



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 11:12 AM GMT

PDB ID : 1O7M  
Title : NAPHTHALENE 1,2-DIOXYGENASE, BINARY COMPLEX WITH DIOXYGEN  
Authors : Karlsson, A.; Parales, J.V.; Parales, R.E.; Gibson, D.T.; Eklund, H.; Ramaswamy, S.  
Deposited on : 2002-11-11  
Resolution : 1.75 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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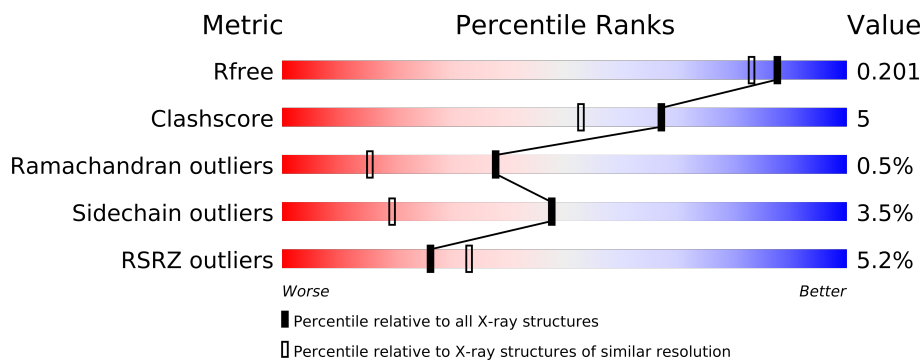
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 1.75 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))
$R_{free}$	66092	1134 (1.76-1.76)
Clashscore	79885	1304 (1.76-1.76)
Ramachandran outliers	78287	1288 (1.76-1.76)
Sidechain outliers	78261	1288 (1.76-1.76)
RSRZ outliers	66119	1135 (1.76-1.76)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	449	
2	B	194	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	EDO	A	1449	-	X
3	EDO	A	1450	-	X
3	EDO	B	1699	-	X
4	OXY	A	1451	X	-
5	SO4	B	1697	-	X

## 2 Entry composition i

There are 8 unique types of molecules in this entry. The entry contains 5694 atoms, of which 0 are hydrogen and 0 are deuterium.

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 NAPHTHALENE 1,2-DIOXYGENASE ALPHA SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	448	Total	C	N	O	S	0	2	1
			3487	2205	599	666	17			

- Molecule 2 is a protein called NAPHTHALENE 1,2-DIOXYGENASE BETA SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	193	Total	C	N	O	S	0	0	0
			1608	1007	302	293	6			

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



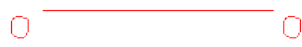
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		

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

- Molecule 4 is OXYGEN MOLECULE (three-letter code: OXY) (formula: O<sub>2</sub>).



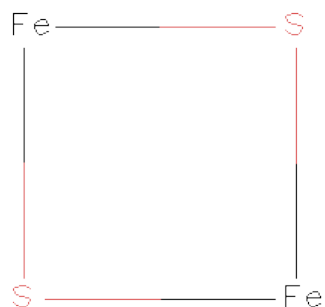
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total O 2 2	0	0

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 6 is FE2/S2 (INORGANIC) CLUSTER (three-letter code: FES) (formula:  $\text{Fe}_2\text{S}_2$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	Fe	S	0	0
			4	2	2		

- Molecule 7 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	1	Total 1	Fe 1	0	0

- Molecule 8 is water.

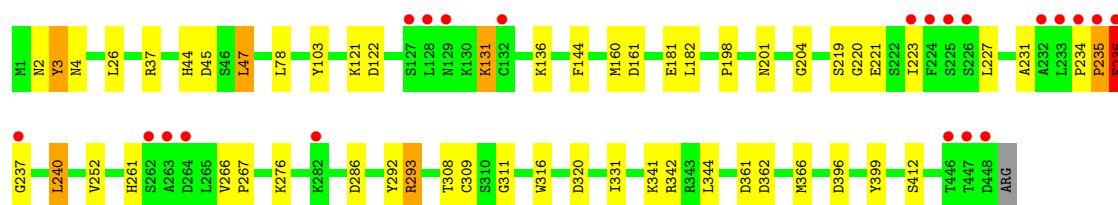
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	401	Total 401	O 401	0	0
8	B	152	Total 152	O 152	0	0

### 3 Residue-property plots

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.

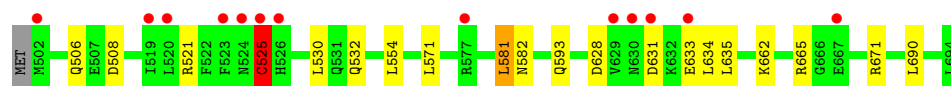
- Molecule 1: NAPHTHALENE 1,2-DIOXYGENASE ALPHA SUBUNIT

Chain A: 



- Molecule 2: NAPHTHALENE 1,2-DIOXYGENASE BETA SUBUNIT

Chain B: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	140.28Å 140.28Å 208.75Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 1.75 32.06 – 1.56	Depositor EDS
% Data completeness (in resolution range)	98.9 (30.00-1.75) 84.8 (32.06-1.56)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.53 (at 1.57Å)	Xtriage
Refinement program	REFMAC 5.0.36	Depositor
R, $R_{free}$	0.178 , 0.187 0.193 , 0.201	Depositor DCC
$R_{free}$ test set	771 reflections (0.99%)	DCC
Wilson B-factor (Å <sup>2</sup> )	16.5	Xtriage
Anisotropy	0.124	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.41 , 38.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	1 of 94013 reflections (0.001%)	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	5694	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.03% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.



## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, EDO, FE, FES, OXY

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	0.66	0/3589	0.86	10/4862 (0.2%)
2	B	0.61	0/1638	0.89	1/2209 (0.0%)
All	All	0.64	0/5227	0.87	11/7071 (0.2%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	286	ASP	CB-CG-OD2	7.18	124.76	118.30
1	A	320	ASP	CB-CG-OD2	6.65	124.28	118.30
1	A	293	ARG	NE-CZ-NH1	-6.50	117.05	120.30
1	A	37	ARG	NE-CZ-NH2	6.46	123.53	120.30
1	A	361	ASP	CB-CG-OD2	6.19	123.87	118.30
1	A	161	ASP	CB-CG-OD2	5.72	123.44	118.30
1	A	45	ASP	CB-CG-OD2	5.59	123.33	118.30
2	B	628	ASP	CB-CG-OD2	5.55	123.30	118.30
1	A	3	TYR	CB-CG-CD2	5.36	124.22	121.00
1	A	122	ASP	CB-CG-OD2	5.15	122.93	118.30
1	A	362	ASP	CB-CG-OD2	5.12	122.91	118.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the

chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3487	0	3323	36	0
2	B	1608	0	1583	14	0
3	A	12	0	18	0	0
3	B	12	0	18	4	0
4	A	2	0	0	0	0
5	A	5	0	0	1	0
5	B	10	0	0	0	0
6	A	4	0	0	0	0
7	A	1	0	0	0	0
8	A	401	0	0	3	0
8	B	152	0	0	0	0
All	All	5694	0	4942	50	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 5.

All (50) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:240:LEU:HD13	1:A:252:VAL:HB	1.40	1.04
1:A:131:LYS:HD3	5:A:1452:SO4:O2	1.68	0.93
1:A:342:ARG:HH11	1:A:342:ARG:HG2	1.34	0.89
1:A:227:LEU:CD2	1:A:234:PRO:HD3	2.14	0.77
1:A:342:ARG:HG2	1:A:342:ARG:NH1	1.99	0.77
2:B:581:LEU:HD13	3:B:1696:EDO:H21	1.66	0.75
1:A:47:LEU:HD22	1:A:182:LEU:HD23	1.80	0.62
1:A:240:LEU:CD1	1:A:252:VAL:HB	2.24	0.62
2:B:525:CYS:O	2:B:525:CYS:SG	2.59	0.61
2:B:581:LEU:HD13	3:B:1696:EDO:C2	2.31	0.60
1:A:227:LEU:HD21	1:A:234:PRO:HD3	1.84	0.59
1:A:235:PRO:HB2	1:A:236:GLU:OE1	2.03	0.59
1:A:227:LEU:HD22	1:A:234:PRO:HD3	1.84	0.59
1:A:220:GLY:HA3	8:A:2228:HOH:O	2.03	0.59
2:B:581:LEU:CD1	3:B:1696:EDO:H21	2.33	0.58
2:B:634:LEU:HD11	2:B:662:LYS:HB3	1.87	0.56
2:B:631:ASP:OD1	2:B:633:GLU:HB2	2.07	0.55
1:A:366[B]:MET:HE1	8:A:2219:HOH:O	2.06	0.55
1:A:240:LEU:HD13	1:A:252:VAL:CB	2.28	0.54
1:A:220:GLY:O	1:A:221:GLU:HB2	2.09	0.53
1:A:266:VAL:HB	1:A:267:PRO:HD3	1.91	0.52
1:A:3:TYR:CE2	1:A:26:LEU:HD13	2.45	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:144:PHE:HD1	1:A:160:MET:HE1	1.75	0.51
2:B:665:ARG:HG2	2:B:671:ARG:HG2	1.92	0.50
1:A:231:ALA:HB2	8:A:2230:HOH:O	2.12	0.50
1:A:144:PHE:CD1	1:A:160:MET:HE1	2.48	0.49
1:A:204:GLY:HA3	1:A:412:SER:HB2	1.96	0.48
1:A:2:ASN:OD1	1:A:4:ASN:HB2	2.12	0.48
2:B:665:ARG:HA	2:B:671:ARG:HA	1.96	0.47
1:A:236:GLU:HB2	1:A:237:GLY:H	1.36	0.47
1:A:44:HIS:HB2	1:A:182:LEU:HD22	1.97	0.47
2:B:581:LEU:HD13	3:B:1696:EDO:C1	2.45	0.46
2:B:634:LEU:HD21	2:B:662:LYS:HD3	1.96	0.46
1:A:198:PRO:HB2	1:A:316:TRP:CE2	2.52	0.45
1:A:234:PRO:HA	1:A:235:PRO:HD2	1.79	0.45
1:A:160:MET:HB2	1:A:160:MET:HE3	1.65	0.45
1:A:261:HIS:NE2	1:A:344:LEU:HD11	2.32	0.45
2:B:581:LEU:O	2:B:582:ASN:C	2.56	0.44
1:A:181:GLU:OE1	1:A:341:LYS:NZ	2.40	0.44
2:B:530:LEU:HD21	2:B:635:LEU:HD13	2.01	0.43
1:A:311:GLY:HA2	1:A:331:ILE:HG13	2.00	0.43
1:A:236:GLU:H	1:A:236:GLU:HG3	1.46	0.43
1:A:78:LEU:HD22	1:A:136:LYS:HD3	2.02	0.42
1:A:276:LYS:HD2	1:A:292:TYR:CZ	2.55	0.41
1:A:261:HIS:NE2	1:A:344:LEU:CD1	2.84	0.41
1:A:396:ASP:HB3	1:A:399:TYR:O	2.21	0.40
1:A:308:THR:O	1:A:309:CYS:HB2	2.20	0.40
2:B:631:ASP:C	2:B:633:GLU:H	2.23	0.40
2:B:631:ASP:C	2:B:633:GLU:N	2.75	0.40
1:A:223:ILE:HA	1:A:223:ILE:HD12	1.89	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	448/449 (100%)	431 (96%)	15 (3%)	2 (0%)	43 21

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
2	B	191/194 (98%)	184 (96%)	6 (3%)	1 (0%)	38 15
All	All	639/643 (99%)	615 (96%)	21 (3%)	3 (0%)	38 15

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	236	GLU
2	B	525	CYS
1	A	235	PRO

### 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	369/369 (100%)	360 (98%)	9 (2%)	61 34
2	B	172/173 (99%)	162 (94%)	10 (6%)	28 7
All	All	541/542 (100%)	522 (96%)	19 (4%)	48 19

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	47	LEU
1	A	103	TYR
1	A	121	LYS
1	A	131	LYS
1	A	201	ASN
1	A	219	SER
1	A	236	GLU
1	A	240	LEU
1	A	293	ARG
2	B	506	GLN
2	B	508	ASP
2	B	521	ARG
2	B	525	CYS
2	B	532	GLN
2	B	554	LEU

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Mol	Chain	Res	Type
2	B	571	LEU
2	B	581	LEU
2	B	593	GLN
2	B	690	LEU

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	115	GLN
1	A	201	ASN
2	B	506	GLN
2	B	514	HIS
2	B	526	HIS

### 5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

Of 12 ligands modelled in this entry, 1 is monoatomic - leaving 11 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	1448	-	3,3,3	0.56	0	2,2,2	0.25	0
3	EDO	A	1449	-	3,3,3	0.58	0	2,2,2	0.69	0
3	EDO	A	1450	-	3,3,3	0.49	0	2,2,2	0.44	0
4	OXY	A	1451	7	1,1,1	2.56	1 (100%)	0,0,0	0.00	-
5	SO4	A	1452	-	4,4,4	0.26	0	6,6,6	0.48	0
6	FES	A	1453	1	0,4,4	0.00	-	0,4,4	0.00	-
3	EDO	B	1695	-	3,3,3	0.43	0	2,2,2	0.44	0
3	EDO	B	1696	-	3,3,3	0.47	0	2,2,2	0.40	0
5	SO4	B	1697	-	4,4,4	0.12	0	6,6,6	0.08	0
5	SO4	B	1698	-	4,4,4	0.33	0	6,6,6	0.25	0
3	EDO	B	1699	-	3,3,3	0.54	0	2,2,2	0.43	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EDO	A	1448	-	-	0/1/1/1	0/0/0/0
3	EDO	A	1449	-	-	0/1/1/1	0/0/0/0
3	EDO	A	1450	-	-	0/1/1/1	0/0/0/0
4	OXY	A	1451	7	-	0/0/0/0	0/0/0/0
5	SO4	A	1452	-	-	0/0/0/0	0/0/0/0
6	FES	A	1453	1	-	0/0/4/4	0/0/1/1
3	EDO	B	1695	-	-	0/1/1/1	0/0/0/0
3	EDO	B	1696	-	-	0/1/1/1	0/0/0/0
5	SO4	B	1697	-	-	0/0/0/0	0/0/0/0
5	SO4	B	1698	-	-	0/0/0/0	0/0/0/0
3	EDO	B	1699	-	-	0/1/1/1	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1451	OXY	O2-O1	2.56	1.45	1.16

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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	448/449 (99%)	-0.18	21 (4%) 30 37	7, 13, 30, 44	0
2	B	193/194 (99%)	0.06	13 (6%) 17 24	8, 14, 38, 46	0
All	All	641/643 (99%)	-0.11	34 (5%) 26 32	7, 13, 33, 46	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	447	THR	10.3
2	B	502	MET	9.9
1	A	448	ASP	8.5
2	B	525	CYS	8.4
1	A	446	THR	6.5
2	B	667	GLU	5.5
1	A	235	PRO	5.5
1	A	224	PHE	5.4
1	A	236	GLU	5.1
2	B	577	ARG	5.0
1	A	232	ALA	4.7
2	B	523	PHE	4.6
1	A	234	PRO	4.4
1	A	264	ASP	4.2
2	B	630	ASN	4.2
2	B	629	VAL	4.1
2	B	633	GLU	3.7
1	A	226	SER	3.3
1	A	233	LEU	3.3
1	A	262	SER	3.1
2	B	524	ASN	3.1
1	A	127	SER	3.0
2	B	519	ILE	2.9
2	B	631	ASP	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	282	LYS	2.8
1	A	225	SER	2.6
2	B	520	LEU	2.5
1	A	263	ALA	2.3
1	A	129	ASN	2.3
1	A	237	GLY	2.3
1	A	132	CYS	2.2
2	B	526	HIS	2.2
1	A	223	ILE	2.1
1	A	128	LEU	2.1

## 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 ⓘ

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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
5	SO4	B	1697	5/5	0.31	46.14	45,47,49,50	0
3	EDO	B	1699	4/4	0.11	5.84	12,22,23,26	0
3	EDO	A	1450	4/4	0.22	5.11	41,41,41,43	0
3	EDO	A	1449	4/4	0.11	2.91	15,21,22,22	0
5	SO4	A	1452	5/5	0.23	1.20	51,52,53,53	0
7	FE	A	1454	1/1	0.09	1.13	22,22,22,22	0
4	OXY	A	1451	2/2	0.09	1.10	36,36,36,38	0
3	EDO	B	1696	4/4	0.14	0.34	24,26,27,28	0
3	EDO	A	1448	4/4	0.06	0.01	8,10,11,11	0
5	SO4	B	1698	5/5	0.12	-0.21	36,36,37,37	0
3	EDO	B	1695	4/4	0.06	-1.07	11,12,13,13	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
6	FES	A	1453	4/4	0.03	-2.72	11,12,13,14	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.