	0.89	1 (8%)

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	TYR	A	600	-	-	0/4/8/8	0/1/1/1
2	TYR	B	600	-	-	0/4/8/8	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	600	TYR	CG-CB-CA	-2.40	109.47	114.43

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	600	TYR	7	0
2	B	600	TYR	4	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	526/537 (97%)	0.21	35 (6%) 19 17	24, 40, 70, 80	0
1	B	526/537 (97%)	0.03	23 (4%) 35 33	23, 38, 63, 90	0
All	All	1052/1074 (97%)	0.12	58 (5%) 26 25	23, 39, 69, 90	0

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	297	VAL	9.4
1	B	304	LEU	8.5
1	A	402	GLY	7.1
1	B	296	ASP	6.9
1	B	293	ALA	6.2
1	A	293	ALA	5.5
1	A	11	VAL	5.2
1	A	295	LYS	5.1
1	A	59	ASN	4.9
1	A	344	PHE	4.8
1	B	299	ARG	4.8
1	A	52	PHE	4.7
1	B	295	LYS	4.7
1	A	294	GLY	4.5
1	A	54	GLY	4.5
1	A	51	ILE	4.3
1	B	292	GLU	4.3
1	A	77	VAL	4.0
1	B	300	SER	4.0
1	A	57	GLU	3.7
1	B	294	GLY	3.6
1	A	189	GLY	3.6
1	B	298	GLN	3.6
1	B	11	VAL	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	60	ILE	3.4
1	A	296	ASP	3.3
1	A	350	ILE	3.3
1	A	78	ASP	2.9
1	A	535	ASP	2.9
1	A	174	ARG	2.9
1	B	60	ILE	2.8
1	B	59	ASN	2.8
1	A	186	ALA	2.7
1	B	186	ALA	2.7
1	A	79	LYS	2.7
1	B	288	GLN	2.7
1	A	349	GLU	2.7
1	A	55	ILE	2.7
1	A	393	TYR	2.7
1	A	73	ILE	2.6
1	A	479	ILE	2.6
1	A	484	ASP	2.5
1	B	312	ALA	2.5
1	A	74	TYR	2.5
1	A	297	VAL	2.4
1	A	483	PHE	2.4
1	B	344	PHE	2.4
1	A	176	GLY	2.4
1	A	444	VAL	2.4
1	B	303	TYR	2.3
1	A	81	LYS	2.3
1	A	403	ASP	2.2
1	B	78	ASP	2.2
1	A	56	ALA	2.1
1	B	435	VAL	2.1
1	B	310	LEU	2.1
1	B	444	VAL	2.0
1	B	437	SER	2.0

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

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
1	MDO	A	152	13/14	0.86	0.22	-	37,41,45,50	0
1	MDO	B	152	13/14	0.86	0.21	-	36,40,44,50	0

### 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	TYR	A	600	13/13	0.77	0.71	19.22	31,33,35,35	13
2	TYR	B	600	13/13	0.75	0.60	13.47	24,26,27,28	13

### 6.5 Other polymers [i](#)

There are no such residues in this entry.