



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:09 PM UTC

PDB ID : 1ZZ9 / pdb_00001zz9
Title : Crystal Structure of FeII HppE
Authors : Higgins, L.J.; Yan, F.; Liu, P.; Liu, H.W.; Drennan, C.L.
Deposited on : 2005-06-13
Resolution : 2.40 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
Mogul : 2022.3.0, CSD as543be (2022)
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

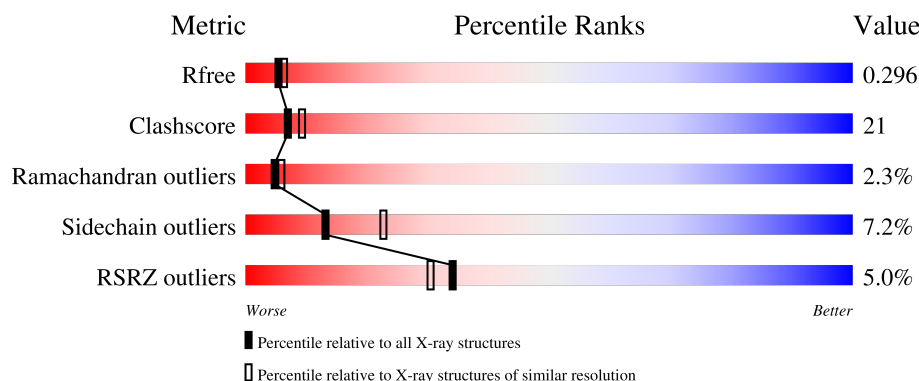
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	198	5% 57% 34% 5% • • •
1	B	198	6% 57% 33% 7% •
1	C	198	4% 63% 28% • • •

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	SO4	A	200	-	-	X	-
3	SO4	B	200	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4442 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 Hydroxypropylphosphonic Acid Epoxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	192	Total	C	N	O	S	0	0	0
			1440	912	248	275	5			
1	B	193	Total	C	N	O	S	0	0	0
			1446	915	250	276	5			
1	C	190	Total	C	N	O	S	0	0	0
			1411	895	242	269	5			

- Molecule 2 is FE (II) ION (CCD ID: FE2) (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		
2	C	1	Total	Fe	0	0
			1	1		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

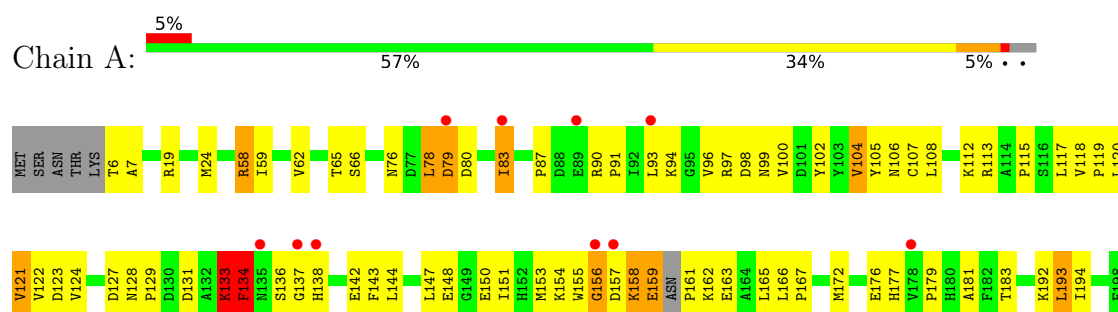
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	34	Total	O	0	0
			34	34		
4	B	49	Total	O	0	0
			49	49		
4	C	44	Total	O	0	0
			44	44		

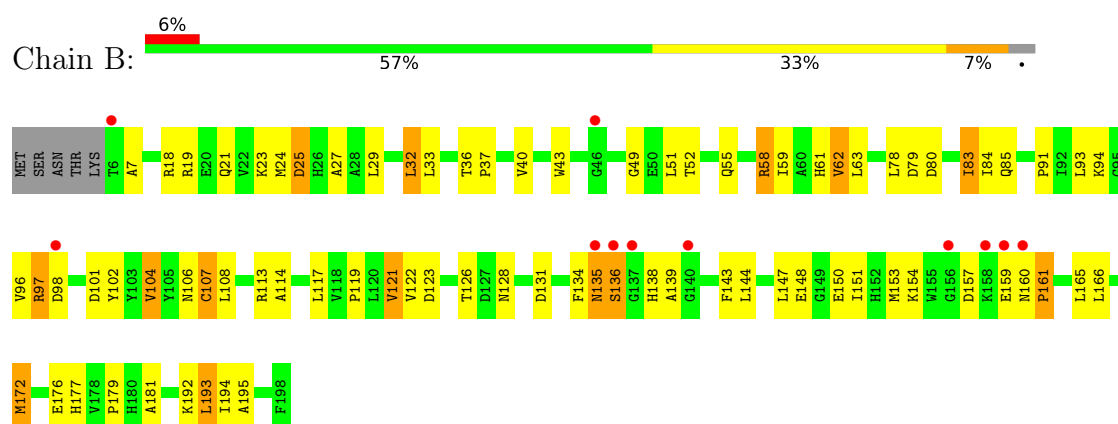
3 Residue-property plots [i](#)

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 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.

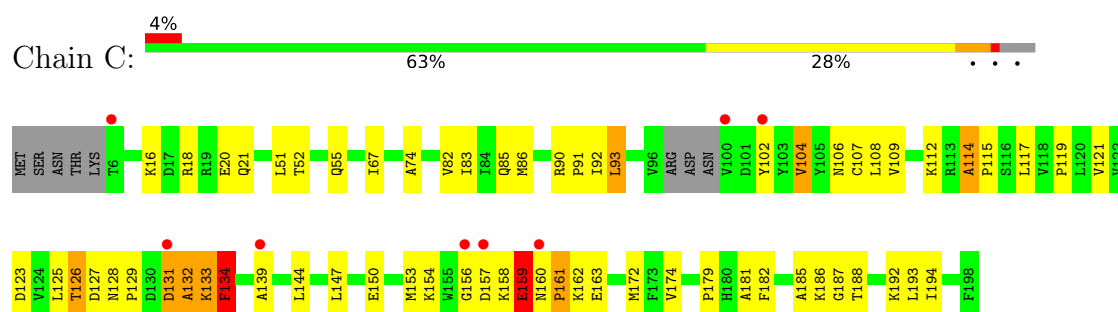
• Molecule 1: Hydroxypropylphosphonic Acid Epoxidase



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4 Data and refinement statistics

Property	Value	Source
Space group	P 42 21 2	Depositor
Cell constants a, b, c, α , β , γ	111.87Å 111.87Å 153.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.34 – 2.40 15.34 – 2.40	Depositor EDS
% Data completeness (in resolution range)	90.5 (15.34-2.40) 90.0 (15.34-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.00 (at 2.20Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.249 , 0.287 0.257 , 0.296	Depositor DCC
R_{free} test set	1737 reflections (3.70%)	wwPDB-VP
Wilson B-factor (Å ²)	40.6	Xtriage
Anisotropy	0.012	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 37.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4442	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 37.67 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.1420e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: SO4, FE2

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	0.51	0/1469	1.04	6/1996 (0.3%)
1	B	0.54	1/1476 (0.1%)	1.04	9/2008 (0.4%)
1	C	0.55	0/1440	1.03	7/1957 (0.4%)
All	All	0.53	1/4385 (0.0%)	1.04	22/5961 (0.4%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	172	MET	SD-CE	-5.41	1.66	1.79

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	7	ALA	N-CA-C	-7.91	102.26	112.23
1	C	139	ALA	N-CA-C	-6.49	95.69	107.75
1	A	7	ALA	N-CA-C	-6.38	104.24	111.14
1	C	83	ILE	CB-CA-C	6.09	119.91	110.28
1	B	83	ILE	CB-CA-C	6.09	119.90	110.28
1	C	114	ALA	CA-C-N	6.04	127.39	119.84
1	C	114	ALA	C-N-CA	6.04	127.39	119.84
1	B	114	ALA	CA-C-N	5.96	125.84	119.28
1	B	114	ALA	C-N-CA	5.96	125.84	119.28
1	A	156	GLY	N-CA-C	-5.94	99.11	113.18
1	C	109	VAL	N-CA-C	5.92	117.97	109.51
1	B	23	LYS	N-CA-C	5.75	119.05	111.28
1	B	121	VAL	N-CA-C	-5.71	99.51	107.80
1	C	125	LEU	N-CA-C	5.71	120.24	113.28
1	B	84	ILE	N-CA-C	5.64	116.00	108.11
1	A	121	VAL	N-CA-C	-5.58	99.12	107.37
1	A	96	VAL	N-CA-C	5.41	116.37	108.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	83	ILE	CB-CA-C	5.33	119.94	110.71
1	B	107	CYS	N-CA-C	5.24	117.29	109.59
1	C	67	ILE	N-CA-C	-5.22	105.30	110.62
1	B	122	VAL	N-CA-C	5.18	115.04	107.37
1	A	133	LYS	N-CA-C	5.15	116.58	110.97

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	1440	0	1419	63	0
1	B	1446	0	1417	72	0
1	C	1411	0	1381	49	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
3	A	5	0	0	2	0
3	B	5	0	0	2	0
3	C	5	0	0	0	0
4	A	34	0	0	2	0
4	B	49	0	0	4	0
4	C	44	0	0	0	0
All	All	4442	0	4217	180	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 21.

All (180) 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:156:GLY:O	1:A:179:PRO:HG2	1.52	1.08
1:A:166:LEU:HD11	1:A:172:MET:HE3	1.41	1.02
1:B:104:VAL:HG22	1:B:123:ASP:HB2	1.43	0.99

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:91:PRO:HG2	1:B:107:CYS:HB2	1.43	0.98
1:B:151:ILE:HD13	1:B:193:LEU:HD23	1.48	0.95
1:A:91:PRO:HG2	1:A:107:CYS:HB2	1.48	0.95
1:C:91:PRO:HG2	1:C:107:CYS:HB2	1.48	0.93
1:A:104:VAL:HG13	1:A:123:ASP:HB2	1.56	0.86
1:C:153:MET:HE3	1:C:172:MET:HE1	1.57	0.86
1:C:104:VAL:HG22	1:C:123:ASP:HB2	1.57	0.86
1:C:154:LYS:HG2	1:C:163:GLU:HG2	1.59	0.84
1:C:52:THR:OG1	1:C:55:GLN:HG3	1.78	0.83
1:C:150:GLU:HG2	1:C:185:ALA:HB2	1.61	0.82
1:B:123:ASP:OD1	1:B:192:LYS:HE3	1.81	0.81
1:A:192:LYS:NZ	3:A:200:SO4:O3	2.14	0.81
1:A:104:VAL:CG1	1:A:123:ASP:HB2	2.11	0.80
1:B:32:LEU:HD13	1:B:62:VAL:HG23	1.62	0.80
1:B:151:ILE:CD1	1:B:193:LEU:HD23	2.12	0.78
1:A:133:LYS:O	1:A:134:PHE:HB2	1.84	0.77
1:A:58:ARG:HD3	1:C:86:MET:HG3	1.65	0.76
1:A:76:ASN:ND2	1:A:113:ARG:HD3	2.01	0.76
1:C:159:GLU:O	1:C:161:PRO:HD3	1.86	0.75
1:A:87:PRO:HA	1:A:90:ARG:HD3	1.69	0.74
1:A:134:PHE:HA	1:A:181:ALA:HB2	1.69	0.74
1:B:166:LEU:HD11	1:B:172:MET:CE	2.21	0.71
1:B:166:LEU:HD11	1:B:172:MET:HE3	1.72	0.70
1:B:192:LYS:NZ	3:B:200:SO4:O4	2.22	0.69
1:B:59:ILE:O	1:B:63:LEU:HD13	1.91	0.69
1:A:123:ASP:OD1	1:A:192:LYS:HE2	1.93	0.69
1:A:78:LEU:HG	1:A:113:ARG:HG3	1.75	0.69
1:C:153:MET:HE3	1:C:172:MET:CE	2.24	0.67
3:A:200:SO4:O4	4:A:230:HOH:O	2.07	0.67
1:B:85:GLN:OE1	1:B:108:LEU:HA	1.96	0.66
1:B:33:LEU:HD21	1:B:59:ILE:HA	1.79	0.63
1:B:62:VAL:HG22	1:B:63:LEU:HD12	1.80	0.62
1:A:193:LEU:HD12	1:A:193:LEU:O	1.98	0.62
1:C:193:LEU:HD12	1:C:193:LEU:O	2.01	0.61
1:C:153:MET:CE	1:C:172:MET:HE1	2.29	0.61
1:C:185:ALA:HB3	1:C:188:THR:HG21	1.82	0.60
1:A:154:LYS:HG2	1:A:163:GLU:HG2	1.83	0.60
1:B:136:SER:O	1:B:179:PRO:HB3	2.02	0.60
1:C:106:ASN:HB2	1:C:121:VAL:HB	1.84	0.60
1:A:93:LEU:HD13	1:A:94:LYS:N	2.17	0.59
1:C:128:ASN:ND2	1:C:131:ASP:H	2.00	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:143:PHE:C	1:A:144:LEU:HD12	2.28	0.59
1:A:158:LYS:O	1:A:159:GLU:CD	2.46	0.58
1:B:93:LEU:HD12	1:B:94:LYS:H	1.68	0.58
1:A:91:PRO:HG2	1:A:107:CYS:CB	2.31	0.58
1:B:193:LEU:C	1:B:193:LEU:HD12	2.28	0.57
1:C:134:PHE:HA	1:C:181:ALA:HB2	1.85	0.57
1:C:158:LYS:HG3	1:C:159:GLU:OE1	2.05	0.57
1:B:147:LEU:HG	1:B:194:ILE:HG13	1.87	0.57
1:A:147:LEU:HB2	1:A:192:LYS:HG2	1.87	0.56
1:B:134:PHE:CZ	1:B:154:LYS:HB3	2.41	0.56
1:B:29:LEU:HD23	1:B:40:VAL:HG11	1.86	0.56
1:B:123:ASP:OD1	1:B:192:LYS:HG3	2.06	0.56
1:B:104:VAL:CG2	1:B:123:ASP:HB2	2.27	0.55
1:B:157:ASP:HB3	1:B:160:ASN:O	2.06	0.55
1:C:92:ILE:C	1:C:93:LEU:HD12	2.32	0.55
1:C:104:VAL:CG2	1:C:123:ASP:OD2	2.54	0.55
1:C:159:GLU:O	1:C:161:PRO:CD	2.53	0.55
1:A:150:GLU:CD	1:A:165:LEU:HD21	2.32	0.55
1:A:157:ASP:O	1:A:158:LYS:C	2.49	0.54
1:B:153:MET:HG2	1:B:172:MET:HE1	1.90	0.54
1:A:142:GLU:HG2	1:A:144:LEU:CD1	2.38	0.54
1:A:127:ASP:O	1:A:129:PRO:HD3	2.08	0.54
1:A:112:LYS:HZ1	1:B:61:HIS:CE1	2.26	0.53
1:C:93:LEU:HD12	1:C:93:LEU:N	2.24	0.53
1:A:166:LEU:CD1	1:A:172:MET:HE3	2.28	0.53
1:C:156:GLY:O	1:C:179:PRO:HG2	2.09	0.53
1:A:112:LYS:NZ	1:B:61:HIS:CE1	2.78	0.52
1:C:150:GLU:CG	1:C:185:ALA:HB2	2.38	0.52
1:B:52:THR:HG23	1:B:55:GLN:OE1	2.09	0.52
1:C:104:VAL:HG22	1:C:123:ASP:CB	2.36	0.52
1:A:108:LEU:HB2	1:A:119:PRO:HG2	1.90	0.52
1:C:102:TYR:HD1	1:C:126:THR:HG21	1.75	0.52
1:B:97:ARG:NH1	1:B:131:ASP:O	2.43	0.52
1:B:58:ARG:O	1:B:62:VAL:HG12	2.09	0.51
1:B:106:ASN:ND2	4:B:242:HOH:O	2.30	0.51
1:B:144:LEU:HD12	1:B:193:LEU:HD13	1.93	0.51
1:B:18:ARG:HA	1:B:21:GLN:HG2	1.93	0.51
1:B:128:ASN:ND2	1:B:128:ASN:C	2.69	0.51
1:A:133:LYS:O	1:A:134:PHE:CB	2.58	0.51
1:C:123:ASP:OD1	1:C:192:LYS:NZ	2.44	0.50
1:A:154:LYS:HE2	1:A:163:GLU:OE1	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:117:LEU:HG	1:A:119:PRO:HD3	1.94	0.50
1:A:157:ASP:O	1:A:159:GLU:N	2.45	0.50
1:A:158:LYS:O	1:A:159:GLU:HB2	2.12	0.49
1:C:114:ALA:N	1:C:115:PRO:HD3	2.27	0.49
1:C:131:ASP:O	1:C:132:ALA:C	2.55	0.49
1:B:128:ASN:C	1:B:128:ASN:HD22	2.18	0.49
1:B:93:LEU:HD12	1:B:94:LYS:N	2.27	0.49
1:B:134:PHE:HA	1:B:181:ALA:HB2	1.95	0.49
1:C:18:ARG:HH21	1:C:21:GLN:NE2	2.11	0.48
1:A:155:TRP:CZ2	1:A:162:LYS:HB3	2.48	0.48
1:A:136:SER:HA	1:A:179:PRO:HB3	1.94	0.48
1:B:78:LEU:CD1	1:B:113:ARG:HG3	2.43	0.48
1:B:25:ASP:OD1	1:B:27:ALA:HB3	2.14	0.48
1:A:158:LYS:O	1:A:159:GLU:CB	2.61	0.48
1:B:33:LEU:CD2	1:B:62:VAL:HG11	2.44	0.48
1:A:98:ASP:O	1:A:99:ASN:HB2	2.14	0.47
1:B:117:LEU:HG	1:B:119:PRO:HD3	1.95	0.47
1:C:90:ARG:HB2	1:C:107:CYS:O	2.14	0.47
1:B:150:GLU:OE2	1:B:165:LEU:HD11	2.14	0.47
1:A:122:VAL:HB	1:A:193:LEU:HD12	1.96	0.47
1:A:128:ASN:HB3	1:A:131:ASP:OD2	2.15	0.47
1:B:134:PHE:HZ	1:B:154:LYS:HB3	1.77	0.47
1:B:166:LEU:HD11	1:B:172:MET:HE1	1.96	0.47
1:C:51:LEU:HA	1:C:55:GLN:OE1	2.15	0.47
1:C:108:LEU:HD21	1:C:121:VAL:HG23	1.96	0.47
1:A:106:ASN:HB2	1:A:121:VAL:HB	1.96	0.47
1:B:97:ARG:HG3	1:B:102:TYR:CD1	2.50	0.47
1:A:93:LEU:HD21	1:B:25:ASP:OD2	2.15	0.47
1:A:124:VAL:HG11	1:A:183:THR:HA	1.96	0.47
1:B:101:ASP:HB2	4:B:239:HOH:O	2.14	0.47
1:A:151:ILE:HG13	1:A:193:LEU:HD23	1.95	0.46
1:B:139:ALA:HA	1:B:177:HIS:CE1	2.50	0.46
1:A:97:ARG:O	1:A:98:ASP:HB2	2.15	0.46
1:B:143:PHE:C	1:B:144:LEU:HD22	2.40	0.46
1:B:79:ASP:OD1	1:B:79:ASP:C	2.58	0.46
1:B:18:ARG:HA	1:B:21:GLN:HE21	1.81	0.46
1:C:127:ASP:OD2	1:C:187:GLY:HA2	2.15	0.46
1:C:172:MET:HE3	1:C:174:VAL:CG2	2.45	0.46
1:C:16:LYS:HE2	1:C:20:GLU:OE1	2.16	0.45
1:C:147:LEU:HD12	1:C:194:ILE:HG13	1.98	0.45
1:C:154:LYS:HA	1:C:162:LYS:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:74:ALA:HB1	1:C:112:LYS:HB3	1.99	0.45
1:A:87:PRO:O	1:A:90:ARG:HG2	2.16	0.45
1:B:19:ARG:HH11	1:B:19:ARG:HG2	1.82	0.45
1:B:106:ASN:HB2	1:B:121:VAL:HB	1.98	0.45
1:A:59:ILE:O	1:A:62:VAL:HG22	2.17	0.45
1:A:78:LEU:CG	1:A:113:ARG:HG3	2.45	0.45
1:B:135:ASN:O	1:B:135:ASN:CG	2.59	0.44
1:B:36:THR:O	1:B:37:PRO:C	2.61	0.44
1:B:135:ASN:O	1:B:136:SER:C	2.60	0.44
1:C:117:LEU:CD1	1:C:119:PRO:HD3	2.47	0.44
1:B:43:TRP:CE2	1:B:51:LEU:HD21	2.52	0.44
1:A:79:ASP:O	1:A:80:ASP:HB2	2.17	0.44
1:B:176:GLU:O	1:B:177:HIS:HB2	2.18	0.44
1:B:62:VAL:HG13	1:B:63:LEU:HD13	2.00	0.44
1:A:65:THR:HG23	1:A:66:SER:N	2.33	0.43
1:B:19:ARG:HG2	1:B:24:MET:O	2.17	0.43
1:A:115:PRO:HA	4:B:231:HOH:O	2.17	0.43
1:B:33:LEU:HD21	1:B:62:VAL:HG11	2.00	0.43
1:B:193:LEU:HD12	1:B:193:LEU:O	2.19	0.43
1:A:100:VAL:O	1:A:102:TYR:HD1	2.02	0.43
1:B:138:HIS:O	1:B:177:HIS:N	2.51	0.43
1:C:90:ARG:HA	1:C:91:PRO:HD2	1.84	0.43
1:B:148:GLU:HB2	4:B:220:HOH:O	2.19	0.43
1:A:153:MET:HG2	1:A:172:MET:HE1	2.00	0.43
1:A:176:GLU:HG3	1:A:177:HIS:CD2	2.54	0.43
1:C:128:ASN:O	1:C:131:ASP:N	2.52	0.43
1:A:105:TYR:HD2	1:A:122:VAL:HG22	1.84	0.42
1:A:19:ARG:HG2	1:A:24:MET:O	2.18	0.42
1:A:148:GLU:HB2	1:A:192:LYS:HB3	2.02	0.42
1:C:185:ALA:O	1:C:186:LYS:C	2.62	0.42
1:A:143:PHE:O	1:A:144:LEU:HD12	2.19	0.42
1:B:135:ASN:O	1:B:135:ASN:OD1	2.38	0.42
1:C:144:LEU:HD23	1:C:193:LEU:HD22	2.00	0.42
1:C:93:LEU:N	1:C:93:LEU:CD1	2.83	0.42
1:A:118:VAL:HG11	1:B:21:GLN:O	2.19	0.42
1:A:159:GLU:HA	1:A:159:GLU:OE1	2.20	0.42
1:B:63:LEU:HD12	1:B:63:LEU:N	2.34	0.41
1:A:94:LYS:NZ	4:A:201:HOH:O	2.53	0.41
1:A:159:GLU:O	1:A:161:PRO:HD3	2.20	0.41
1:C:144:LEU:HD22	1:C:172:MET:HE2	2.02	0.41
1:C:153:MET:HE1	1:C:182:PHE:CD1	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:85:GLN:OE1	1:C:108:LEU:HA	2.20	0.41
1:A:150:GLU:HG3	1:A:167:PRO:HA	2.02	0.41
1:B:192:LYS:NZ	3:B:200:SO4:S	2.82	0.41
1:C:157:ASP:OD2	1:C:159:GLU:HB2	2.21	0.41
1:A:120:LEU:O	1:A:194:ILE:HG13	2.21	0.41
1:C:129:PRO:O	1:C:154:LYS:NZ	2.54	0.41
1:C:153:MET:SD	1:C:172:MET:HE1	2.61	0.41
1:B:104:VAL:CG2	1:B:104:VAL:O	2.67	0.40
1:B:121:VAL:CG1	1:B:192:LYS:HE2	2.51	0.40
1:B:94:LYS:HD3	1:B:101:ASP:OD1	2.20	0.40
1:B:144:LEU:HD13	1:B:195:ALA:HA	2.04	0.40
1:B:29:LEU:HD13	1:B:63:LEU:HD11	2.03	0.40
1:B:97:ARG:O	1:B:98:ASP:C	2.64	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	188/198 (95%)	171 (91%)	12 (6%)	5 (3%)	4	4
1	B	191/198 (96%)	181 (95%)	7 (4%)	3 (2%)	7	11
1	C	186/198 (94%)	173 (93%)	8 (4%)	5 (3%)	4	4
All	All	565/594 (95%)	525 (93%)	27 (5%)	13 (2%)	5	6

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	136	SER
1	C	133	LYS
1	C	159	GLU
1	A	78	LEU

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Mol	Chain	Res	Type
1	A	158	LYS
1	C	134	PHE
1	A	79	ASP
1	A	134	PHE
1	C	132	ALA
1	A	137	GLY
1	B	49	GLY
1	B	161	PRO
1	C	160	ASN

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	150/160 (94%)	141 (94%)	9 (6%)	17	31
1	B	149/160 (93%)	135 (91%)	14 (9%)	8	13
1	C	145/160 (91%)	136 (94%)	9 (6%)	16	29
All	All	444/480 (92%)	412 (93%)	32 (7%)	13	23

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	THR
1	A	58	ARG
1	A	83	ILE
1	A	104	VAL
1	A	133	LYS
1	A	134	PHE
1	A	138	HIS
1	A	159	GLU
1	A	193	LEU
1	B	25	ASP
1	B	32	LEU
1	B	58	ARG
1	B	62	VAL

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Mol	Chain	Res	Type
1	B	80	ASP
1	B	83	ILE
1	B	96	VAL
1	B	97	ARG
1	B	104	VAL
1	B	126	THR
1	B	135	ASN
1	B	159	GLU
1	B	161	PRO
1	B	193	LEU
1	C	82	VAL
1	C	93	LEU
1	C	104	VAL
1	C	126	THR
1	C	131	ASP
1	C	133	LYS
1	C	134	PHE
1	C	159	GLU
1	C	161	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	GLN
1	A	76	ASN
1	B	21	GLN
1	B	45	ASN
1	B	128	ASN
1	B	152	HIS
1	B	177	HIS
1	C	21	GLN
1	C	141	ASN
1	C	197	ASN

5.3.3 RNA ⓘ

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 3 are monoatomic - leaving 3 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	SO4	B	200	-	4,4,4	0.65	0	6,6,6	0.64	0
3	SO4	A	200	-	4,4,4	0.62	0	6,6,6	0.63	0
3	SO4	C	200	-	4,4,4	0.39	0	6,6,6	0.56	0

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.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	200	SO4	2	0
3	A	200	SO4	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	192/198 (96%)	0.31	10 (5%) 33 29	19, 41, 62, 70	0
1	B	193/198 (97%)	0.25	11 (5%) 29 25	17, 38, 57, 66	0
1	C	190/198 (95%)	0.21	8 (4%) 40 36	17, 38, 64, 74	0
All	All	575/594 (96%)	0.26	29 (5%) 34 30	17, 39, 61, 74	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	131	ASP	4.8
1	B	136	SER	4.8
1	C	100	VAL	4.5
1	A	135	ASN	3.4
1	C	160	ASN	3.4
1	B	135	ASN	3.3
1	A	156	GLY	3.3
1	A	83	ILE	3.1
1	B	98	ASP	2.7
1	B	140	GLY	2.7
1	C	102	TYR	2.7
1	B	160	ASN	2.6
1	A	137	GLY	2.5
1	A	178	VAL	2.4
1	A	93	LEU	2.4
1	B	137	GLY	2.3
1	A	138	HIS	2.2
1	A	89	GLU	2.2
1	C	6	THR	2.2
1	A	157	ASP	2.2
1	B	158	LYS	2.2
1	B	46	GLY	2.1
1	B	156	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	156	GLY	2.1
1	B	159	GLU	2.1
1	A	79	ASP	2.1
1	C	139	ALA	2.0
1	B	6	THR	2.0
1	C	157	ASP	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 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($\text{\AA}^2$)	Q<0.9
3	SO4	C	200	5/5	0.73	0.15	97,97,98,99	0
3	SO4	B	200	5/5	0.87	0.10	71,73,73,74	0
3	SO4	A	200	5/5	0.90	0.09	65,66,67,68	0
2	FE2	A	199	1/1	0.95	0.06	45,45,45,45	0
2	FE2	C	199	1/1	0.99	0.02	47,47,47,47	0
2	FE2	B	199	1/1	0.99	0.02	34,34,34,34	0

6.5 Other polymers [i](#)

There are no such residues in this entry.