



Full wwPDB X-ray Structure Validation Report ⓘ

May 24, 2020 – 04:55 pm BST

PDB ID : 1EMS
Title : CRYSTAL STRUCTURE OF THE C. ELEGANS NITFHIT PROTEIN
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Deposited on : 2000-03-17
Resolution : 2.80 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

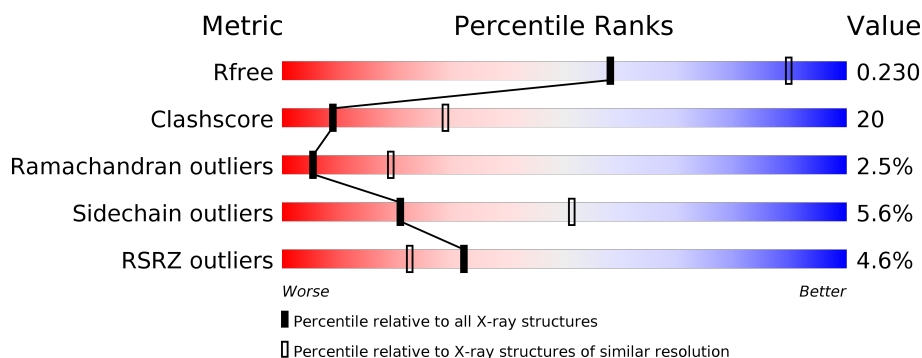
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 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	440	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, green 62%, yellow 27%, orange 2%, grey 6%);"></div> <div style="display: flex; justify-content: space-between; padding: 0 5px;"> % 62% 27% • • 6% </div> </div>
1	B	440	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 7%, green 53%, yellow 35%, orange 5%, grey 7%);"></div> <div style="display: flex; justify-content: space-between; padding: 0 5px;"> 7% 53% 35% • 7% </div> </div>

2 Entry composition [i](#)

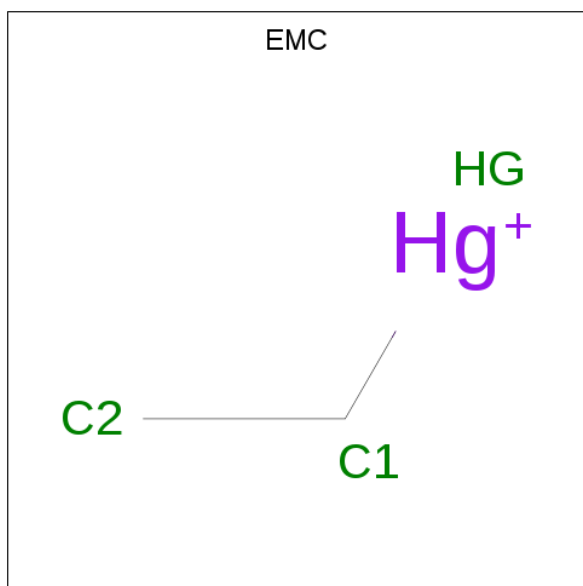
There are 5 unique types of molecules in this entry. The entry contains 6708 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 NIT-FRAGILE HISTIDINE TRIAD FUSION PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	412	Total	C	N	O	S	0	0	0
			3274	2066	581	604	23			
1	B	410	Total	C	N	O	S	0	0	0
			3246	2044	579	600	23			

- Molecule 2 is ETHYL MERCURY ION (three-letter code: EMC) (formula: $\text{C}_2\text{H}_5\text{Hg}$).



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

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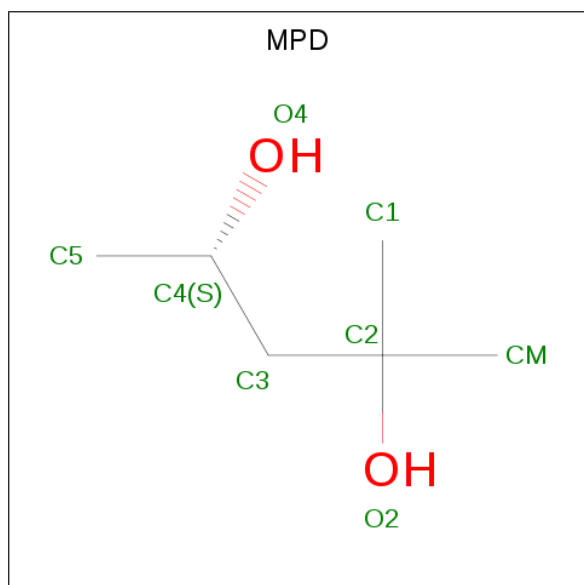
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	C	Hg	0	0
			3	2	1		

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Na	0	0
			1	1		
3	A	3	Total	Na	0	0
			3	3		

- Molecule 4 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: C₆H₁₄O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			8	6	2		

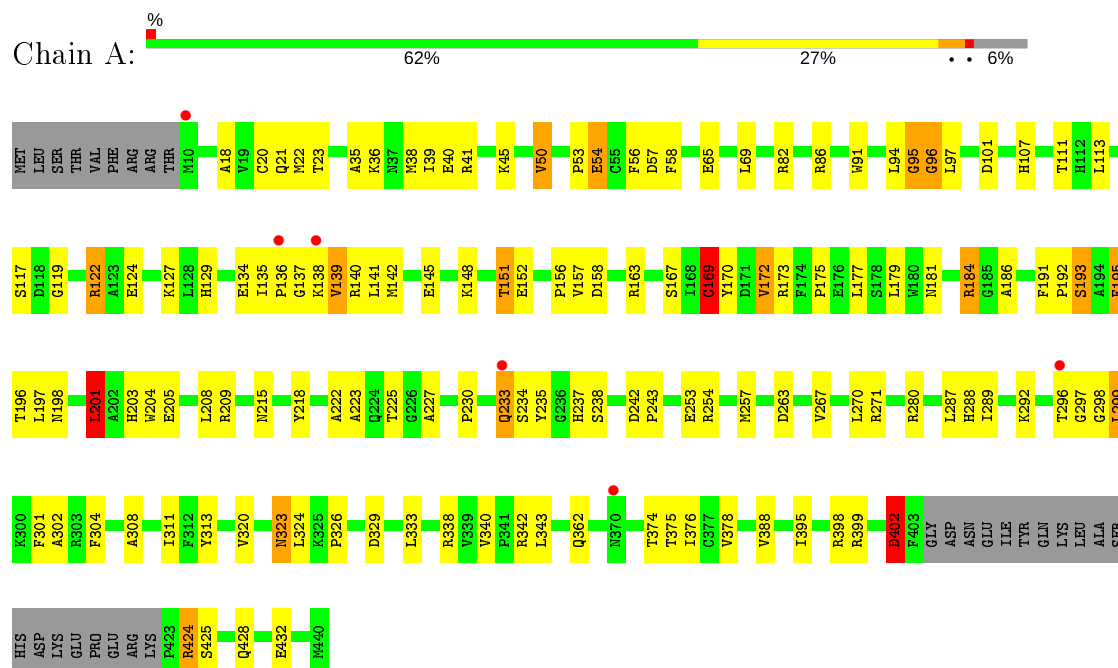
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	100	Total	O	0	0
			100	100		
5	B	61	Total	O	0	0
			61	61		

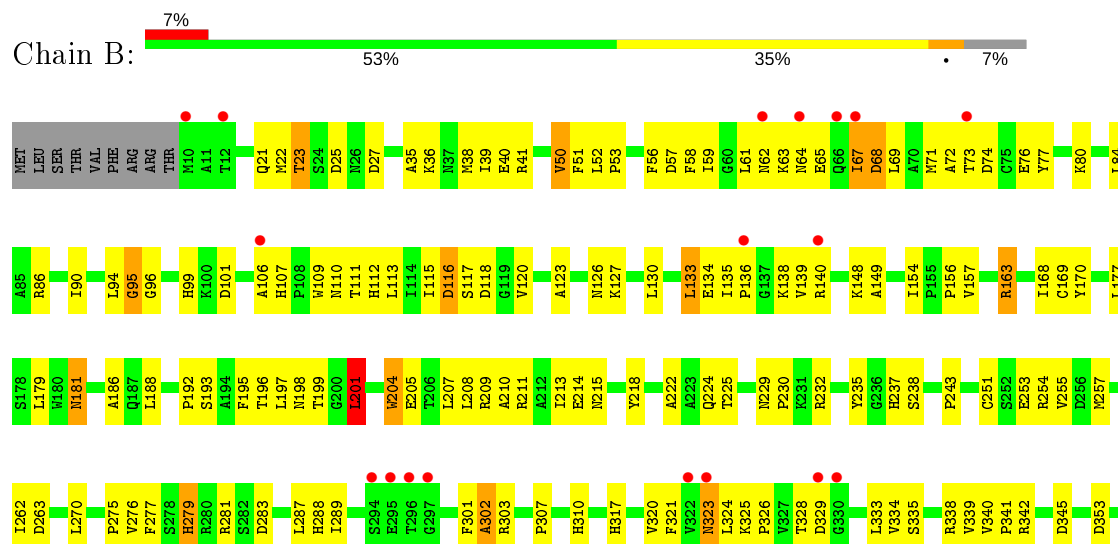
3 Residue-property plots [i](#)

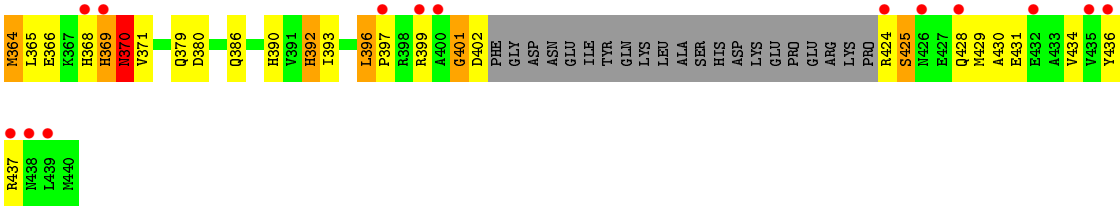
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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: NIT-FRAGILE HISTIDINE TRIAD FUSION PROTEIN



• Molecule 1: NIT-FRAGILE HISTIDINE TRIAD FUSION PROTEIN





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	68.75Å 100.44Å 158.65Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.80 39.66 – 2.80	Depositor EDS
% Data completeness (in resolution range)	94.8 (30.00-2.80) 96.6 (39.66-2.80)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	8.28 (at 2.81Å)	Xtriage
Refinement program	CNS 0.9	Depositor
R, R_{free}	0.190 , 0.231 0.192 , 0.230	Depositor DCC
R_{free} test set	3344 reflections (6.70%)	wwPDB-VP
Wilson B-factor (Å ²)	27.1	Xtriage
Anisotropy	0.489	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 54.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6708	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.56% 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: NA, MPD, EMC

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	0.38	0/3351	0.69	3/4537 (0.1%)
1	B	0.36	0/3320	0.64	1/4496 (0.0%)
All	All	0.37	0/6671	0.67	4/9033 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	201	LEU	CA-CB-CG	6.03	129.18	115.30
1	B	201	LEU	CA-CB-CG	5.62	128.22	115.30
1	A	169	CYS	CA-CB-SG	5.21	123.39	114.00
1	A	95	GLY	N-CA-C	5.09	125.84	113.10

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	3274	0	3210	112	0
1	B	3246	0	3184	152	0
2	A	9	0	0	0	0
2	B	6	0	0	0	0
3	A	3	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	1	0	0	0	0
4	A	8	0	14	0	0
5	A	100	0	0	3	0
5	B	61	0	0	1	0
All	All	6708	0	6408	262	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 20.

All (262) 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:169:CYS:HB3	1:A:193:SER:CB	1.67	1.23
1:A:169:CYS:HB3	1:A:193:SER:HB3	1.30	1.09
1:B:329:ASP:HB2	1:B:399:ARG:HH22	1.20	1.05
1:A:54:GLU:HB2	1:A:193:SER:HA	1.40	1.03
1:B:67:ILE:HD11	1:B:106:ALA:HA	1.43	0.98
1:B:329:ASP:HB2	1:B:399:ARG:NH2	1.82	0.94
1:A:169:CYS:HB3	1:A:193:SER:HB2	1.50	0.93
1:B:135:ILE:HG21	1:B:138:LYS:HD2	1.54	0.87
1:A:156:PRO:HD3	1:A:181:ASN:HD22	1.39	0.86
1:A:169:CYS:CB	1:A:193:SER:HB3	2.05	0.86
1:A:243:PRO:HB2	1:A:270:LEU:HD12	1.58	0.85
1:B:116:ASP:HB3	1:B:118:ASP:H	1.44	0.83
1:B:323:ASN:HD22	1:B:324:LEU:H	1.27	0.83
1:B:38:MET:HA	1:B:257:MET:HE1	1.63	0.78
1:A:135:ILE:HB	1:A:139:VAL:HG23	1.67	0.76
1:B:99:HIS:HB3	1:B:109:TRP:HB2	1.66	0.76
1:A:38:MET:HA	1:A:257:MET:HE1	1.67	0.76
1:A:169:CYS:O	1:A:172:VAL:HG22	1.87	0.75
1:B:23:THR:HB	1:B:225:THR:O	1.87	0.75
1:B:324:LEU:HD22	1:B:424:ARG:HB2	1.71	0.72
1:B:25:ASP:HB3	1:B:27:ASP:H	1.56	0.71
1:A:21:GLN:NE2	1:A:238:SER:H	1.89	0.71
1:B:243:PRO:HB2	1:B:270:LEU:HD12	1.73	0.71
1:A:184:ARG:HB3	1:A:184:ARG:HH11	1.57	0.70
1:B:364:MET:HE2	1:B:365:LEU:HA	1.74	0.69
1:B:72:ALA:O	1:B:73:THR:HB	1.93	0.69
1:B:116:ASP:HB2	1:B:120:VAL:H	1.55	0.68
1:B:134:GLU:HG3	1:B:140:ARG:HG3	1.73	0.68
1:B:329:ASP:CB	1:B:399:ARG:HH22	2.02	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:430:ALA:O	1:B:434:VAL:HG13	1.93	0.68
1:B:323:ASN:ND2	1:B:324:LEU:H	1.90	0.68
1:A:172:VAL:HG13	1:A:191:PHE:CD2	2.29	0.68
1:A:184:ARG:HB3	1:A:184:ARG:NH1	2.09	0.67
1:B:112:HIS:HB2	1:B:168:ILE:HD11	1.75	0.67
1:A:36:LYS:O	1:A:40:GLU:HG3	1.95	0.67
1:B:425:SER:OG	1:B:428:GLN:HB3	1.94	0.66
1:B:41:ARG:NE	1:B:257:MET:HE3	2.10	0.65
1:B:71:MET:SD	1:B:77:TYR:HD2	2.19	0.65
1:B:323:ASN:HD22	1:B:324:LEU:N	1.93	0.64
1:B:287:LEU:HG	1:B:289:ILE:CD1	2.27	0.64
1:B:364:MET:HG3	1:B:365:LEU:N	2.11	0.64
1:A:399:ARG:O	1:A:402:ASP:HB2	1.99	0.62
1:B:133:LEU:CD2	1:B:135:ILE:HG13	2.29	0.62
1:B:323:ASN:ND2	1:B:324:LEU:N	2.48	0.62
1:A:297:GLY:HA2	1:A:308:ALA:HB2	1.83	0.61
1:B:23:THR:HG22	1:B:255:VAL:CG2	2.30	0.61
1:B:287:LEU:HG	1:B:289:ILE:HD11	1.81	0.61
1:B:379:GLN:O	1:B:386:GLN:HB2	2.01	0.61
1:B:116:ASP:HB2	1:B:120:VAL:N	2.16	0.61
1:A:18:ALA:HB3	1:A:50:VAL:HG12	1.83	0.60
1:B:396:LEU:HD23	1:B:396:LEU:N	2.15	0.60
1:A:288:HIS:C	1:A:289:ILE:HD12	2.20	0.60
1:B:107:HIS:HB3	1:B:148:LYS:HB2	1.83	0.60
1:B:138:LYS:O	1:B:139:VAL:HG23	2.02	0.60
1:B:340:VAL:CG1	1:B:345:ASP:HB2	2.32	0.60
1:A:151:THR:HG22	1:A:152:GLU:HG3	1.82	0.60
1:A:156:PRO:HD3	1:A:181:ASN:ND2	2.12	0.60
1:A:280:ARG:NH2	5:A:535:HOH:O	2.35	0.60
1:B:64:ASN:O	1:B:68:ASP:HB2	2.02	0.60
1:A:297:GLY:CA	1:A:308:ALA:HB2	2.32	0.59
1:B:50:VAL:HG22	1:B:90:ILE:CD1	2.33	0.59
1:B:110:ASN:HB3	1:B:127:LYS:HB2	1.86	0.58
1:B:135:ILE:CG2	1:B:138:LYS:HD2	2.31	0.58
1:B:288:HIS:C	1:B:289:ILE:HD12	2.23	0.58
1:A:41:ARG:HB2	1:A:257:MET:CE	2.34	0.58
1:B:201:LEU:C	1:B:201:LEU:HD23	2.24	0.58
1:B:301:PHE:O	1:B:302:ALA:HB3	2.04	0.58
1:B:320:VAL:HG13	1:B:334:VAL:HG22	1.86	0.57
1:B:67:ILE:HD11	1:B:106:ALA:CA	2.26	0.57
1:B:38:MET:CA	1:B:257:MET:HE1	2.33	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:122:ARG:HH11	1:A:122:ARG:HG3	1.69	0.57
1:A:35:ALA:O	1:A:39:ILE:HG13	2.05	0.57
1:B:368:HIS:C	1:B:370:ASN:H	2.07	0.57
1:B:22:MET:O	1:B:224:GLN:HA	2.05	0.56
1:B:123:ALA:CB	1:B:157:VAL:HG21	2.35	0.56
1:B:36:LYS:HE2	1:B:40:GLU:OE1	2.05	0.56
1:B:434:VAL:HA	1:B:437:ARG:NH1	2.20	0.56
1:B:71:MET:SD	1:B:77:TYR:CD2	3.00	0.55
1:A:140:ARG:HH11	1:A:140:ARG:HG3	1.72	0.55
1:A:167:SER:O	1:A:192:PRO:HD2	2.07	0.55
1:B:364:MET:HE2	1:B:365:LEU:CA	2.37	0.55
1:A:362:GLN:HG3	1:A:374:THR:HG21	1.88	0.55
1:A:21:GLN:HE22	1:A:237:HIS:N	2.04	0.55
1:B:73:THR:HG23	1:B:113:LEU:HD21	1.89	0.55
1:B:431:GLU:O	1:B:434:VAL:HG22	2.06	0.55
1:A:56:PHE:CE2	1:A:94:LEU:HB3	2.42	0.54
1:B:396:LEU:HD23	1:B:396:LEU:H	1.71	0.54
1:A:107:HIS:HB3	1:A:148:LYS:HB2	1.89	0.54
1:B:210:ALA:O	1:B:214:GLU:HG3	2.08	0.53
1:B:324:LEU:HD13	1:B:325:LYS:HG3	1.90	0.53
1:A:96:GLY:HA2	1:A:111:THR:O	2.08	0.53
1:A:97:LEU:CD2	1:A:113:LEU:HD12	2.37	0.53
1:A:97:LEU:HD21	1:A:113:LEU:HD12	1.91	0.53
1:B:339:VAL:HA	1:B:390:HIS:CD2	2.43	0.53
1:B:156:PRO:HD3	1:B:181:ASN:ND2	2.24	0.53
1:B:197:LEU:O	1:B:201:LEU:HB3	2.09	0.53
1:A:22:MET:HG2	1:A:53:PRO:HG3	1.91	0.53
1:A:173:ARG:HD2	5:A:527:HOH:O	2.09	0.52
1:B:41:ARG:CZ	1:B:257:MET:HE3	2.40	0.52
1:B:62:ASN:ND2	1:B:63:LYS:H	2.08	0.52
1:A:184:ARG:CB	1:A:184:ARG:HH11	2.22	0.52
1:A:91:TRP:CH2	1:A:122:ARG:HG2	2.44	0.52
1:B:156:PRO:HD3	1:B:181:ASN:HD22	1.73	0.52
1:A:196:THR:HG22	1:A:198:ASN:H	1.74	0.52
1:B:205:GLU:O	1:B:209:ARG:HG3	2.10	0.52
1:B:434:VAL:HA	1:B:437:ARG:HH12	1.75	0.52
1:A:302:ALA:HB2	1:A:388:VAL:HG23	1.92	0.52
1:A:142:MET:HB3	1:A:145:GLU:HG3	1.92	0.52
1:A:195:PHE:O	1:A:234:SER:HA	2.10	0.51
1:A:41:ARG:NE	1:A:257:MET:HE3	2.24	0.51
1:A:54:GLU:CB	1:A:193:SER:HA	2.27	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:235:TYR:O	1:B:237:HIS:HD2	1.94	0.51
1:B:63:LYS:C	1:B:65:GLU:H	2.13	0.51
1:A:142:MET:CB	1:A:145:GLU:HG3	2.40	0.50
1:A:20:CYS:O	1:A:222:ALA:HB1	2.11	0.50
1:A:173:ARG:O	1:A:175:PRO:HD3	2.11	0.50
1:A:201:LEU:HA	1:A:235:TYR:CE1	2.47	0.50
1:B:369:HIS:O	1:B:371:VAL:HG13	2.11	0.50
1:A:242:ASP:HB2	1:A:243:PRO:CD	2.42	0.50
1:B:207:LEU:O	1:B:211:ARG:HG2	2.12	0.50
1:B:169:CYS:HA	1:B:193:SER:OG	2.12	0.50
1:B:364:MET:HE1	1:B:436:TYR:CD1	2.47	0.50
1:B:340:VAL:HG11	1:B:345:ASP:HB2	1.92	0.50
1:A:41:ARG:HB2	1:A:257:MET:HE1	1.92	0.50
1:B:275:PRO:HG2	1:B:279:HIS:CE1	2.47	0.50
1:B:23:THR:HG22	1:B:255:VAL:HG22	1.93	0.49
1:A:263:ASP:OD2	1:B:254:ARG:NH2	2.35	0.49
1:B:364:MET:HE2	1:B:365:LEU:N	2.27	0.49
1:B:52:LEU:HD12	1:B:52:LEU:N	2.28	0.49
1:A:323:ASN:HD22	1:A:324:LEU:N	2.11	0.49
1:A:173:ARG:HH22	1:A:203:HIS:CE1	2.28	0.49
1:B:364:MET:CE	1:B:365:LEU:HD23	2.43	0.49
1:B:50:VAL:HG22	1:B:90:ILE:HD11	1.93	0.49
1:A:201:LEU:HD23	1:A:201:LEU:C	2.33	0.49
1:B:424:ARG:NH2	1:B:429:MET:HG2	2.28	0.49
1:B:218:TYR:CD2	1:B:262:ILE:HG22	2.48	0.49
1:A:158:ASP:OD1	1:A:163:ARG:HD3	2.13	0.49
1:A:205:GLU:O	1:A:209:ARG:HG3	2.13	0.49
1:A:342:ARG:HG3	1:A:342:ARG:HH11	1.78	0.48
1:B:209:ARG:O	1:B:213:ILE:HG13	2.13	0.48
1:B:41:ARG:HB2	1:B:257:MET:CE	2.42	0.48
1:A:197:LEU:O	1:A:201:LEU:HB3	2.13	0.48
1:B:366:GLU:O	1:B:370:ASN:HA	2.13	0.48
1:B:163:ARG:HB3	1:B:186:ALA:HA	1.95	0.48
1:B:56:PHE:CZ	1:B:94:LEU:HB3	2.49	0.48
1:B:364:MET:C	1:B:364:MET:HE2	2.33	0.48
1:B:36:LYS:HG3	1:B:84:LEU:HD11	1.96	0.48
1:B:63:LYS:C	1:B:65:GLU:N	2.66	0.48
1:B:303:ARG:HG2	1:B:303:ARG:HH11	1.78	0.48
1:A:138:LYS:O	1:A:139:VAL:HG13	2.14	0.47
1:A:163:ARG:HB3	1:A:186:ALA:HA	1.96	0.47
1:A:38:MET:HG2	1:A:257:MET:HE2	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:113:LEU:HD23	1:A:124:GLU:HG2	1.96	0.47
1:B:324:LEU:O	1:B:424:ARG:HG3	2.13	0.47
1:B:51:PHE:C	1:B:52:LEU:HD12	2.35	0.47
1:A:298:GLY:O	1:A:299:LEU:HB2	2.14	0.47
1:B:188:LEU:HD12	1:B:188:LEU:N	2.29	0.47
1:B:179:LEU:HD23	1:B:215:ASN:OD1	2.15	0.47
1:A:425:SER:OG	1:A:428:GLN:HG3	2.13	0.47
1:A:56:PHE:CZ	1:A:94:LEU:HB3	2.50	0.47
1:A:45:LYS:HE2	5:A:597:HOH:O	2.15	0.47
1:B:58:PHE:CZ	1:B:69:LEU:HB2	2.49	0.47
1:B:198:ASN:O	1:B:201:LEU:HD22	2.14	0.47
1:B:229:ASN:HB2	1:B:230:PRO:HD2	1.97	0.47
1:A:267:VAL:O	1:A:271:ARG:HG2	2.15	0.46
1:B:369:HIS:HD2	1:B:399:ARG:NH2	2.13	0.46
1:B:72:ALA:C	1:B:74:ASP:H	2.19	0.46
1:A:227:ALA:HA	1:A:233:GLN:NE2	2.31	0.46
1:B:328:THR:OG1	1:B:399:ARG:NH1	2.49	0.46
1:A:86:ARG:HA	1:A:117:SER:O	2.16	0.46
1:B:335:SER:OG	1:B:392:HIS:HB3	2.16	0.46
1:A:301:PHE:O	1:A:304:PHE:HB2	2.16	0.46
1:A:196:THR:HG22	1:A:198:ASN:N	2.31	0.45
1:B:59:ILE:HG23	1:B:232:ARG:HH11	1.81	0.45
1:B:111:THR:OG1	1:B:126:ASN:ND2	2.49	0.45
1:B:366:GLU:HG2	1:B:371:VAL:HG23	1.98	0.45
1:B:38:MET:HA	1:B:257:MET:CE	2.42	0.45
1:B:51:PHE:HB3	5:B:471:HOH:O	2.16	0.45
1:B:253:GLU:O	1:B:254:ARG:HB3	2.16	0.45
1:A:254:ARG:NH2	1:B:263:ASP:OD2	2.39	0.45
1:B:321:PHE:CE1	1:B:333:LEU:HD12	2.51	0.45
1:A:18:ALA:HB3	1:A:50:VAL:CG1	2.45	0.45
1:B:181:ASN:HA	1:B:181:ASN:HD22	1.64	0.45
1:A:338:ARG:NH1	1:A:340:VAL:HG11	2.32	0.45
1:B:342:ARG:HA	1:B:380:ASP:OD2	2.16	0.45
1:B:57:ASP:OD2	1:B:57:ASP:N	2.50	0.45
1:A:134:GLU:OE2	1:A:140:ARG:NH1	2.50	0.45
1:B:335:SER:HG	1:B:392:HIS:HB3	1.82	0.45
1:A:292:LYS:HE2	1:A:313:TYR:HE1	1.83	0.44
1:A:376:ILE:CD1	1:A:395:ILE:HG12	2.47	0.44
1:B:86:ARG:HA	1:B:117:SER:O	2.18	0.44
1:B:339:VAL:HA	1:B:390:HIS:NE2	2.33	0.44
1:A:140:ARG:HG3	1:A:140:ARG:NH1	2.31	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:204:TRP:CZ2	1:A:208:LEU:HD11	2.52	0.44
1:B:276:VAL:HG23	1:B:277:PHE:HD1	1.81	0.44
1:A:326:PRO:CD	1:A:424:ARG:HD2	2.47	0.44
1:B:324:LEU:O	1:B:324:LEU:HD22	2.16	0.44
1:A:398:ARG:HG3	1:A:398:ARG:HH11	1.82	0.44
1:A:113:LEU:CD2	1:A:124:GLU:HG2	2.48	0.44
1:A:375:THR:OG1	1:A:398:ARG:NH1	2.51	0.44
1:B:301:PHE:O	1:B:302:ALA:CB	2.64	0.44
1:B:368:HIS:C	1:B:370:ASN:N	2.71	0.44
1:B:424:ARG:HH21	1:B:429:MET:HG2	1.83	0.44
1:A:135:ILE:O	1:A:137:GLY:N	2.50	0.44
1:A:65:GLU:HG2	1:A:69:LEU:HD11	2.00	0.44
1:B:196:THR:HG22	1:B:199:THR:H	1.83	0.44
1:B:339:VAL:HG13	1:B:390:HIS:HD2	1.83	0.44
1:A:323:ASN:O	1:A:326:PRO:HD3	2.18	0.43
1:B:101:ASP:OD2	1:B:148:LYS:NZ	2.50	0.43
1:A:122:ARG:HH11	1:A:122:ARG:CG	2.31	0.43
1:B:56:PHE:CE1	1:B:94:LEU:HB3	2.53	0.43
1:A:297:GLY:HA3	1:A:308:ALA:HB2	2.00	0.43
1:B:59:ILE:HG23	1:B:232:ARG:NH1	2.33	0.43
1:B:126:ASN:O	1:B:127:LYS:C	2.56	0.43
1:B:238:SER:HB2	1:B:251:CYS:SG	2.58	0.43
1:A:127:LYS:HE2	1:A:129:HIS:O	2.19	0.43
1:A:38:MET:CA	1:A:257:MET:HE1	2.45	0.43
1:B:61:LEU:HD12	1:B:61:LEU:N	2.34	0.43
1:B:339:VAL:HG22	1:B:390:HIS:CD2	2.53	0.43
1:B:317:HIS:HB2	1:B:353:ASP:OD1	2.18	0.43
1:B:51:PHE:CD1	1:B:51:PHE:N	2.87	0.43
1:B:334:VAL:HB	1:B:393:ILE:HB	2.01	0.42
1:A:82:ARG:HD3	1:A:119:GLY:O	2.19	0.42
1:B:72:ALA:O	1:B:73:THR:CB	2.66	0.42
1:A:23:THR:HG23	1:A:225:THR:O	2.18	0.42
1:A:243:PRO:HB2	1:A:270:LEU:CD1	2.38	0.42
1:B:130:LEU:HD13	1:B:149:ALA:HB2	2.02	0.42
1:A:107:HIS:O	1:A:148:LYS:HE2	2.19	0.42
1:A:243:PRO:CG	1:A:267:VAL:HG22	2.50	0.42
1:B:41:ARG:HB2	1:B:257:MET:HE1	2.02	0.42
1:A:21:GLN:HG3	1:A:223:ALA:O	2.19	0.42
1:A:320:VAL:HA	1:A:333:LEU:O	2.19	0.42
1:B:281:ARG:NH1	1:B:283:ASP:OD1	2.53	0.42
1:B:53:PRO:O	1:B:56:PHE:HB3	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:392:HIS:CD2	1:B:392:HIS:N	2.88	0.42
1:B:401:GLY:O	1:B:402:ASP:C	2.59	0.42
1:B:326:PRO:HD2	1:B:424:ARG:HG3	2.01	0.42
1:A:41:ARG:CZ	1:A:257:MET:HE3	2.50	0.41
1:A:57:ASP:O	1:A:58:PHE:HB3	2.20	0.41
1:B:133:LEU:HD21	1:B:135:ILE:HG13	2.00	0.41
1:A:101:ASP:OD2	1:A:148:LYS:NZ	2.53	0.41
1:B:204:TRP:CE2	1:B:208:LEU:HD11	2.55	0.41
1:B:192:PRO:HA	1:B:222:ALA:O	2.21	0.41
1:A:343:LEU:HD11	1:A:378:VAL:HG11	2.02	0.41
1:B:115:ILE:CG2	1:B:116:ASP:N	2.83	0.41
1:A:299:LEU:HD12	1:A:311:ILE:HD13	2.02	0.41
1:A:41:ARG:HB2	1:A:257:MET:HE3	2.02	0.41
1:A:57:ASP:OD2	1:A:58:PHE:N	2.54	0.41
1:B:21:GLN:HB2	1:B:238:SER:OG	2.20	0.41
1:A:253:GLU:O	1:A:254:ARG:HB3	2.21	0.41
1:B:307:PRO:HD2	1:B:310:HIS:ND1	2.35	0.41
1:B:396:LEU:HA	1:B:397:PRO:HD2	1.89	0.40
1:A:227:ALA:HB2	1:A:233:GLN:HE22	1.86	0.40
1:B:35:ALA:O	1:B:39:ILE:HG13	2.21	0.40
1:A:179:LEU:HD23	1:A:215:ASN:OD1	2.21	0.40
1:A:218:TYR:CD2	1:A:242:ASP:HA	2.56	0.40
1:A:424:ARG:NH2	1:A:432:GLU:OE1	2.48	0.40
1:A:21:GLN:NE2	1:A:237:HIS:N	2.69	0.40
1:A:323:ASN:HD22	1:A:324:LEU:H	1.67	0.40
1:B:198:ASN:HA	1:B:201:LEU:HD13	2.04	0.40
1:B:229:ASN:HB2	1:B:230:PRO:CD	2.51	0.40
1:B:95:GLY:HA2	1:B:192:PRO:HG3	2.04	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	408/440 (93%)	376 (92%)	23 (6%)	9 (2%)	6	22
1	B	406/440 (92%)	359 (88%)	36 (9%)	11 (3%)	5	17
All	All	814/880 (92%)	735 (90%)	59 (7%)	20 (2%)	5	19

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	169	CYS
1	A	402	ASP
1	B	136	PRO
1	A	54	GLU
1	A	95	GLY
1	A	424	ARG
1	B	170	TYR
1	B	302	ALA
1	A	96	GLY
1	B	67	ILE
1	B	95	GLY
1	B	370	ASN
1	B	401	GLY
1	A	136	PRO
1	A	170	TYR
1	A	299	LEU
1	B	204	TRP
1	B	96	GLY
1	B	369	HIS
1	B	341	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	357/383 (93%)	338 (95%)	19 (5%)	22	54
1	B	353/383 (92%)	332 (94%)	21 (6%)	19	49
All	All	710/766 (93%)	670 (94%)	40 (6%)	21	51

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	50	VAL
1	A	122	ARG
1	A	139	VAL
1	A	141	LEU
1	A	151	THR
1	A	157	VAL
1	A	172	VAL
1	A	177	LEU
1	A	184	ARG
1	A	193	SER
1	A	195	PHE
1	A	201	LEU
1	A	230	PRO
1	A	233	GLN
1	A	287	LEU
1	A	296	THR
1	A	323	ASN
1	A	329	ASP
1	A	402	ASP
1	B	23	THR
1	B	50	VAL
1	B	68	ASP
1	B	76	GLU
1	B	80	LYS
1	B	116	ASP
1	B	133	LEU
1	B	154	ILE
1	B	163	ARG
1	B	177	LEU
1	B	181	ASN
1	B	195	PHE
1	B	201	LEU
1	B	279	HIS
1	B	323	ASN
1	B	338	ARG
1	B	364	MET
1	B	370	ASN
1	B	392	HIS
1	B	396	LEU
1	B	425	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	GLN
1	A	37	ASN
1	A	126	ASN
1	A	181	ASN
1	A	233	GLN
1	A	323	ASN
1	A	370	ASN
1	A	438	ASN
1	B	62	ASN
1	B	64	ASN
1	B	66	GLN
1	B	126	ASN
1	B	181	ASN
1	B	237	HIS
1	B	323	ASN
1	B	390	HIS
1	B	392	HIS
1	B	438	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 4 are monoatomic - leaving 6 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the

expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	EMC	B	452	1	1,2,2	0.61	0	-		
2	EMC	B	455	1	1,2,2	0.51	0	-		
4	MPD	A	502	-	7,7,7	0.47	0	9,10,10	0.53	0
2	EMC	A	453	1	1,2,2	0.55	0	-		
2	EMC	A	451	1	1,2,2	0.72	0	-		
2	EMC	A	454	1	1,2,2	0.75	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
4	MPD	A	502	-	-	0/5/5/5	-

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, 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	412/440 (93%)	-0.38	6 (1%) 73 68	7, 18, 41, 63	0
1	B	410/440 (93%)	0.11	32 (7%) 13 7	9, 33, 69, 88	0
All	All	822/880 (93%)	-0.14	38 (4%) 32 22	7, 23, 64, 88	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	296	THR	5.4
1	B	329	ASP	4.0
1	B	368	HIS	3.9
1	B	10	MET	3.9
1	B	399	ARG	3.7
1	B	369	HIS	3.7
1	B	400	ALA	3.4
1	B	64	ASN	3.4
1	B	297	GLY	3.3
1	A	136	PRO	3.3
1	B	66	GLN	3.2
1	B	294	SER	3.1
1	A	296	THR	3.0
1	A	370	ASN	3.0
1	B	435	VAL	3.0
1	B	136	PRO	2.7
1	A	138	LYS	2.6
1	B	295	GLU	2.6
1	B	12	THR	2.5
1	B	67	ILE	2.5
1	B	426	ASN	2.5
1	B	397	PRO	2.4
1	B	424	ARG	2.4
1	A	10	MET	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	438	ASN	2.3
1	B	330	GLY	2.3
1	B	62	ASN	2.3
1	B	140	ARG	2.2
1	B	436	TYR	2.2
1	B	106	ALA	2.2
1	B	73	THR	2.2
1	B	439	LEU	2.1
1	B	428	GLN	2.1
1	B	323	ASN	2.1
1	A	233	GLN	2.1
1	B	322	VAL	2.0
1	B	437	ARG	2.0
1	B	432	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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
4	MPD	A	502	8/8	0.87	0.22	40,43,45,45	0
3	NA	B	458	1/1	0.94	0.26	20,20,20,20	0
3	NA	A	459	1/1	0.95	0.32	11,11,11,11	0
3	NA	A	456	1/1	0.97	0.10	16,16,16,16	0
2	EMC	A	453	3/3	0.99	0.10	25,25,25,25	3
2	EMC	B	455	3/3	0.99	0.09	24,24,24,24	3
3	NA	A	457	1/1	0.99	0.13	19,19,19,19	0
2	EMC	A	454	3/3	0.99	0.11	25,25,25,25	3
2	EMC	A	451	3/3	1.00	0.09	24,24,24,24	3

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	EMC	B	452	3/3	1.00	0.10	23,23,23,23	3

6.5 Other polymers [i](#)

There are no such residues in this entry.