



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 11:50 PM UTC

PDB ID : 1Q0O / pdb_00001q0o
Title : CRYSTAL STRUCTURE OF HOMOPROTOCATECHUATE 2,3-DIOXYGENASE FROM BREVIBACTERIUM FUSCUM (FULL LENGTH PROTEIN)
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Deposited on : 2003-07-16
Resolution : 2.30 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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

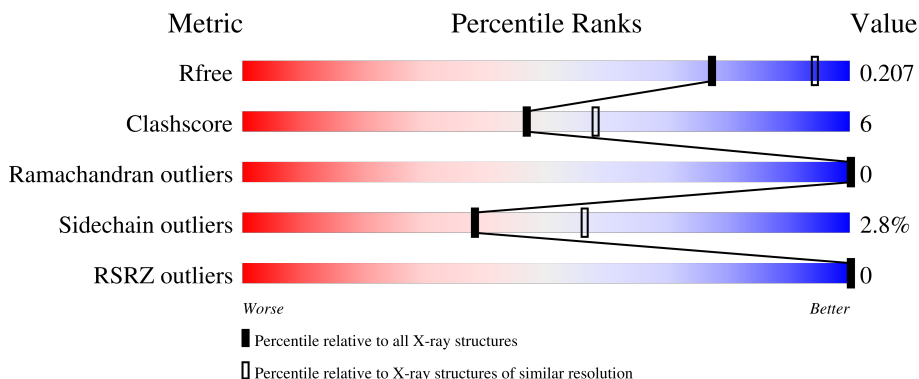
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	365	
1	B	365	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5955 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 homoprotocatechuate 2,3-dioxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	356	Total	C	N	O	S	0	0	0
			2879	1815	506	551	7			
1	B	356	Total	C	N	O	S	0	0	0
			2879	1815	506	551	7			

- Molecule 2 is FE (III) ION (CCD ID: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Fe	0	0
			1	1		
2	B	1	Total	Fe	0	0
			1	1		

- Molecule 3 is water.

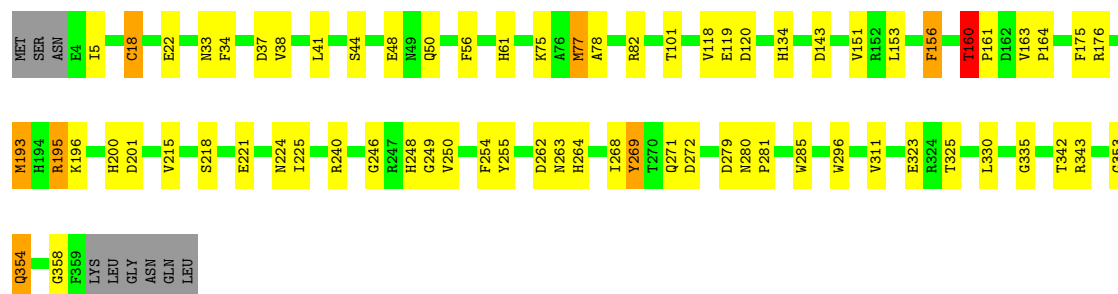
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	91	Total	O	0	0
			91	91		
3	B	104	Total	O	0	0
			104	104		

3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: homoprotocatechuate 2,3-dioxygenase

Chain A: 



- Molecule 1: homoprotocatechuate 2,3-dioxygenase

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	118.90Å 118.90Å 110.30Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.30 20.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	80.8 (20.00-2.30) 80.8 (20.00-2.30)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.07 (at 2.31Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.161 , 0.208 0.160 , 0.207	Depositor DCC
R_{free} test set	1613 reflections (4.42%)	wwPDB-VP
Wilson B-factor (Å ²)	15.5	Xtriage
Anisotropy	0.307	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 33.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.039 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5955	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.05% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.74	0/2955	1.16	31/4014 (0.8%)
1	B	0.74	1/2955 (0.0%)	1.15	27/4014 (0.7%)
All	All	0.74	1/5910 (0.0%)	1.15	58/8028 (0.7%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	193	MET	SD-CE	5.06	1.92	1.79

All (58) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	5	ILE	CA-C-N	8.81	128.75	119.85
1	B	5	ILE	C-N-CA	8.81	128.75	119.85
1	B	201	ASP	N-CA-C	-8.61	103.29	113.88
1	A	160	THR	CA-C-N	7.66	126.94	118.97
1	A	160	THR	C-N-CA	7.66	126.94	118.97
1	B	342	THR	N-CA-C	-7.46	102.76	112.68
1	B	271	GLN	N-CA-C	7.33	121.78	111.52
1	A	196	LYS	N-CA-C	6.94	121.60	113.20
1	B	196	LYS	N-CA-C	6.71	121.32	113.20
1	A	18	CYS	N-CA-C	-6.60	101.19	110.50
1	A	201	ASP	N-CA-C	-6.57	105.80	113.88
1	B	198	THR	N-CA-C	-6.50	100.33	109.69
1	A	271	GLN	N-CA-C	6.43	120.53	111.52
1	A	342	THR	N-CA-C	-6.39	104.18	112.68
1	A	82	ARG	N-CA-C	6.35	118.28	111.36
1	A	44	SER	N-CA-C	-6.28	104.67	112.90
1	A	264	HIS	N-CA-C	-6.19	101.12	110.28
1	A	255	TYR	N-CA-C	6.09	118.22	109.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	151	VAL	N-CA-C	6.00	120.81	112.35
1	A	153	LEU	N-CA-C	-6.00	99.50	109.46
1	B	18	CYS	N-CA-C	-5.97	101.45	110.28
1	B	139	HIS	N-CA-C	5.96	119.65	112.38
1	B	61	HIS	N-CA-C	5.96	118.54	111.33
1	B	156	PHE	N-CA-C	-5.93	100.74	109.81
1	B	78	ALA	N-CA-C	5.92	118.90	109.24
1	B	151	VAL	N-CA-C	5.91	120.68	112.35
1	A	77	MET	N-CA-C	-5.90	98.80	108.41
1	B	82	ARG	N-CA-C	5.89	117.70	111.28
1	A	56	PHE	N-CA-C	5.79	118.34	111.33
1	B	268	ILE	N-CA-C	-5.77	100.12	108.71
1	A	119	GLU	N-CA-C	-5.74	97.07	107.75
1	B	56	PHE	N-CA-C	5.60	118.11	111.33
1	B	12	ALA	CA-C-N	-5.58	114.84	120.31
1	B	12	ALA	C-N-CA	-5.58	114.84	120.31
1	A	269	TYR	N-CA-CB	-5.57	101.16	111.53
1	B	215	VAL	N-CA-C	-5.56	100.44	108.89
1	A	215	VAL	N-CA-C	-5.55	100.30	108.85
1	A	254	PHE	N-CA-C	-5.53	101.46	110.20
1	B	270	THR	N-CA-C	5.52	117.48	108.76
1	A	5	ILE	CA-C-N	5.50	125.41	119.85
1	A	5	ILE	C-N-CA	5.50	125.41	119.85
1	A	272	ASP	N-CA-C	-5.45	101.41	109.86
1	A	200	HIS	N-CA-C	5.44	116.80	108.42
1	B	153	LEU	N-CA-C	-5.43	100.83	109.96
1	B	254	PHE	N-CA-C	-5.43	101.61	110.20
1	A	61	HIS	N-CA-C	5.43	117.90	111.33
1	B	200	HIS	N-CA-C	5.38	117.02	108.79
1	B	249	GLY	N-CA-C	-5.32	100.56	113.18
1	A	143	ASP	N-CA-C	-5.30	106.65	113.01
1	A	78	ALA	N-CA-C	5.28	118.08	109.59
1	A	195	ARG	N-CA-C	-5.27	106.17	113.18
1	A	249	GLY	N-CA-C	-5.27	100.68	113.18
1	A	156	PHE	N-CA-C	-5.17	101.69	109.85
1	B	9	VAL	N-CA-C	-5.13	105.11	112.35
1	A	358	GLY	N-CA-C	5.13	122.40	115.32
1	B	264	HIS	N-CA-C	-5.12	102.27	109.96
1	A	50	GLN	N-CA-C	5.08	116.80	109.07
1	B	149	GLU	N-CA-C	5.08	117.71	110.24

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	2879	0	2724	34	0
1	B	2879	0	2724	33	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	91	0	0	1	0
3	B	104	0	0	1	0
All	All	5955	0	5448	64	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 6.

All (64) 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:221:GLU:H	1:A:224:ASN:HD22	1.26	0.83
1:A:176:ARG:HG2	3:A:556:HOH:O	1.82	0.80
1:A:48:GLU:H	1:A:48:GLU:CD	1.91	0.77
1:A:101:THR:HG22	1:A:118:VAL:HG12	1.73	0.71
1:B:221:GLU:H	1:B:224:ASN:HD22	1.40	0.67
1:A:335:GLY:O	1:A:353:GLY:HA2	1.99	0.63
1:B:218:SER:HA	1:B:269:TYR:O	1.99	0.63
1:A:221:GLU:H	1:A:224:ASN:ND2	1.97	0.62
1:A:77:MET:HE3	1:A:156:PHE:CD2	2.39	0.58
1:A:250:VAL:HG22	1:A:285:TRP:CD1	2.39	0.58
1:B:225:ILE:HG12	1:B:268:ILE:CG2	2.35	0.56
1:B:183:ASP:CG	1:B:187:THR:HG23	2.32	0.54
1:B:183:ASP:OD1	1:B:187:THR:HG23	2.08	0.54
1:B:221:GLU:H	1:B:224:ASN:ND2	2.04	0.54
1:B:4:GLU:HG2	1:B:5:ILE:N	2.22	0.53
1:B:183:ASP:C	1:B:185:GLU:H	2.17	0.52
1:B:183:ASP:OD2	1:B:187:THR:HG23	2.10	0.52
1:A:175:PHE:HB3	1:A:193:MET:HB2	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:279:ASP:OD2	1:B:134:HIS:HE1	1.92	0.52
1:A:48:GLU:CD	1:A:48:GLU:N	2.67	0.51
1:B:225:ILE:HG12	1:B:268:ILE:HG22	1.92	0.51
1:A:218:SER:HA	1:A:269:TYR:O	2.11	0.50
1:A:33:ASN:O	1:A:37:ASP:HB2	2.11	0.50
1:B:47:ASP:OD1	1:B:47:ASP:C	2.55	0.49
1:B:162:ASP:OD1	1:B:164:PRO:HD2	2.13	0.49
1:A:330:LEU:C	1:A:330:LEU:HD23	2.37	0.49
1:A:262:ASP:O	1:A:263:ASN:HB2	2.14	0.48
1:A:134:HIS:HE1	1:B:279:ASP:OD2	1.97	0.48
1:B:4:GLU:CG	1:B:5:ILE:N	2.77	0.47
1:B:356:SER:O	1:B:357:LYS:HB2	2.13	0.47
1:A:246:GLY:HA2	1:A:296:TRP:CZ3	2.48	0.47
1:B:18:CYS:HB2	1:B:156:PHE:CE2	2.49	0.47
1:B:48:GLU:H	1:B:48:GLU:CD	2.21	0.47
1:B:77:MET:HE3	1:B:156:PHE:CD2	2.50	0.47
1:A:176:ARG:NH1	1:A:195:ARG:HA	2.29	0.47
1:A:163:VAL:HB	1:A:164:PRO:HD3	1.96	0.46
1:A:41:LEU:HD23	1:A:41:LEU:HA	1.81	0.46
1:B:246:GLY:HA2	1:B:296:TRP:CZ3	2.50	0.46
1:A:193:MET:HE3	1:A:193:MET:HB3	1.86	0.46
1:B:330:LEU:C	1:B:330:LEU:HD23	2.41	0.46
1:A:22:GLU:CD	1:A:75:LYS:HD3	2.41	0.45
1:B:101:THR:HB	1:B:116:LEU:HD11	1.97	0.45
1:B:203:ALA:HB1	3:B:603:HOH:O	2.18	0.44
1:A:225:ILE:HG12	1:A:268:ILE:HG22	1.99	0.44
1:A:248:HIS:HB3	1:A:250:VAL:O	2.18	0.43
1:B:262:ASP:O	1:B:263:ASN:HB2	2.18	0.43
1:B:101:THR:HG22	1:B:118:VAL:HG12	1.99	0.43
1:B:28:LEU:HD12	1:B:28:LEU:HA	1.92	0.42
1:A:225:ILE:HG12	1:A:268:ILE:CG2	2.49	0.42
1:B:225:ILE:HG12	1:B:268:ILE:HG21	2.00	0.42
1:B:5:ILE:HA	1:B:6:PRO:HD3	1.63	0.42
1:A:160:THR:HA	1:A:161:PRO:HD3	1.88	0.41
1:A:101:THR:CG2	1:A:118:VAL:HG12	2.48	0.41
1:B:58:GLU:CD	1:B:60:ILE:HG12	2.45	0.41
1:A:18:CYS:HB2	1:A:156:PHE:CE2	2.56	0.41
1:A:240:ARG:HA	1:A:240:ARG:HD2	1.89	0.41
1:A:354:GLN:HE21	1:A:354:GLN:HB3	1.57	0.41
1:B:294:ASP:OD2	1:B:298:ASN:HB2	2.21	0.41
1:A:279:ASP:OD2	1:B:134:HIS:CE1	2.73	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:34:PHE:O	1:A:38:VAL:HB	2.20	0.41
1:A:22:GLU:HG2	1:A:75:LYS:HB3	2.03	0.41
1:B:119:GLU:HG3	1:B:314:LEU:HG	2.03	0.41
1:A:280:ASN:HA	1:A:281:PRO:HD3	1.94	0.40
1:B:250:VAL:HG23	1:B:292:ARG:O	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	354/365 (97%)	344 (97%)	10 (3%)	0	100	100
1	B	354/365 (97%)	343 (97%)	11 (3%)	0	100	100
All	All	708/730 (97%)	687 (97%)	21 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	305/313 (97%)	297 (97%)	8 (3%)	40	59
1	B	305/313 (97%)	296 (97%)	9 (3%)	36	53
All	All	610/626 (97%)	593 (97%)	17 (3%)	38	56

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	120	ASP
1	A	160	THR
1	A	193	MET
1	A	311	VAL
1	A	323	GLU
1	A	325	THR
1	A	343	ARG
1	A	354	GLN
1	B	120	ASP
1	B	184	ASP
1	B	187	THR
1	B	193	MET
1	B	258	ILE
1	B	323	GLU
1	B	325	THR
1	B	328	SER
1	B	332	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	33	ASN
1	A	95	GLN
1	A	134	HIS
1	A	158	GLN
1	A	224	ASN
1	A	354	GLN
1	B	33	ASN
1	B	95	GLN
1	B	134	HIS
1	B	224	ASN
1	B	271	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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, 95th 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 ²)	Q<0.9
1	A	356/365 (97%)	-0.58	0 100 100	4, 13, 35, 56	0
1	B	356/365 (97%)	-0.63	0 100 100	4, 13, 31, 49	0
All	All	712/730 (97%)	-0.61	0 100 100	4, 13, 34, 56	0

There are no RSRZ outliers to report.

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 oligosaccharides 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. The B-factors column lists the minimum, median, 95th 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	B-factors(Å ²)	Q<0.9
2	FE	A	500	1/1	0.99	0.06	19,19,19,19	0
2	FE	B	500	1/1	0.99	0.04	14,14,14,14	0

6.5 Other polymers [i](#)

There are no such residues in this entry.