Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:47 PM GMT

PDB ID	:	1YKO
Title	:	Protocatechuate 3,4-Dioxygenase Y408H mutant
Authors	:	Brown, C.K.; Ohlendorf, D.H.
Deposited on	:	2005-01-18
Resolution	:	2.54 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
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

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

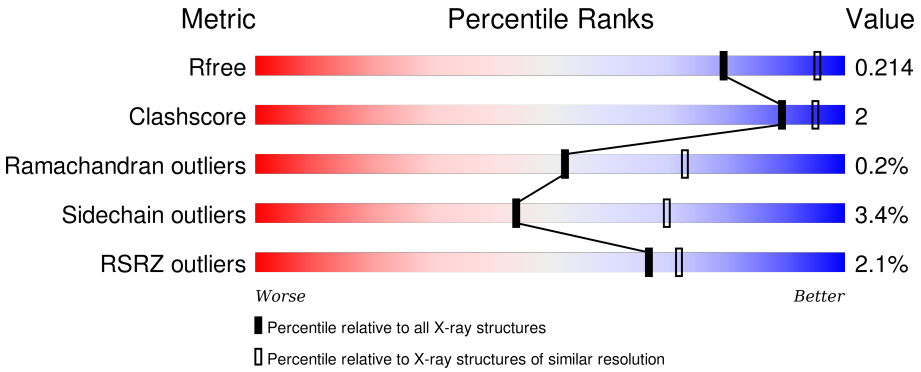
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.54 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	4549 (2.58-2.50)
Clashscore	102246	5292 (2.58-2.50)
Ramachandran outliers	100387	5194 (2.58-2.50)
Sidechain outliers	100360	5196 (2.58-2.50)
RSRZ outliers	91569	4561 (2.58-2.50)

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 >=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 <=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	200	2% 94% 5%
1	C	200	2% 94% 5%
1	E	200	2% 97%
1	G	200	4% 94% 5%
1	I	200	2% 93% 6%

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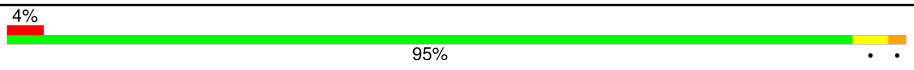
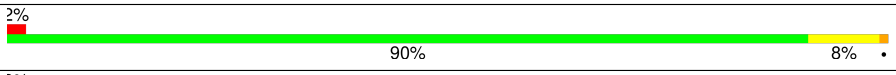
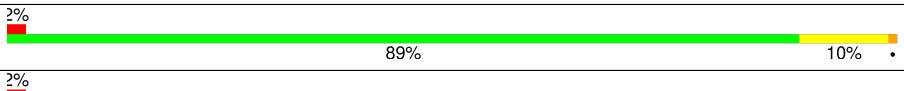
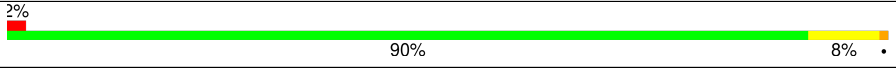
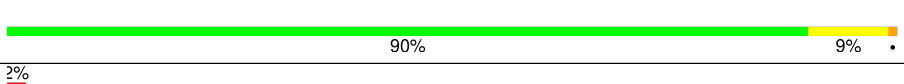
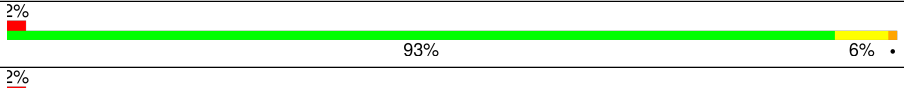
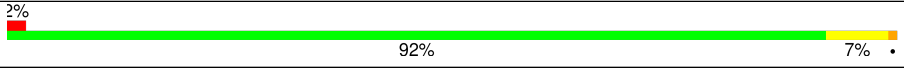
Mol	Chain	Length	Quality of chain
1	K	200	
2	B	238	
2	D	238	
2	F	238	
2	H	238	
2	J	238	
2	L	238	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 21444 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 Protocatechuate 3,4-dioxygenase alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	200	Total	C	N	O	S	0	0	0
			1571	993	276	299	3			
1	C	200	Total	C	N	O	S	0	0	0
			1571	993	276	299	3			
1	E	200	Total	C	N	O	S	0	0	0
			1571	993	276	299	3			
1	G	200	Total	C	N	O	S	0	0	0
			1571	993	276	299	3			
1	I	200	Total	C	N	O	S	0	0	0
			1571	993	276	299	3			
1	K	200	Total	C	N	O	S	0	0	0
			1571	993	276	299	3			

- Molecule 2 is a protein called Protocatechuate 3,4-dioxygenase beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	238	Total	C	N	O	S	0	0	0
			1877	1187	344	337	9			
2	D	238	Total	C	N	O	S	0	0	0
			1877	1187	344	337	9			
2	F	238	Total	C	N	O	S	0	0	0
			1877	1187	344	337	9			
2	H	238	Total	C	N	O	S	0	0	0
			1877	1187	344	337	9			
2	J	238	Total	C	N	O	S	0	0	0
			1877	1187	344	337	9			
2	L	238	Total	C	N	O	S	0	0	0
			1877	1187	344	337	9			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	408	HIS	TYR	ENGINEERED	UNP P00437

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Chain	Residue	Modelled	Actual	Comment	Reference
B	429	CME	CYS	MODIFIED RESIDUE	UNP P00437
D	408	HIS	TYR	ENGINEERED	UNP P00437
D	429	CME	CYS	MODIFIED RESIDUE	UNP P00437
F	408	HIS	TYR	ENGINEERED	UNP P00437
F	429	CME	CYS	MODIFIED RESIDUE	UNP P00437
H	408	HIS	TYR	ENGINEERED	UNP P00437
H	429	CME	CYS	MODIFIED RESIDUE	UNP P00437
J	408	HIS	TYR	ENGINEERED	UNP P00437
J	429	CME	CYS	MODIFIED RESIDUE	UNP P00437
L	408	HIS	TYR	ENGINEERED	UNP P00437
L	429	CME	CYS	MODIFIED RESIDUE	UNP P00437

- Molecule 3 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	J	1	Total Fe 1 1	0	0
3	D	1	Total Fe 1 1	0	0
3	H	1	Total Fe 1 1	0	0
3	B	1	Total Fe 1 1	0	0
3	L	1	Total Fe 1 1	0	0
3	F	1	Total Fe 1 1	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	47	Total O 47 47	0	0
4	B	77	Total O 77 77	0	0
4	C	44	Total O 44 44	0	0
4	D	83	Total O 83 83	0	0
4	E	45	Total O 45 45	0	0
4	F	78	Total O 78 78	0	0

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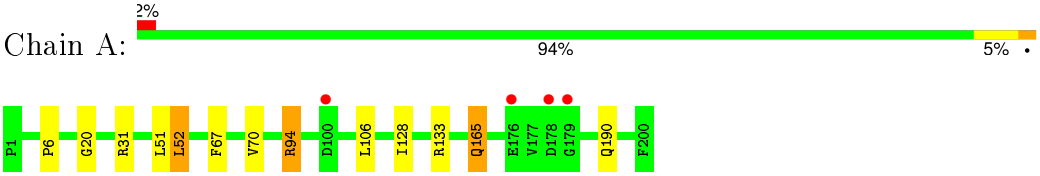
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	G	45	Total O 45 45	0	0
4	H	81	Total O 81 81	0	0
4	I	43	Total O 43 43	0	0
4	J	83	Total O 83 83	0	0
4	K	46	Total O 46 46	0	0
4	L	78	Total O 78 78	0	0

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.

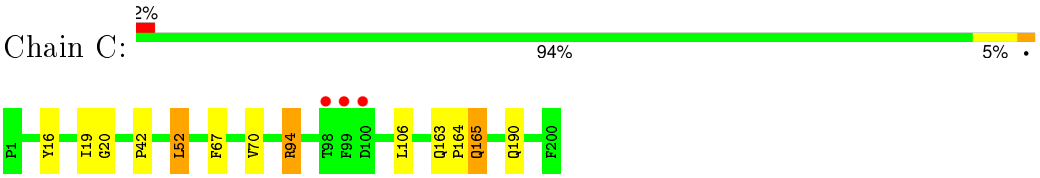
- Molecule 1: Protocatechuate 3,4-dioxygenase alpha chain

Chain A:



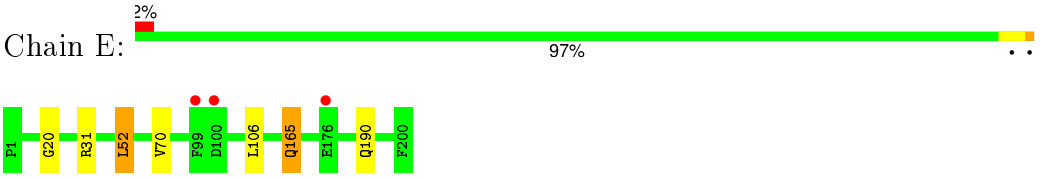
- Molecule 1: Protocatechuate 3,4-dioxygenase alpha chain

Chain C:



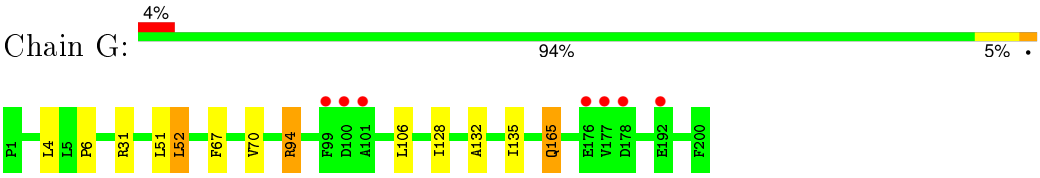
- Molecule 1: Protocatechuate 3,4-dioxygenase alpha chain

Chain E:



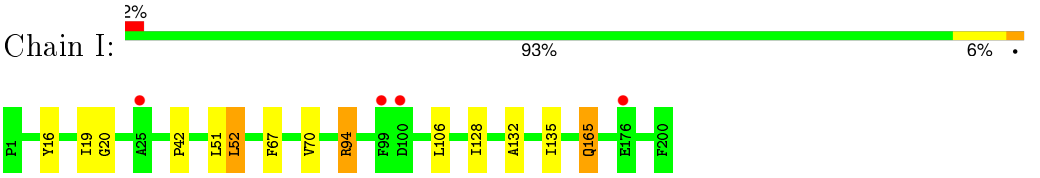
- Molecule 1: Protocatechuate 3,4-dioxygenase alpha chain

Chain G:



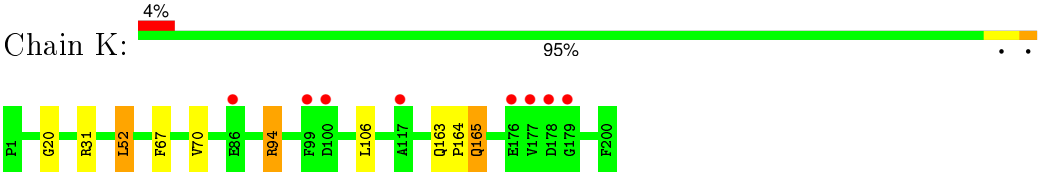
- Molecule 1: Protocatechuate 3,4-dioxygenase alpha chain

Chain I:



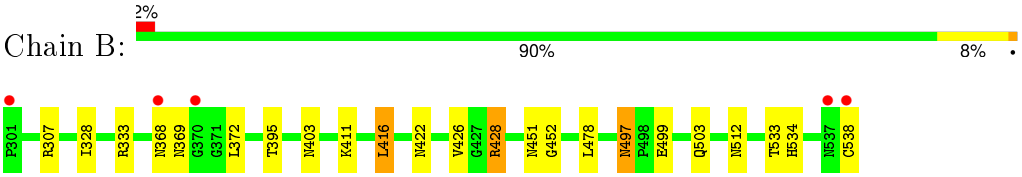
- Molecule 1: Protocatechuate 3,4-dioxygenase alpha chain

Chain K:



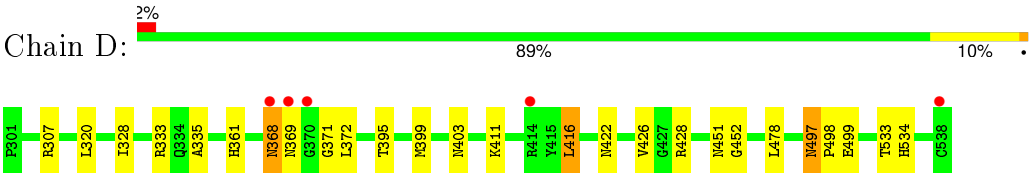
• Molecule 2: Protocatechuate 3,4-dioxygenase beta chain

Chain B:



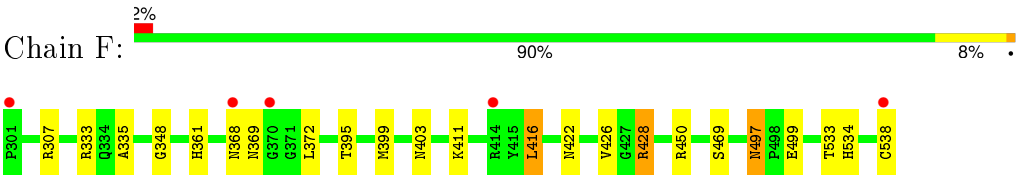
• Molecule 2: Protocatechuate 3,4-dioxygenase beta chain

Chain D:



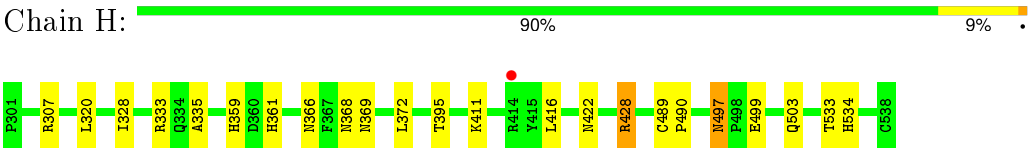
• Molecule 2: Protocatechuate 3,4-dioxygenase beta chain

Chain F:



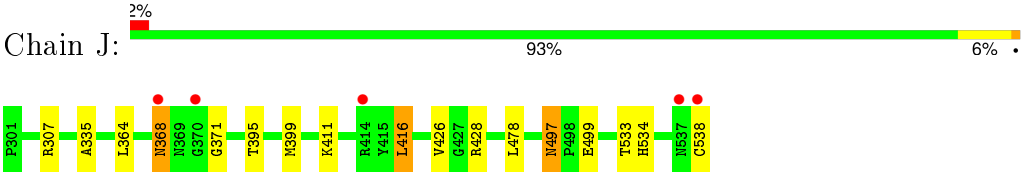
• Molecule 2: Protocatechuate 3,4-dioxygenase beta chain

Chain H:



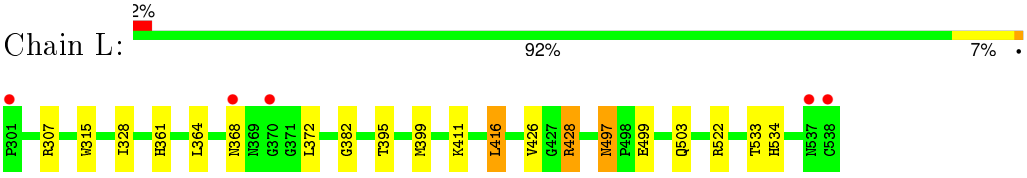
• Molecule 2: Protocatechuate 3,4-dioxygenase beta chain

Chain J:



• Molecule 2: Protocatechuate 3,4-dioxygenase beta chain

Chain L:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	196.59Å 127.63Å 134.45Å 90.00° 97.60° 90.00°	Depositor
Resolution (Å)	37.27 – 2.54 37.27 – 1.95	Depositor EDS
% Data completeness (in resolution range)	76.1 (37.27-2.54) 76.2 (37.27-1.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	0.00 (at 1.95Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.183 , 0.217 0.180 , 0.214	Depositor DCC
R_{free} test set	1053 reflections (1.29%)	DCC
Wilson B-factor (Å^2)	24.0	Xtriage
Anisotropy	0.398	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.26 , 23.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Outliers	0 of 189293 reflections	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	21444	wwPDB-VP
Average B, all atoms (Å^2)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.87% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: CME, 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	0.62	0/1611	0.75	1/2195 (0.0%)
1	C	0.61	0/1611	0.76	1/2195 (0.0%)
1	E	0.63	0/1611	0.74	1/2195 (0.0%)
1	G	0.61	0/1611	0.74	1/2195 (0.0%)
1	I	0.64	0/1611	0.74	1/2195 (0.0%)
1	K	0.66	0/1611	0.73	1/2195 (0.0%)
2	B	0.61	0/1922	0.77	1/2615 (0.0%)
2	D	0.61	0/1922	0.78	0/2615
2	F	0.61	0/1922	0.76	0/2615
2	H	0.62	0/1922	0.76	0/2615
2	J	0.63	0/1922	0.76	0/2615
2	L	0.64	0/1922	0.77	0/2615
All	All	0.62	0/21198	0.76	7/28860 (0.0%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	52	LEU	CA-CB-CG	7.37	132.24	115.30
1	C	52	LEU	CA-CB-CG	7.27	132.01	115.30
1	G	52	LEU	CA-CB-CG	7.10	131.62	115.30
1	E	52	LEU	CA-CB-CG	7.00	131.40	115.30
1	K	52	LEU	CA-CB-CG	6.95	131.28	115.30
1	I	52	LEU	CA-CB-CG	6.78	130.90	115.30
2	B	478	LEU	CA-CB-CG	5.05	126.91	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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	1571	0	1499	10	0
1	C	1571	0	1499	8	0
1	E	1571	0	1499	5	0
1	G	1571	0	1499	11	0
1	I	1571	0	1499	8	0
1	K	1571	0	1499	7	0
2	B	1877	0	1821	12	0
2	D	1877	0	1821	14	0
2	F	1877	0	1821	13	0
2	H	1877	0	1821	14	0
2	J	1877	0	1821	6	0
2	L	1877	0	1821	10	0
3	B	1	0	0	0	0
3	D	1	0	0	0	0
3	F	1	0	0	0	0
3	H	1	0	0	0	0
3	J	1	0	0	0	0
3	L	1	0	0	0	0
4	A	47	0	0	0	0
4	B	77	0	0	1	0
4	C	44	0	0	0	0
4	D	83	0	0	1	0
4	E	45	0	0	0	0
4	F	78	0	0	1	0
4	G	45	0	0	0	0
4	H	81	0	0	0	0
4	I	43	0	0	0	0
4	J	83	0	0	0	0
4	K	46	0	0	0	0
4	L	78	0	0	0	0
All	All	21444	0	19920	99	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 2.

All (99) 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:C:165:GLN:H	1:C:165:GLN:HE21	1.29	0.81
1:I:165:GLN:H	1:I:165:GLN:HE21	1.41	0.66
2:B:497:ASN:HD22	2:B:499:GLU:H	1.45	0.62
2:F:497:ASN:HD22	2:F:499:GLU:H	1.48	0.60
2:L:497:ASN:HD22	2:L:499:GLU:H	1.49	0.59
1:I:165:GLN:H	1:I:165:GLN:NE2	2.00	0.59
2:H:369:ASN:H	2:H:422:ASN:HD22	1.50	0.59
1:K:165:GLN:NE2	1:K:165:GLN:H	2.01	0.59
1:E:165:GLN:NE2	1:E:165:GLN:H	2.01	0.59
2:H:497:ASN:HD22	2:H:499:GLU:H	1.51	0.58
2:B:307:ARG:HG2	2:B:533:THR:HG22	1.86	0.58
2:D:307:ARG:HG2	2:D:533:THR:HG22	1.86	0.57
1:C:165:GLN:H	1:C:165:GLN:NE2	2.01	0.57
1:E:31:ARG:NH1	2:F:428:ARG:HG2	2.20	0.57
1:K:20:GLY:HA2	2:L:426:VAL:HG13	1.87	0.57
1:I:51:LEU:HD12	1:I:106:LEU:HD23	1.87	0.56
1:A:6:PRO:HB2	2:B:503:GLN:HE22	1.70	0.56
2:J:497:ASN:HD22	2:J:499:GLU:H	1.53	0.55
1:G:6:PRO:HB2	2:H:503:GLN:HE22	1.74	0.53
2:D:368:ASN:ND2	2:D:371:GLY:H	2.05	0.53
2:J:416:LEU:H	2:J:416:LEU:HD23	1.74	0.53
1:G:165:GLN:NE2	1:G:165:GLN:H	2.05	0.53
1:I:67:PHE:HZ	1:I:94:ARG:HD2	1.74	0.52
1:I:20:GLY:HA2	2:J:426:VAL:HG13	1.90	0.52
2:D:416:LEU:H	2:D:416:LEU:HD23	1.74	0.52
2:F:416:LEU:HD23	2:F:416:LEU:H	1.75	0.52
1:K:67:PHE:HZ	1:K:94:ARG:HD2	1.74	0.52
1:A:165:GLN:H	1:A:165:GLN:NE2	2.07	0.51
2:D:369:ASN:H	2:D:422:ASN:HD22	1.59	0.51
1:G:31:ARG:NH1	2:H:428:ARG:HG2	2.25	0.50
2:H:335:ALA:HB2	2:L:328:ILE:HD12	1.94	0.50
2:L:307:ARG:HG2	2:L:533:THR:HG22	1.92	0.50
1:A:20:GLY:HA2	2:B:426:VAL:HG13	1.94	0.50
1:C:67:PHE:HZ	1:C:94:ARG:HD2	1.77	0.50
2:L:416:LEU:H	2:L:416:LEU:HD23	1.76	0.49
2:B:369:ASN:H	2:B:422:ASN:HD22	1.58	0.49
1:A:67:PHE:HZ	1:A:94:ARG:HD2	1.78	0.49
1:C:20:GLY:HA2	2:D:426:VAL:HG13	1.94	0.49
2:B:416:LEU:H	2:B:416:LEU:HD23	1.77	0.49
2:B:328:ILE:HD12	2:D:335:ALA:HB2	1.95	0.48
2:H:369:ASN:H	2:H:422:ASN:ND2	2.12	0.48
2:B:497:ASN:ND2	2:B:499:GLU:H	2.11	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:190:GLN:HG3	2:F:333:ARG:HG2	1.96	0.48
1:E:20:GLY:HA2	2:F:426:VAL:HG13	1.96	0.47
2:H:307:ARG:HG2	2:H:533:THR:HG22	1.95	0.47
1:G:165:GLN:HE21	1:G:165:GLN:H	1.63	0.47
1:K:67:PHE:CZ	1:K:94:ARG:HD2	2.48	0.47
1:G:67:PHE:HZ	1:G:94:ARG:HD2	1.78	0.46
1:A:70:VAL:HG21	1:A:106:LEU:HD21	1.98	0.46
1:E:70:VAL:HG21	1:E:106:LEU:HD21	1.98	0.45
2:J:307:ARG:HG2	2:J:533:THR:HG22	1.98	0.45
2:F:307:ARG:HG2	2:F:533:THR:HG22	1.99	0.45
2:F:497:ASN:ND2	2:F:499:GLU:H	2.14	0.45
2:F:403:ASN:HB2	4:F:2620:HOH:O	2.15	0.44
1:G:132:ALA:HB3	1:G:135:ILE:HD12	1.99	0.44
2:D:320:LEU:HD13	2:D:333:ARG:HH12	1.81	0.44
1:G:51:LEU:HD12	1:G:106:LEU:HD23	1.98	0.44
1:K:31:ARG:NH1	2:L:428:ARG:HG2	2.33	0.44
1:K:70:VAL:HG21	1:K:106:LEU:HD21	2.00	0.44
1:I:70:VAL:HG12	1:I:128:ILE:HG12	2.00	0.44
1:A:70:VAL:HG12	1:A:128:ILE:HG12	2.00	0.44
2:F:361:HIS:H	2:F:361:HIS:CD2	2.36	0.43
1:C:190:GLN:HG3	2:D:333:ARG:HG2	2.01	0.43
1:A:31:ARG:NH1	2:B:428:ARG:HG2	2.34	0.43
1:I:16:TYR:O	1:I:19:ILE:HG12	2.18	0.43
2:H:489:CYS:HA	2:H:490:PRO:HD3	1.92	0.43
2:F:369:ASN:H	2:F:422:ASN:HD22	1.66	0.43
1:C:16:TYR:O	1:C:19:ILE:HG12	2.17	0.43
2:H:416:LEU:H	2:H:416:LEU:HD23	1.83	0.43
2:B:403:ASN:HB2	4:B:620:HOH:O	2.19	0.43
1:G:6:PRO:HB2	2:H:503:GLN:NE2	2.34	0.43
1:A:67:PHE:CZ	1:A:94:ARG:HD2	2.54	0.43
2:L:497:ASN:ND2	2:L:499:GLU:H	2.15	0.43
2:L:315:TRP:HZ2	2:L:503:GLN:HE21	1.66	0.42
2:H:359:HIS:O	2:H:366:ASN:HB3	2.19	0.42
1:G:67:PHE:CZ	1:G:94:ARG:HD2	2.54	0.42
2:J:368:ASN:ND2	2:J:371:GLY:H	2.16	0.42
1:K:163:GLN:HA	1:K:164:PRO:HD3	1.91	0.42
2:D:497:ASN:ND2	2:D:499:GLU:H	2.18	0.41
2:H:320:LEU:HD13	2:H:333:ARG:HH22	1.85	0.41
1:G:70:VAL:HG21	1:G:106:LEU:HD21	2.01	0.41
1:C:163:GLN:HA	1:C:164:PRO:HD3	1.92	0.41
2:F:348:GLY:HA3	2:F:469:SER:HA	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:361:HIS:CD2	2:D:361:HIS:H	2.37	0.41
2:H:361:HIS:H	2:H:361:HIS:CD2	2.38	0.41
2:D:497:ASN:HD22	2:D:498:PRO:N	2.18	0.41
2:D:403:ASN:HB2	4:D:1620:HOH:O	2.20	0.41
1:C:70:VAL:HG21	1:C:106:LEU:HD21	2.01	0.41
2:L:382:GLY:HA2	2:L:522:ARG:HE	1.85	0.41
2:L:361:HIS:CD2	2:L:361:HIS:H	2.39	0.41
1:G:70:VAL:HG12	1:G:128:ILE:HG12	2.02	0.41
1:A:190:GLN:HG3	2:B:333:ARG:HG2	2.01	0.41
2:B:451:ASN:HB3	2:B:452:GLY:H	1.76	0.41
2:F:333:ARG:HE	2:F:333:ARG:HA	1.86	0.40
2:D:451:ASN:HB3	2:D:452:GLY:H	1.71	0.40
1:A:51:LEU:HD12	1:A:106:LEU:HD23	2.03	0.40
1:I:132:ALA:HB3	1:I:135:ILE:HD12	2.03	0.40
2:D:328:ILE:HD12	2:F:335:ALA:HB2	2.03	0.40
2:H:328:ILE:HD12	2:J:335:ALA:HB2	2.03	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	198/200 (99%)	194 (98%)	4 (2%)	0	100	100
1	C	198/200 (99%)	191 (96%)	7 (4%)	0	100	100
1	E	198/200 (99%)	192 (97%)	6 (3%)	0	100	100
1	G	198/200 (99%)	192 (97%)	6 (3%)	0	100	100
1	I	198/200 (99%)	193 (98%)	5 (2%)	0	100	100
1	K	198/200 (99%)	192 (97%)	6 (3%)	0	100	100
2	B	235/238 (99%)	227 (97%)	7 (3%)	1 (0%)	39	60

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
2	D	235/238 (99%)	229 (97%)	5 (2%)	1 (0%)	39 60
2	F	235/238 (99%)	228 (97%)	6 (3%)	1 (0%)	39 60
2	H	235/238 (99%)	225 (96%)	9 (4%)	1 (0%)	39 60
2	J	235/238 (99%)	230 (98%)	4 (2%)	1 (0%)	39 60
2	L	235/238 (99%)	226 (96%)	8 (3%)	1 (0%)	39 60
All	All	2598/2628 (99%)	2519 (97%)	73 (3%)	6 (0%)	52 73

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	368	ASN
2	D	368	ASN
2	F	368	ASN
2	H	368	ASN
2	J	368	ASN
2	L	368	ASN

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	162/163 (99%)	158 (98%)	4 (2%)	55 80
1	C	162/163 (99%)	158 (98%)	4 (2%)	55 80
1	E	162/163 (99%)	160 (99%)	2 (1%)	78 92
1	G	162/163 (99%)	158 (98%)	4 (2%)	55 80
1	I	162/163 (99%)	158 (98%)	4 (2%)	55 80
1	K	162/163 (99%)	159 (98%)	3 (2%)	65 86
2	B	199/201 (99%)	190 (96%)	9 (4%)	34 56
2	D	199/201 (99%)	190 (96%)	9 (4%)	34 56
2	F	199/201 (99%)	189 (95%)	10 (5%)	30 51
2	H	199/201 (99%)	193 (97%)	6 (3%)	48 74

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
2	J	199/201 (99%)	189 (95%)	10 (5%)	30 51
2	L	199/201 (99%)	190 (96%)	9 (4%)	34 56
All	All	2166/2184 (99%)	2092 (97%)	74 (3%)	44 70

All (74) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	52	LEU
1	A	94	ARG
1	A	133	ARG
1	A	165	GLN
2	B	372	LEU
2	B	395	THR
2	B	411	LYS
2	B	416	LEU
2	B	428	ARG
2	B	497	ASN
2	B	512	ASN
2	B	534	HIS
2	B	538	CYS
1	C	42	PRO
1	C	52	LEU
1	C	94	ARG
1	C	165	GLN
2	D	372	LEU
2	D	395	THR
2	D	399	MET
2	D	411	LYS
2	D	416	LEU
2	D	428	ARG
2	D	478	LEU
2	D	497	ASN
2	D	534	HIS
1	E	52	LEU
1	E	165	GLN
2	F	372	LEU
2	F	395	THR
2	F	399	MET
2	F	411	LYS
2	F	416	LEU
2	F	428	ARG

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Mol	Chain	Res	Type
2	F	450	ARG
2	F	497	ASN
2	F	534	HIS
2	F	538	CYS
1	G	4	LEU
1	G	52	LEU
1	G	94	ARG
1	G	165	GLN
2	H	372	LEU
2	H	395	THR
2	H	411	LYS
2	H	428	ARG
2	H	497	ASN
2	H	534	HIS
1	I	42	PRO
1	I	52	LEU
1	I	94	ARG
1	I	165	GLN
2	J	364	LEU
2	J	395	THR
2	J	399	MET
2	J	411	LYS
2	J	416	LEU
2	J	428	ARG
2	J	478	LEU
2	J	497	ASN
2	J	534	HIS
2	J	538	CYS
1	K	52	LEU
1	K	94	ARG
1	K	165	GLN
2	L	364	LEU
2	L	372	LEU
2	L	395	THR
2	L	399	MET
2	L	411	LYS
2	L	416	LEU
2	L	428	ARG
2	L	497	ASN
2	L	534	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (37) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	165	GLN
2	B	361	HIS
2	B	412	ASN
2	B	422	ASN
2	B	497	ASN
2	B	503	GLN
1	C	165	GLN
2	D	361	HIS
2	D	368	ASN
2	D	412	ASN
2	D	422	ASN
2	D	497	ASN
2	D	503	GLN
1	E	165	GLN
2	F	361	HIS
2	F	412	ASN
2	F	422	ASN
2	F	497	ASN
2	F	503	GLN
1	G	165	GLN
2	H	361	HIS
2	H	412	ASN
2	H	422	ASN
2	H	497	ASN
2	H	503	GLN
1	I	165	GLN
2	J	361	HIS
2	J	368	ASN
2	J	412	ASN
2	J	422	ASN
2	J	497	ASN
1	K	165	GLN
2	L	361	HIS
2	L	412	ASN
2	L	422	ASN
2	L	497	ASN
2	L	503	GLN

5.3.3 RNA

There are no RNA molecules in this entry.

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

6 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	$\#\|Z\|>2$
2	CME	B	429	2	8,9,10	0.90	1 (12%)	6,9,11	1.44	1 (16%)
2	CME	D	429	2	8,9,10	0.86	1 (12%)	6,9,11	1.30	1 (16%)
2	CME	F	429	2	8,9,10	0.92	1 (12%)	6,9,11	1.22	1 (16%)
2	CME	H	429	2	8,9,10	0.87	1 (12%)	6,9,11	1.47	1 (16%)
2	CME	J	429	2	8,9,10	0.90	1 (12%)	6,9,11	1.44	1 (16%)
2	CME	L	429	2	8,9,10	0.88	1 (12%)	6,9,11	1.35	1 (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
2	CME	B	429	2	-	0/5/8/10	0/0/0/0
2	CME	D	429	2	-	0/5/8/10	0/0/0/0
2	CME	F	429	2	-	0/5/8/10	0/0/0/0
2	CME	H	429	2	-	0/5/8/10	0/0/0/0
2	CME	J	429	2	-	0/5/8/10	0/0/0/0
2	CME	L	429	2	-	0/5/8/10	0/0/0/0

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	429	CME	OH-CZ	-2.21	1.30	1.42
2	D	429	CME	OH-CZ	-2.20	1.30	1.42
2	J	429	CME	OH-CZ	-2.20	1.30	1.42
2	L	429	CME	OH-CZ	-2.19	1.30	1.42
2	F	429	CME	OH-CZ	-2.19	1.30	1.42
2	H	429	CME	OH-CZ	-2.13	1.30	1.42

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	F	429	CME	OH-CZ-CE	2.18	120.37	110.83
2	D	429	CME	OH-CZ-CE	2.29	120.87	110.83
2	L	429	CME	OH-CZ-CE	2.40	121.34	110.83
2	H	429	CME	OH-CZ-CE	2.45	121.57	110.83
2	B	429	CME	OH-CZ-CE	2.48	121.71	110.83
2	J	429	CME	OH-CZ-CE	2.53	121.92	110.83

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

Of 6 ligands modelled in this entry, 6 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

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^{th} 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	200/200 (100%)	-0.49	4 (2%)	68	73	8, 27, 54, 72	0
1	C	200/200 (100%)	-0.54	3 (1%)	76	80	8, 28, 54, 71	0
1	E	200/200 (100%)	-0.52	3 (1%)	76	80	8, 28, 54, 72	0
1	G	200/200 (100%)	-0.53	7 (3%)	48	54	9, 28, 54, 73	0
1	I	200/200 (100%)	-0.50	4 (2%)	68	73	10, 30, 55, 73	0
1	K	200/200 (100%)	-0.23	8 (4%)	42	48	11, 31, 55, 73	0
2	B	237/238 (99%)	-0.70	5 (2%)	67	72	9, 19, 51, 71	0
2	D	237/238 (99%)	-0.67	5 (2%)	67	72	8, 19, 51, 71	0
2	F	237/238 (99%)	-0.69	5 (2%)	67	72	9, 20, 52, 69	0
2	H	237/238 (99%)	-0.79	1 (0%)	93	94	9, 19, 51, 69	0
2	J	237/238 (99%)	-0.75	5 (2%)	67	72	10, 21, 52, 70	0
2	L	237/238 (99%)	-0.67	5 (2%)	67	72	11, 21, 52, 71	0
All	All	2622/2628 (99%)	-0.60	55 (2%)	67	72	8, 24, 54, 73	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	L	368	ASN	5.5
2	B	301	PRO	5.1
2	L	538	CYS	5.0
2	J	370	GLY	4.9
2	B	370	GLY	4.7
2	L	370	GLY	4.6
1	G	100	ASP	4.4
2	B	368	ASN	4.3
2	D	370	GLY	4.3
1	G	99	PHE	4.2
2	F	370	GLY	4.1

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Mol	Chain	Res	Type	RSRZ
1	I	99	PHE	4.1
2	F	368	ASN	3.8
1	E	99	PHE	3.4
1	C	100	ASP	3.4
2	D	368	ASN	3.4
1	K	177	VAL	3.3
1	A	178	ASP	3.3
1	A	176	GLU	3.3
1	E	100	ASP	3.2
1	K	100	ASP	3.2
1	I	100	ASP	3.2
1	A	179	GLY	3.1
2	B	538	CYS	3.1
1	K	99	PHE	3.1
1	C	98	THR	3.1
1	G	177	VAL	3.0
1	K	86	GLU	2.9
2	L	301	PRO	2.9
1	I	176	GLU	2.9
1	C	99	PHE	2.9
1	I	25	ALA	2.8
2	B	537	ASN	2.7
1	K	178	ASP	2.7
2	J	368	ASN	2.7
2	D	538	CYS	2.7
1	K	179	GLY	2.7
1	A	100	ASP	2.6
2	D	414	ARG	2.6
2	F	414	ARG	2.5
2	H	414	ARG	2.5
2	J	537	ASN	2.4
1	G	101	ALA	2.4
1	K	117	ALA	2.3
2	J	414	ARG	2.2
1	K	176	GLU	2.2
2	D	369	ASN	2.2
2	J	538	CYS	2.1
1	E	176	GLU	2.1
1	G	192	GLU	2.1
2	F	538	CYS	2.0
1	G	178	ASP	2.0
1	G	176	GLU	2.0

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Mol	Chain	Res	Type	RSRZ
2	F	301	PRO	2.0
2	L	537	ASN	2.0

6.2 Non-standard residues in protein, DNA, RNA chains (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^{th} 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^2)	Q<0.9
2	CME	F	429	10/11	0.95	0.12	-	19,22,37,41	0
2	CME	D	429	10/11	0.96	0.07	-	19,24,36,38	0
2	CME	B	429	10/11	0.96	0.09	-	18,22,35,37	0
2	CME	L	429	10/11	0.96	0.09	-	22,25,37,39	0
2	CME	J	429	10/11	0.94	0.09	-	23,25,37,39	0
2	CME	H	429	10/11	0.96	0.09	-	19,22,36,38	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^{th} 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^2)	Q<0.9
3	FE	B	600	1/1	0.97	0.11	0.83	41,41,41,41	0
3	FE	L	600	1/1	0.99	0.07	-1.44	37,37,37,37	0
3	FE	D	600	1/1	0.99	0.09	-1.49	40,40,40,40	0
3	FE	F	600	1/1	0.98	0.07	-1.82	42,42,42,42	0
3	FE	H	600	1/1	0.99	0.06	-3.57	40,40,40,40	0
3	FE	J	600	1/1	0.99	0.07	-	51,51,51,51	0

6.5 Other polymers (i)

There are no such residues in this entry.