Full wwPDB X-ray Structure Validation Report (i)

Jun 12, 2024 – 11:12 PM EDT

PDB ID	:	1JMX
Title	:	crystal structure of a quinohemoprotein amine dehydrogenase from pseudomonas putida
Authors	:	Satoh, A.; Miyahara, I.; Hirotsu, K.
Deposited on	:	2001-07-20
Resolution	:	1.90 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 (i) 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 (1)) were used in the production of this report:

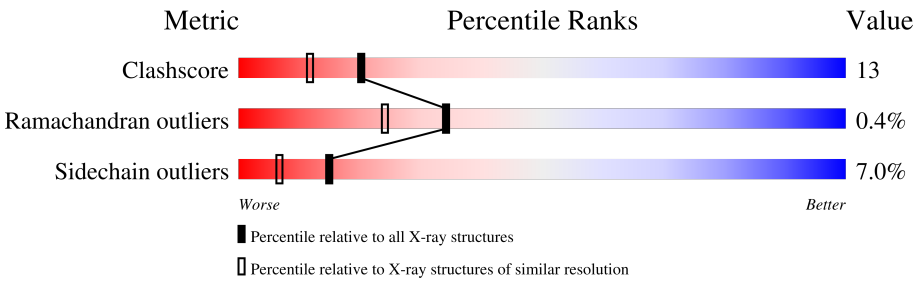
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

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 1.90 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))
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)

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 <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	494	76% 20% •
2	B	349	75% 20% • •
3	G	79	59% 27% 10% • •

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 7611 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 Amine Dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	493	3790	2372	688	713	17	10	0	0

- Molecule 2 is a protein called Amine Dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	339	2689	1724	453	499	13	11	0	0

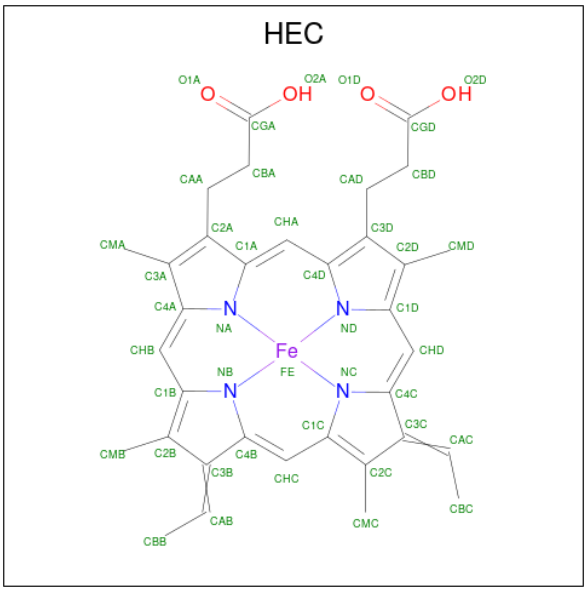
- Molecule 3 is a protein called Amine Dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	G	77	588	365	94	122	7	0	0	0

- Molecule 4 is NICKEL (II) ION (three-letter code: NI) (formula: Ni).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Ni		
4	A	1	1	1	0	0

- Molecule 5 is HEME C (three-letter code: HEC) (formula: $C_{34}H_{34}FeN_4O_4$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	Fe	N	O		
5	A	1	43	34	1	4	4	0	0
			Total	C	Fe	N	O		
5	A	1	43	34	1	4	4	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
6	A	247	247	247	0	0
			Total	O		
6	B	159	159	159	0	0
			Total	O		
6	G	51	51	51	0	0

3 Residue-property plots (i)

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. 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. 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.

Note EDS was not executed.

• Molecule 1: Amine Dehydrogenase

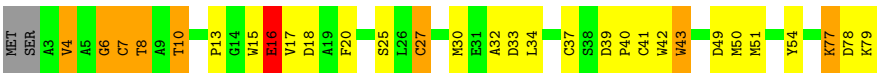
Chain A:

• Molecule 2: Amine Dehydrogenase

Chain B:

• Molecule 3: Amine Dehydrogenase

Chain G:



4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	167.21Å 92.37Å 79.30Å 90.00° 112.00° 90.00°	Depositor
Resolution (Å)	10.00 – 1.90	Depositor
% Data completeness (in resolution range)	(Not available) (10.00-1.90)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.209 , 0.267	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7611	wwPDB-VP
Average B, all atoms (Å^2)	20.0	wwPDB-VP

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: HEC, TRQ, NI

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	0.41	0/3872	0.68	2/5240 (0.0%)
2	B	0.42	0/2761	0.74	1/3752 (0.0%)
3	G	1.39	4/588 (0.7%)	1.49	14/802 (1.7%)
All	All	0.56	4/7221 (0.1%)	0.80	17/9794 (0.2%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	G	6	GLY	C-N	17.83	1.75	1.34
3	G	16	GLU	C-N	16.70	1.72	1.34
3	G	7	CYS	CB-SG	-14.04	1.58	1.82
3	G	15	TRP	C-N	11.35	1.60	1.34

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	G	7	CYS	O-C-N	16.80	149.58	122.70
3	G	7	CYS	CA-C-N	-12.95	88.71	117.20
3	G	41	CYS	CA-CB-SG	-11.89	92.59	114.00
3	G	7	CYS	CA-CB-SG	-10.39	95.30	114.00
3	G	16	GLU	N-CA-CB	-8.80	94.75	110.60
3	G	6	GLY	O-C-N	-8.56	109.00	122.70
3	G	16	GLU	CA-C-N	-8.19	99.18	117.20
3	G	6	GLY	C-N-CA	7.76	141.09	121.70
3	G	15	TRP	O-C-N	-7.17	111.23	122.70
3	G	16	GLU	CA-CB-CG	6.87	128.51	113.40
3	G	7	CYS	C-N-CA	-6.81	104.67	121.70
2	B	19	TYR	N-CA-C	5.98	127.14	111.00
3	G	15	TRP	C-N-CA	5.49	135.41	121.70
3	G	49	ASP	N-CA-C	5.33	125.39	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	407	LEU	CA-CB-CG	5.24	127.34	115.30
1	A	391	GLY	N-CA-C	-5.17	100.19	113.10
3	G	6	GLY	CA-C-N	5.01	128.23	117.20

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	3790	0	3728	92	0
2	B	2689	0	2645	53	0
3	G	588	0	510	39	0
4	A	1	0	0	0	0
5	A	86	0	60	7	0
6	A	247	0	0	5	0
6	B	159	0	0	0	0
6	G	51	0	0	1	0
All	All	7611	0	6943	175	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 13.

All (175) 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)
3:G:7:CYS:SG	3:G:16:GLU:HG2	1.50	1.52
3:G:7:CYS:CB	3:G:16:GLU:HG2	1.40	1.46
3:G:16:GLU:C	3:G:17:VAL:N	1.72	1.42
3:G:6:GLY:C	3:G:7:CYS:N	1.75	1.37
1:A:484:THR:HG21	3:G:16:GLU:O	1.33	1.21
3:G:7:CYS:HB3	3:G:16:GLU:HG2	1.20	1.09
3:G:7:CYS:SG	3:G:16:GLU:HG3	1.94	1.07
3:G:7:CYS:CB	3:G:16:GLU:CG	2.37	0.95
1:A:469:THR:HG22	1:A:476:THR:HG22	1.48	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:50:LEU:HD11	5:A:1002:HEC:HBB2	1.54	0.90
1:A:271:GLU:HG2	1:A:389:VAL:HG13	1.54	0.87
2:B:314:THR:HG22	2:B:315:PHE:H	1.37	0.87
5:A:1001:HEC:HMD1	5:A:1001:HEC:HBD2	1.58	0.84
2:B:314:THR:HG22	2:B:315:PHE:N	1.93	0.82
3:G:7:CYS:HB3	3:G:16:GLU:CG	2.02	0.81
1:A:347:ARG:NH1	1:A:361:VAL:HG21	1.98	0.79
3:G:16:GLU:CA	3:G:17:VAL:N	2.47	0.76
1:A:251:GLN:HG2	1:A:264:MET:HG3	1.67	0.75
1:A:86:LEU:O	3:G:4:VAL:HG13	1.86	0.74
1:A:220:ASP:OD2	1:A:222:THR:HG22	1.91	0.71
2:B:18:ASN:HD22	2:B:20:PRO:HD2	1.56	0.70
1:A:484:THR:CG2	3:G:16:GLU:O	2.27	0.70
3:G:16:GLU:N	3:G:17:VAL:N	2.41	0.69
2:B:18:ASN:ND2	2:B:20:PRO:HD2	2.08	0.67
2:B:202:PRO:HD2	2:B:343:THR:HG21	1.75	0.66
1:A:288:ALA:HA	1:A:357:VAL:HG11	1.76	0.66
1:A:241:ASN:HD22	1:A:250:ARG:HD2	1.62	0.65
3:G:16:GLU:C	3:G:17:VAL:CA	2.63	0.63
2:B:270:ILE:HB	2:B:282:LEU:HD21	1.80	0.63
1:A:416:TRP:CZ3	1:A:468:ALA:HB2	2.34	0.62
3:G:7:CYS:CA	3:G:16:GLU:HG2	2.27	0.62
1:A:271:GLU:CG	1:A:389:VAL:HG13	2.28	0.61
2:B:208:HIS:HB3	2:B:241:LYS:NZ	2.14	0.61
1:A:350:ALA:HB2	1:A:355:LYS:HD3	1.81	0.61
1:A:241:ASN:HD21	1:A:250:ARG:HH11	1.50	0.60
1:A:26:ARG:NH2	6:A:2003:HOH:O	2.34	0.59
3:G:13:PRO:O	3:G:18:ASP:HA	2.02	0.59
2:B:208:HIS:HB3	2:B:241:LYS:HZ3	1.66	0.58
1:A:98:GLU:HG2	1:A:138:GLN:NE2	2.19	0.57
1:A:241:ASN:ND2	1:A:250:ARG:HH11	2.02	0.57
2:B:242:THR:OG1	2:B:244:LYS:HD3	2.04	0.57
1:A:190:MET:HG3	1:A:273:GLY:O	2.04	0.57
1:A:328:THR:HB	1:A:329:PRO:HD2	1.86	0.57
1:A:218:TYR:CD2	1:A:222:THR:HG23	2.40	0.57
1:A:241:ASN:ND2	1:A:250:ARG:NH1	2.53	0.57
1:A:193:LYS:NZ	1:A:222:THR:HG21	2.21	0.56
1:A:63:TYR:O	1:A:67:LYS:HE2	2.06	0.56
3:G:8:THR:HG22	3:G:10:THR:H	1.70	0.56
1:A:209:THR:HG22	1:A:231:ILE:HG12	1.88	0.56
2:B:204:GLN:H	2:B:345:GLN:NE2	2.03	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:259:GLU:HG2	1:A:261:LYS:HE2	1.86	0.56
1:A:261:LYS:HE3	6:A:2108:HOH:O	2.06	0.55
1:A:2:GLU:HB2	1:A:7:LEU:HG	1.88	0.55
2:B:262:ARG:HD2	2:B:267:PRO:O	2.07	0.55
2:B:19:TYR:O	2:B:20:PRO:C	2.44	0.55
2:B:314:THR:CG2	2:B:315:PHE:N	2.65	0.55
2:B:339:MET:HB3	2:B:342:THR:HG22	1.88	0.54
1:A:279:VAL:HG23	1:A:287:LEU:HD21	1.90	0.54
1:A:214:VAL:O	1:A:215:LYS:HD2	2.06	0.54
1:A:145:LEU:HB3	1:A:146:PRO:HD3	1.88	0.54
1:A:343:LYS:O	1:A:347:ARG:NH2	2.40	0.54
1:A:4:GLY:N	1:A:5:PRO:HD2	2.22	0.54
1:A:251:GLN:HG2	1:A:264:MET:CG	2.38	0.53
2:B:248:GLN:O	2:B:248:GLN:HG3	2.08	0.53
1:A:285:ARG:HB3	1:A:285:ARG:NH1	2.24	0.53
1:A:394:GLU:HG2	1:A:411:TYR:CG	2.44	0.53
2:B:298:TYR:OH	3:G:33:ASP:HA	2.08	0.53
1:A:196:VAL:HG11	1:A:274:LEU:HD13	1.91	0.52
1:A:197:ARG:HD3	6:A:2219:HOH:O	2.09	0.52
1:A:317:ALA:O	1:A:347:ARG:NH1	2.42	0.52
2:B:18:ASN:HD21	2:B:336:GLY:HA3	1.75	0.52
1:A:386:VAL:HG21	1:A:486:GLN:HG2	1.91	0.52
1:A:396:GLU:HG3	1:A:408:ARG:HH11	1.74	0.52
2:B:250:PHE:O	2:B:251:ALA:HB2	2.09	0.52
1:A:63:TYR:CE2	1:A:67:LYS:HE3	2.45	0.51
1:A:127:LEU:HD13	2:B:122:VAL:HG11	1.92	0.51
2:B:28:VAL:O	2:B:308:LYS:NZ	2.44	0.51
1:A:456:MET:O	1:A:457:MET:HG2	2.11	0.51
3:G:50:MET:HG2	3:G:54:TYR:O	2.11	0.51
2:B:314:THR:HG23	3:G:32:ALA:HB1	1.93	0.51
2:B:110:VAL:O	2:B:112:PRO:HD3	2.11	0.51
1:A:418:VAL:HG12	1:A:431:VAL:HG22	1.92	0.50
2:B:112:PRO:HG2	2:B:123:LYS:HB2	1.94	0.50
3:G:39:ASP:O	3:G:51:MET:HB2	2.12	0.50
1:A:99:THR:HG21	1:A:152:VAL:HG21	1.94	0.50
1:A:279:VAL:HG11	1:A:285:ARG:HG3	1.93	0.50
1:A:250:ARG:HD3	1:A:267:ALA:HB2	1.94	0.49
2:B:30:SER:O	2:B:32:THR:HG23	2.13	0.49
3:G:8:THR:HG21	3:G:43:TRQ:HD1	1.95	0.49
1:A:241:ASN:HD21	1:A:250:ARG:NH1	2.11	0.49
1:A:127:LEU:HB3	2:B:122:VAL:HG21	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:176:LYS:HE3	1:A:178:ASP:CG	2.32	0.48
2:B:217:ALA:HB1	2:B:229:ALA:HB1	1.95	0.48
2:B:234:GLY:HA2	2:B:249:GLU:HA	1.95	0.48
3:G:16:GLU:H	3:G:17:VAL:N	2.11	0.48
3:G:27:CYS:HB2	6:G:88:HOH:O	2.12	0.48
2:B:24:HIS:HD2	2:B:36:SER:OG	1.97	0.47
1:A:177:ALA:HB1	1:A:210:PHE:CZ	2.50	0.47
2:B:263:SER:HB3	2:B:266:ASP:O	2.15	0.47
2:B:316:ASN:HB3	2:B:337:GLY:O	2.15	0.47
1:A:99:THR:HG22	5:A:1001:HEC:HBB1	1.96	0.47
1:A:153:VAL:HB	1:A:154:PRO:HD3	1.97	0.46
2:B:339:MET:HB3	2:B:342:THR:CG2	2.44	0.46
1:A:485:VAL:HG11	3:G:30:MET:HB2	1.97	0.46
1:A:103:CYS:HB2	6:A:2013:HOH:O	2.15	0.46
1:A:177:ALA:HB1	1:A:210:PHE:HZ	1.80	0.46
3:G:79:LYS:HZ2	3:G:79:LYS:N	2.13	0.46
3:G:37:CYS:HB2	3:G:43:TRQ:HZ3	1.68	0.46
1:A:102:ARG:HH11	3:G:10:THR:HG21	1.81	0.46
1:A:396:GLU:HG3	1:A:408:ARG:NH1	2.31	0.46
2:B:20:PRO:O	2:B:21:ASN:HB2	2.15	0.45
1:A:417:LYS:NZ	1:A:467:ILE:HG21	2.30	0.45
2:B:18:ASN:CB	2:B:24:HIS:HE1	2.28	0.45
1:A:155:ASP:O	1:A:159:ARG:HG3	2.16	0.45
5:A:1002:HEC:HAD1	2:B:119:HIS:HA	1.98	0.45
2:B:109:THR:HG21	2:B:150:VAL:HG11	1.99	0.45
3:G:39:ASP:HB2	3:G:40:PRO:HD3	1.98	0.45
1:A:313:PRO:HA	1:A:350:ALA:O	2.17	0.45
2:B:296:THR:O	2:B:314:THR:HG21	2.17	0.45
5:A:1001:HEC:HBA1	5:A:1001:HEC:HMA2	1.98	0.44
3:G:7:CYS:HB3	3:G:16:GLU:CB	2.47	0.44
2:B:18:ASN:HD21	2:B:337:GLY:H	1.65	0.44
1:A:115:PRO:HD3	1:A:162:LEU:HG	2.00	0.44
3:G:7:CYS:HA	3:G:16:GLU:CD	2.38	0.44
1:A:350:ALA:HA	1:A:354:LEU:O	2.17	0.43
1:A:284:ALA:HB1	1:A:354:LEU:HD11	1.99	0.43
1:A:17:ILE:O	1:A:17:ILE:HG13	2.18	0.43
2:B:74:THR:O	2:B:75:CYS:HB2	2.18	0.43
3:G:79:LYS:HZ2	3:G:79:LYS:H	1.66	0.43
1:A:39:MET:O	1:A:43:ARG:HG3	2.19	0.43
2:B:190:LYS:HD2	2:B:190:LYS:HA	1.81	0.43
1:A:172:LYS:HA	1:A:172:LYS:HE2	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:215:LYS:HE3	1:A:225:ASN:HB3	2.01	0.43
1:A:365:VAL:HG11	1:A:470:LEU:CD2	2.49	0.43
2:B:28:VAL:HG13	2:B:310:TYR:OH	2.18	0.43
2:B:313:GLY:O	2:B:314:THR:HB	2.18	0.43
1:A:241:ASN:ND2	1:A:250:ARG:HD2	2.32	0.42
1:A:43:ARG:NH2	5:A:1002:HEC:HBA2	2.34	0.42
1:A:193:LYS:CE	1:A:222:THR:HG21	2.49	0.42
3:G:20:PHE:CD1	3:G:25:SER:HB2	2.54	0.42
3:G:77:LYS:HD2	3:G:78:ASP:H	1.85	0.42
1:A:218:TYR:HD2	1:A:222:THR:HG23	1.82	0.42
2:B:260:GLY:HA2	2:B:271:TYR:O	2.19	0.42
2:B:326:LEU:HD23	2:B:326:LEU:HA	1.80	0.42
1:A:114:ARG:HB3	1:A:118:GLU:HB2	2.01	0.42
3:G:20:PHE:HD1	3:G:25:SER:HB2	1.84	0.42
1:A:17:ILE:HA	1:A:18:PRO:HD3	1.80	0.42
1:A:396:GLU:OE2	1:A:408:ARG:NH1	2.53	0.42
2:B:48:ALA:HB2	2:B:59:VAL:HG22	2.01	0.42
2:B:314:THR:CG2	3:G:32:ALA:HB1	2.49	0.42
1:A:312:LYS:HA	1:A:313:PRO:HD3	1.88	0.42
1:A:96:LEU:HG	1:A:159:ARG:NH2	2.34	0.41
1:A:50:LEU:CD1	5:A:1002:HEC:HBB2	2.37	0.41
1:A:469:THR:HG23	6:A:2091:HOH:O	2.19	0.41
3:G:8:THR:CG2	3:G:10:THR:HG23	2.51	0.41
1:A:148:ALA:O	1:A:153:VAL:HG23	2.19	0.41
1:A:222:THR:HA	1:A:223:PRO:HD3	1.92	0.41
1:A:416:TRP:CH2	1:A:468:ALA:HB2	2.55	0.41
1:A:17:ILE:HD12	1:A:18:PRO:O	2.20	0.41
1:A:269:HIS:HB3	1:A:271:GLU:OE2	2.21	0.41
2:B:250:PHE:HB3	2:B:282:LEU:HD13	2.02	0.41
1:A:60:LEU:HD23	1:A:60:LEU:HA	1.95	0.41
2:B:19:TYR:HB2	2:B:338:ASP:O	2.20	0.41
2:B:273:VAL:CG1	2:B:300:VAL:HG13	2.51	0.41
2:B:318:LEU:HB2	2:B:332:ILE:HB	2.03	0.41
2:B:124:PRO:HA	2:B:125:PRO:HD3	1.95	0.41
1:A:136:GLN:HG2	3:G:42:TRP:CE2	2.56	0.40
1:A:271:GLU:HG2	1:A:389:VAL:CG1	2.37	0.40
2:B:19:TYR:O	2:B:20:PRO:O	2.39	0.40
3:G:16:GLU:C	3:G:17:VAL:HG23	2.42	0.40
1:A:303:THR:HB	1:A:333:ARG:HG2	2.02	0.40
2:B:235:TYR:O	2:B:247:THR:HA	2.22	0.40
2:B:288:ILE:HG13	2:B:289:LYS:N	2.35	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:267:ALA:HB3	1:A:268:GLU:OE2	2.22	0.40
1:A:348:GLU:OE2	1:A:355:LYS:HB3	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	491/494 (99%)	471 (96%)	19 (4%)	1 (0%)	47 38
2	B	335/349 (96%)	318 (95%)	14 (4%)	3 (1%)	17 7
3	G	74/79 (94%)	68 (92%)	6 (8%)	0	100 100
All	All	900/922 (98%)	857 (95%)	39 (4%)	4 (0%)	34 24

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	21	ASN
1	A	21	ASN
2	B	274	LEU
2	B	314	THR

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	387/387 (100%)	358 (92%)	29 (8%)	13 5
2	B	291/300 (97%)	275 (94%)	16 (6%)	21 12
3	G	61/63 (97%)	54 (88%)	7 (12%)	5 2
All	All	739/750 (98%)	687 (93%)	52 (7%)	15 7

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	SER
1	A	9	GLN
1	A	13	MET
1	A	17	ILE
1	A	21	ASN
1	A	50	LEU
1	A	67	LYS
1	A	127	LEU
1	A	162	LEU
1	A	163	GLU
1	A	172	LYS
1	A	191	LEU
1	A	197	ARG
1	A	206	GLN
1	A	215	LYS
1	A	232	LEU
1	A	264	MET
1	A	271	GLU
1	A	274	LEU
1	A	299	GLU
1	A	303	THR
1	A	323	GLU
1	A	348	GLU
1	A	407	LEU
1	A	451	ASN
1	A	470	LEU
1	A	484	THR
1	A	486	GLN
1	A	490	ASN
2	B	19	TYR
2	B	60	LEU
2	B	92	ARG
2	B	109	THR
2	B	143	THR

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Mol	Chain	Res	Type
2	B	199	TYR
2	B	233	TYR
2	B	236	LEU
2	B	248	GLN
2	B	277	LEU
2	B	286	LYS
2	B	288	ILE
2	B	309	LEU
2	B	326	LEU
2	B	343	THR
2	B	345	GLN
3	G	4	VAL
3	G	8	THR
3	G	10	THR
3	G	16	GLU
3	G	27	CYS
3	G	34	LEU
3	G	77	LYS

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	138	GLN
1	A	241	ASN
1	A	382	ASN
1	A	451	ASN
1	A	489	ASN
2	B	18	ASN
2	B	24	HIS
2	B	80	HIS
2	B	111	ASN
2	B	345	GLN
3	G	55	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	$\#\|Z\| > 2$
3	TRQ	G	43	3	13,17,18	3.73	3 (23%)	14,24,26	3.17	7 (50%)

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	TRQ	G	43	3	-	0/4/19/21	0/2/2/2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	G	43	TRQ	CH2-CZ2	-10.15	1.42	1.54
3	G	43	TRQ	CE2-CZ2	-6.30	1.41	1.50
3	G	43	TRQ	CZ3-CE3	5.62	1.44	1.34

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	G	43	TRQ	CZ2-CE2-NE1	6.32	130.03	119.94
3	G	43	TRQ	O7-CZ2-CH2	5.63	125.62	119.00
3	G	43	TRQ	O7-CZ2-CE2	-5.61	115.90	121.84
3	G	43	TRQ	O6-CH2-CZ2	4.00	121.23	118.51
3	G	43	TRQ	CD2-CE3-CZ3	-2.54	118.02	121.09
3	G	43	TRQ	O6-CH2-CZ3	-2.35	117.25	121.51
3	G	43	TRQ	CZ3-CH2-CZ2	2.05	121.52	118.72

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	G	43	TRQ	2	0

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 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	$\#\|Z\| > 2$
5	HEC	A	1002	1	32,50,50	1.66	3 (9%)	24,82,82	1.40	3 (12%)
5	HEC	A	1001	1	32,50,50	1.72	4 (12%)	24,82,82	1.85	6 (25%)

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
5	HEC	A	1002	1	-	2/10/54/54	-
5	HEC	A	1001	1	-	3/10/54/54	-

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1001	HEC	C3C-C2C	-6.50	1.34	1.40
5	A	1002	HEC	C3C-C2C	-6.37	1.34	1.40
5	A	1001	HEC	C2B-C3B	-3.83	1.36	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1002	HEC	C2B-C3B	-3.46	1.37	1.40
5	A	1001	HEC	O2A-CGA	-2.68	1.21	1.30
5	A	1001	HEC	C4B-C3B	2.32	1.47	1.43
5	A	1002	HEC	C3C-C4C	2.24	1.47	1.43

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	A	1001	HEC	CBA-CAA-C2A	-4.75	104.59	112.60
5	A	1001	HEC	C1D-C2D-C3D	-4.01	104.21	107.00
5	A	1002	HEC	C1D-C2D-C3D	-3.29	104.71	107.00
5	A	1001	HEC	CBD-CAD-C3D	3.04	117.81	112.62
5	A	1002	HEC	CMB-C2B-C3B	2.33	128.56	125.82
5	A	1001	HEC	CMD-C2D-C3D	2.23	129.15	124.94
5	A	1001	HEC	CMB-C2B-C3B	2.11	128.30	125.82
5	A	1002	HEC	CBA-CAA-C2A	-2.06	109.13	112.60
5	A	1001	HEC	CMC-C2C-C1C	-2.04	125.33	128.46

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	1001	HEC	C2D-C3D-CAD-CBD
5	A	1001	HEC	C4D-C3D-CAD-CBD
5	A	1001	HEC	C3D-CAD-CBD-CGD
5	A	1002	HEC	CAD-CBD-CGD-O2D
5	A	1002	HEC	CAD-CBD-CGD-O1D

There are no ring outliers.

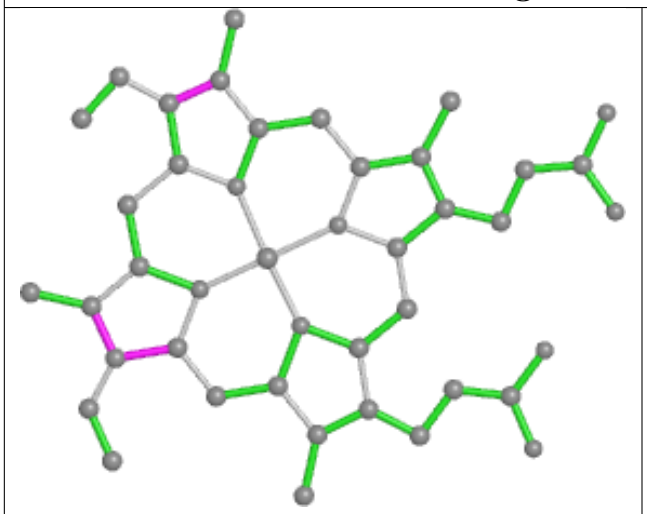
2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1002	HEC	4	0
5	A	1001	HEC	3	0

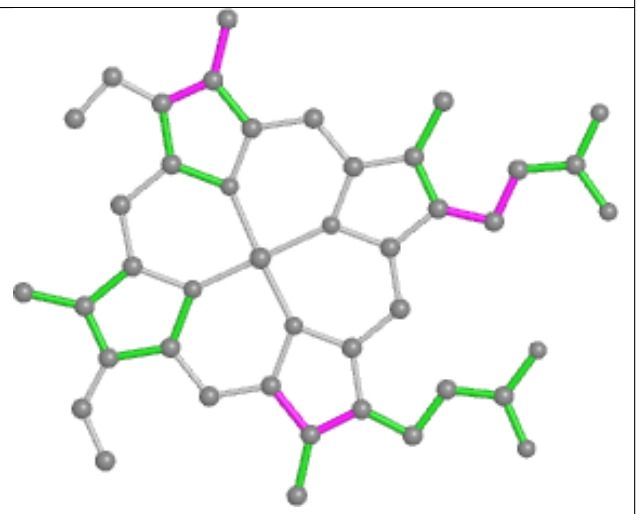
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be

highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

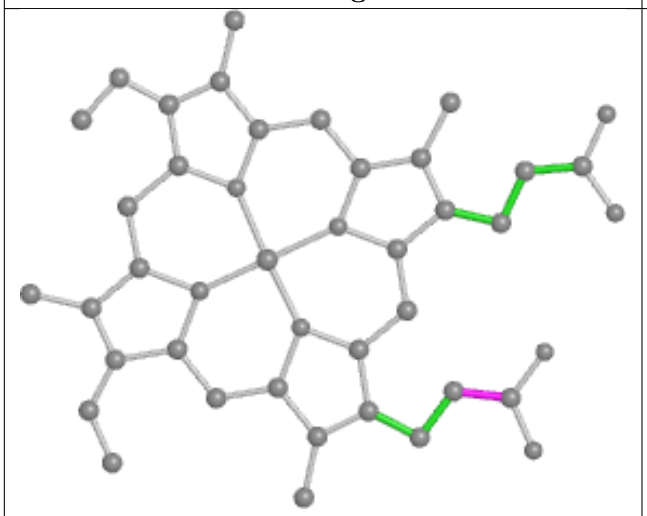
Ligand HEC A 1002



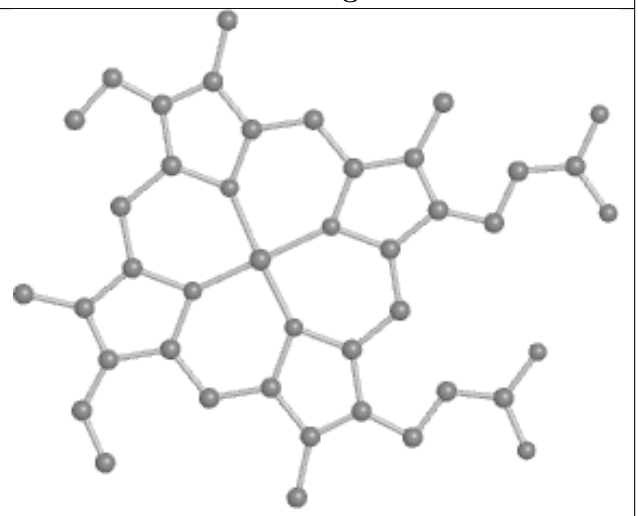
Bond lengths



Bond angles

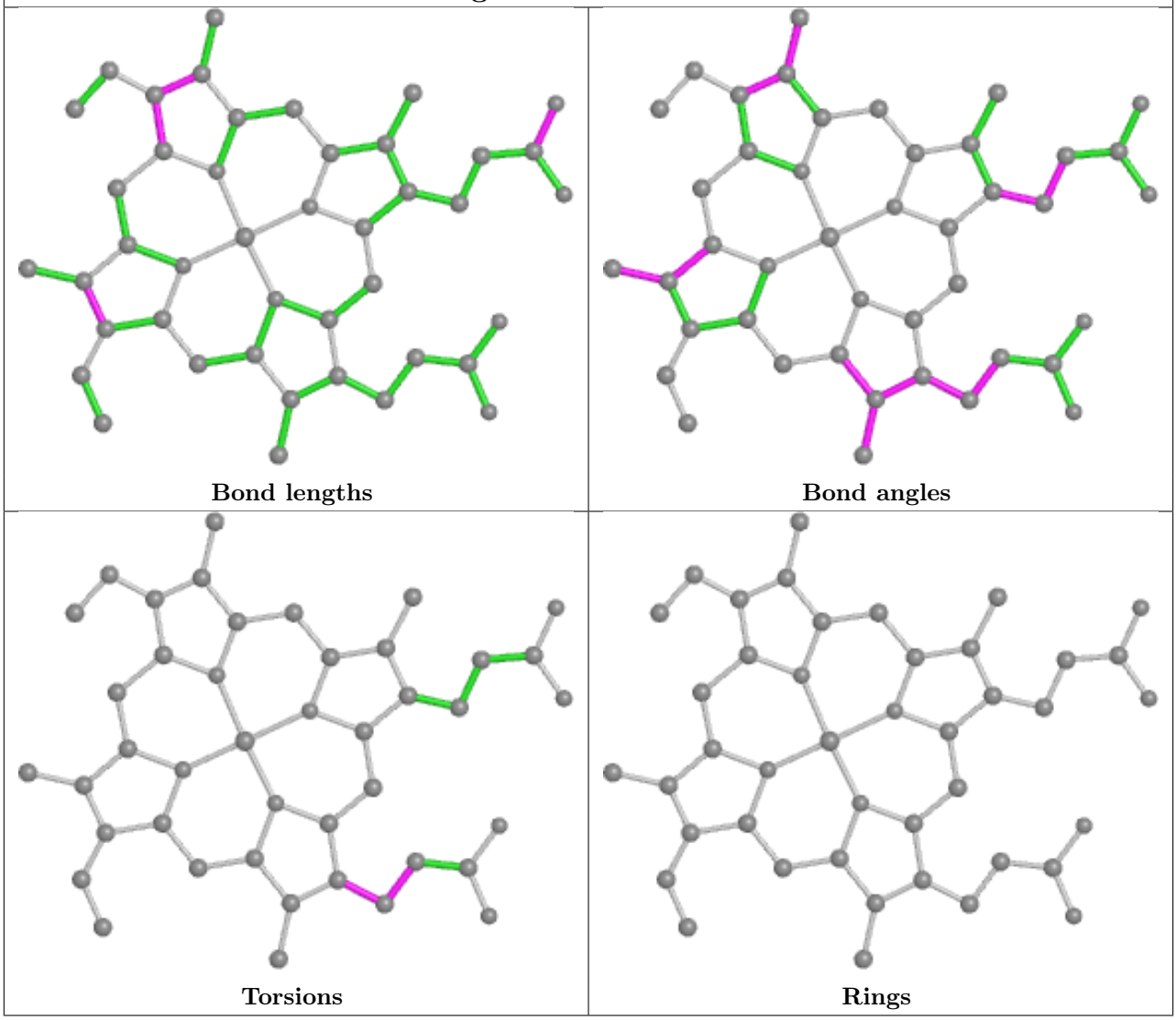


Torsions



Rings

Ligand HEC A 1001



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
3	G	2

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	G	6:GLY	C	7:CYS	N	1.75
1	G	16:GLU	C	17:VAL	N	1.72

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.