



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 02:14 AM GMT

PDB ID : 3PQ3
Title : Structure of I274C variant of E. coli KatE[] - Images 7-12
Authors : Loewen, P.C.; Jha, V.; Louis, S.; Chelikani, P.; Carpena, X.; Fita, I.
Deposited on : 2010-11-25
Resolution : 1.79 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

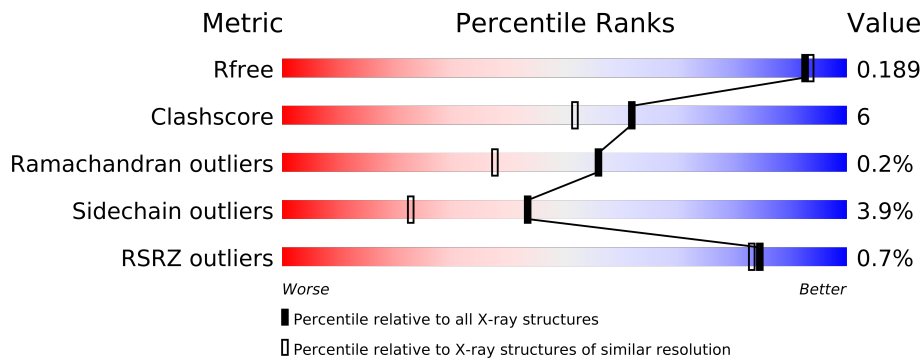
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.79 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}	66092	3513 (1.80-1.80)
Clashscore	79885	4461 (1.80-1.80)
Ramachandran outliers	78287	4404 (1.80-1.80)
Sidechain outliers	78261	4403 (1.80-1.80)
RSRZ outliers	66119	3515 (1.80-1.80)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	753	
1	B	753	
1	C	753	
1	D	753	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
4	H2S	A	754	-	X
4	H2S	B	754	-	X
4	H2S	C	754	-	X
4	H2S	D	754	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 26828 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

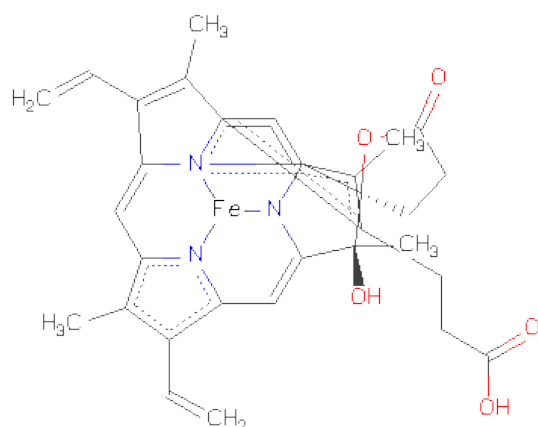
- Molecule 1 is a protein called Catalase HP11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	726	Total	C	N	O	S	0	5	0
			5755	3652	1009	1083	11			
1	B	726	Total	C	N	O	S	0	5	0
			5757	3654	1009	1083	11			
1	C	726	Total	C	N	O	S	0	2	0
			5747	3649	1007	1080	11			
1	D	726	Total	C	N	O	S	0	9	0
			5771	3662	1013	1085	11			

There are 12 discrepancies between the modelled and reference sequences:

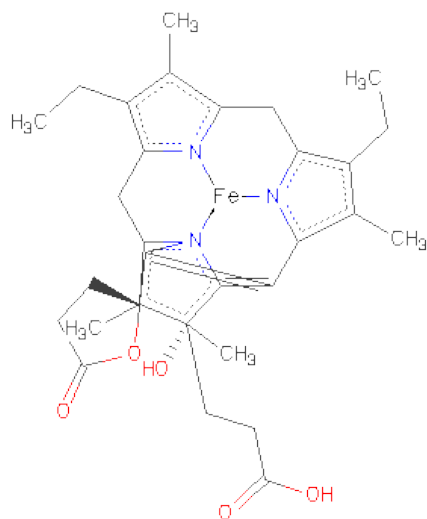
Chain	Residue	Modelled	Actual	Comment	Reference
A	274	CYS	ILE	ENGINEERED MUTATION	UNP P21179
A	438	ALA	CYS	ENGINEERED MUTATION	UNP P21179
A	669	ALA	CYS	ENGINEERED MUTATION	UNP P21179
B	274	CYS	ILE	ENGINEERED MUTATION	UNP P21179
B	438	ALA	CYS	ENGINEERED MUTATION	UNP P21179
B	669	ALA	CYS	ENGINEERED MUTATION	UNP P21179
C	274	CYS	ILE	ENGINEERED MUTATION	UNP P21179
C	438	ALA	CYS	ENGINEERED MUTATION	UNP P21179
C	669	ALA	CYS	ENGINEERED MUTATION	UNP P21179
D	274	CYS	ILE	ENGINEERED MUTATION	UNP P21179
D	438	ALA	CYS	ENGINEERED MUTATION	UNP P21179
D	669	ALA	CYS	ENGINEERED MUTATION	UNP P21179

- Molecule 2 is CIS-HEME D HYDROXYCHLORIN GAMMA-SPIROLACTONE (three-letter code: HDD) (formula: C₃₄H₃₂FeN₄O₅).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	1
			44	34	1	4	5		
2	B	1	Total	C	Fe	N	O	0	1
			44	34	1	4	5		
2	C	1	Total	C	Fe	N	O	0	1
			44	34	1	4	5		
2	D	1	Total	C	Fe	N	O	0	1
			44	34	1	4	5		

- Molecule 3 is CIS-HEME D HYDROXYCHLORIN GAMMA-SPIROLACTONE 17R, 18S (three-letter code: HDE) (formula: $C_{34}H_{38}FeN_4O_5$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 44	C 34	Fe 1	N 4	O 5	0	1
3	B	1	Total 44	C 34	Fe 1	N 4	O 5	0	1
3	C	1	Total 44	C 34	Fe 1	N 4	O 5	0	1
3	D	1	Total 44	C 34	Fe 1	N 4	O 5	0	1

- Molecule 4 is HYDROSULFURIC ACID (three-letter code: H₂S) (formula: H₂S).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total 1	S 1	0	0
4	B	1	Total 1	S 1	0	0
4	C	1	Total 1	S 1	0	0
4	D	1	Total 1	S 1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	895	Total 895	O 895	0	0
5	B	791	Total 791	O 791	0	0

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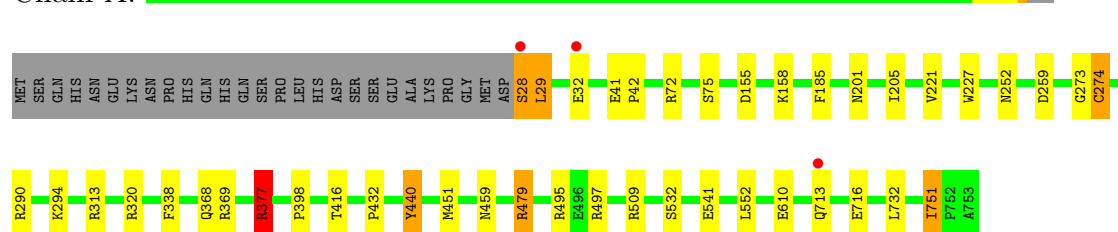
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	842	Total 842	O 842	0	0
5	D	914	Total 914	O 914	0	0

3 Residue-property plots

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.

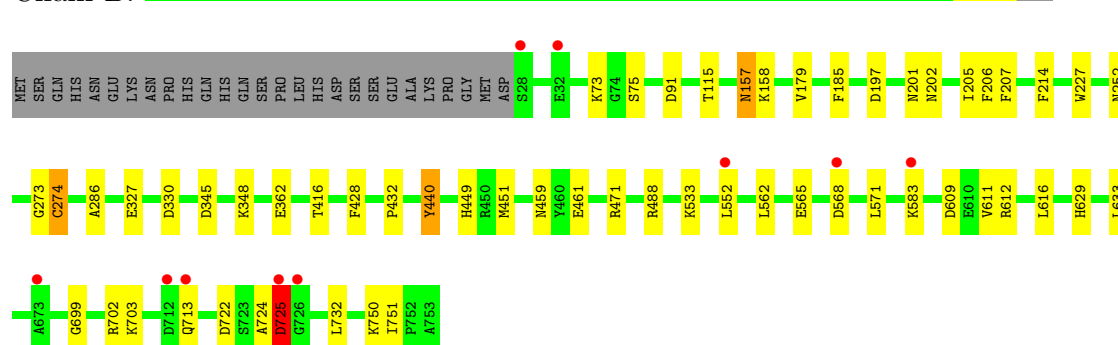
• Molecule 1: Catalase HP1I

Chain A:



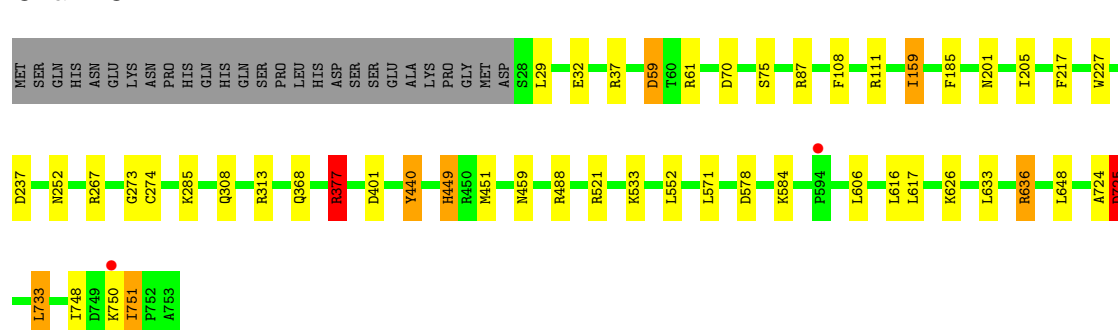
• Molecule 1: Catalase HP1I

Chain B:



• Molecule 1: Catalase HP1I

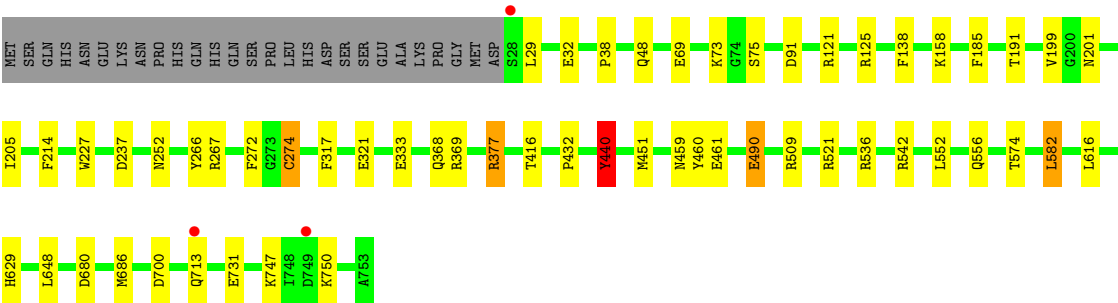
Chain C:



• Molecule 1: Catalase HP1I

Chain D:





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	93.51Å 133.03Å 122.65Å 90.00° 109.39° 90.00°	Depositor
Resolution (Å)	35.23 – 1.79 35.23 – 1.79	Depositor EDS
% Data completeness (in resolution range)	92.3 (35.23-1.79) 92.3 (35.23-1.79)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.16	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.51 (at 1.79Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.142 , 0.189 0.143 , 0.189	Depositor DCC
R_{free} test set	12198 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	11.4	Xtriage
Anisotropy	0.019	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 33.9	EDS
Estimated twinning fraction	0.028 for h,-k,-h-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 243340 reflections (0.000%)	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	26828	wwPDB-VP
Average B, all atoms (Å ²)	13.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.44% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: HDE, HDD, H2S

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	1.17	5/5932 (0.1%)	0.97	11/8064 (0.1%)
1	B	1.12	6/5931 (0.1%)	0.94	4/8062 (0.0%)
1	C	1.13	2/5908 (0.0%)	0.96	16/8033 (0.2%)
1	D	1.19	11/5966 (0.2%)	0.96	11/8109 (0.1%)
All	All	1.15	24/23737 (0.1%)	0.96	42/32268 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	3

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	440	TYR	CE1-CZ	7.19	1.47	1.38
1	B	440	TYR	CE1-CZ	6.39	1.46	1.38
1	B	362	GLU	CB-CG	6.08	1.63	1.52
1	B	327	GLU	CB-CG	5.98	1.63	1.52
1	D	138	PHE	CD1-CE1	5.89	1.51	1.39
1	C	108	PHE	CE2-CZ	5.75	1.48	1.37
1	B	428	PHE	CE2-CZ	5.66	1.48	1.37
1	D	272	PHE	CE2-CZ	5.65	1.48	1.37
1	A	440	TYR	CE1-CZ	5.64	1.45	1.38
1	D	266	TYR	CD2-CE2	5.63	1.47	1.39
1	D	214	PHE	CE2-CZ	5.62	1.48	1.37
1	D	199	VAL	CB-CG1	5.59	1.64	1.52
1	B	179	VAL	CB-CG2	5.58	1.64	1.52
1	D	333	GLU	CD-OE2	-5.54	1.19	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	317	PHE	CE2-CZ	5.46	1.47	1.37
1	D	69	GLU	CG-CD	5.29	1.59	1.51
1	B	286	ALA	CA-CB	5.28	1.63	1.52
1	D	460	TYR	CD1-CE1	5.11	1.47	1.39
1	A	221	VAL	CB-CG2	5.09	1.63	1.52
1	C	217	PHE	CB-CG	-5.09	1.42	1.51
1	A	338	PHE	CE2-CZ	5.08	1.47	1.37
1	D	440	TYR	CD2-CE2	5.06	1.47	1.39
1	A	377	ARG	CZ-NH2	5.02	1.39	1.33
1	A	41	GLU	CG-CD	5.01	1.59	1.51

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	636	ARG	NE-CZ-NH2	-8.87	115.87	120.30
1	C	636	ARG	NE-CZ-NH1	8.11	124.36	120.30
1	A	495	ARG	NE-CZ-NH2	-7.90	116.35	120.30
1	A	495	ARG	NE-CZ-NH1	7.59	124.09	120.30
1	C	377	ARG	NE-CZ-NH1	-7.51	116.55	120.30
1	B	609	ASP	CB-CG-OD2	6.88	124.50	118.30
1	C	111	ARG	NE-CZ-NH1	-6.87	116.86	120.30
1	D	686	MET	CG-SD-CE	6.82	111.10	100.20
1	A	320	ARG	NE-CZ-NH2	-6.67	116.97	120.30
1	A	294	LYS	CD-CE-NZ	-6.35	97.09	111.70
1	A	479	ARG	NE-CZ-NH1	6.32	123.46	120.30
1	C	401	ASP	CB-CG-OD2	6.27	123.94	118.30
1	C	733	LEU	CA-CB-CG	6.19	129.53	115.30
1	C	449[A]	HIS	CA-C-O	-6.12	107.26	120.10
1	C	449[B]	HIS	CA-C-O	-6.12	107.26	120.10
1	D	377	ARG	NE-CZ-NH1	-6.05	117.28	120.30
1	C	87	ARG	NE-CZ-NH2	-6.04	117.28	120.30
1	C	37	ARG	NE-CZ-NH2	-6.04	117.28	120.30
1	D	125	ARG	NE-CZ-NH1	6.02	123.31	120.30
1	D	121	ARG	NE-CZ-NH1	5.84	123.22	120.30
1	C	59	ASP	CB-CG-OD1	5.79	123.51	118.30
1	D	582	LEU	CB-CG-CD1	5.75	120.78	111.00
1	C	159	ILE	CB-CG1-CD1	-5.54	98.38	113.90
1	C	70	ASP	CB-CG-OD2	5.54	123.29	118.30
1	D	125	ARG	NE-CZ-NH2	-5.52	117.54	120.30
1	D	536	ARG	NE-CZ-NH1	-5.49	117.56	120.30
1	A	497	ARG	NE-CZ-NH2	-5.48	117.56	120.30
1	D	509	ARG	NE-CZ-NH1	5.48	123.04	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	197	ASP	CB-CG-OD2	5.44	123.19	118.30
1	C	313	ARG	NE-CZ-NH1	5.42	123.01	120.30
1	C	725	ASP	N-CA-C	5.37	125.51	111.00
1	A	290	ARG	NE-CZ-NH1	-5.35	117.62	120.30
1	D	521	ARG	NE-CZ-NH2	-5.34	117.63	120.30
1	B	471	ARG	NE-CZ-NH2	-5.32	117.64	120.30
1	C	313	ARG	NE-CZ-NH2	-5.31	117.64	120.30
1	A	72	ARG	NE-CZ-NH2	-5.29	117.65	120.30
1	A	313	ARG	NE-CZ-NH1	-5.27	117.67	120.30
1	A	479	ARG	NE-CZ-NH2	-5.23	117.69	120.30
1	D	680	ASP	CB-CG-OD2	5.21	122.99	118.30
1	D	542	ARG	NE-CZ-NH2	-5.17	117.71	120.30
1	A	259	ASP	CB-CG-OD2	5.16	122.95	118.30
1	B	609	ASP	CB-CG-OD1	-5.16	113.65	118.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	449[A]	HIS	Mainchain
1	C	724	ALA	Peptide
1	C	725	ASP	Peptide

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5755	0	0	25	0
1	B	5757	0	0	30	0
1	C	5747	0	0	21	1
1	D	5771	0	0	16	0
2	A	44	0	0	5	0
2	B	44	0	0	7	0
2	C	44	0	0	5	0
2	D	44	0	0	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	44	0	36	5	0
3	B	44	0	36	12	0
3	C	44	0	36	11	0
3	D	44	0	36	10	0
4	A	1	0	0	4	0
4	B	1	0	0	5	0
4	C	1	0	0	2	0
4	D	1	0	0	5	0
5	A	895	0	0	11	2
5	B	791	0	0	9	1
5	C	842	0	0	15	0
5	D	914	0	0	14	0
All	All	26828	0	144	133	2

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 6.

All (133) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:274:CYS:CB	4:D:754:H2S:S	2.15	1.34
2:B:760[A]:HDD:CBB	4:B:754:H2S:S	2.17	1.32
1:B:449[A]:HIS:CE1	5:B:3618:HOH:O	1.65	1.28
3:B:761[B]:HDE:CBC	3:B:761[B]:HDE:HMC	1.61	1.27
1:B:274:CYS:CB	4:B:754:H2S:S	2.23	1.26
1:A:274:CYS:CB	4:A:754:H2S:S	2.25	1.25
3:B:761[B]:HDE:HBCA	3:B:761[B]:HDE:CMC	1.65	1.22
3:C:761[B]:HDE:CBC	3:C:761[B]:HDE:HMC	1.75	1.15
1:D:451:MET:SD	5:D:3617:HOH:O	2.05	1.14
1:C:274:CYS:CB	4:C:754:H2S:S	2.36	1.13
1:C:451:MET:SD	5:C:3612:HOH:O	2.07	1.11
3:D:761[B]:HDE:HMC	3:D:761[B]:HDE:HBCB	1.32	1.10
2:D:760[A]:HDD:CBB	4:D:754:H2S:S	2.40	1.09
3:C:761[B]:HDE:HBCB	3:C:761[B]:HDE:HMC	1.06	1.06
3:C:761[B]:HDE:CMC	3:C:761[B]:HDE:HBCB	1.82	1.05
1:B:451:MET:SD	5:B:3614:HOH:O	2.16	1.03
1:A:274:CYS:SG	2:A:760[A]:HDD:CMB	2.53	0.95
1:A:28:SER:O	1:A:28:SER:OG	1.82	0.94
1:B:449[A]:HIS:NE2	5:B:3618:HOH:O	1.82	0.91
3:A:761[B]:HDE:HBCB	3:A:761[B]:HDE:HMC	1.52	0.91
3:B:761[B]:HDE:HBCA	3:B:761[B]:HDE:HMC	0.90	0.90
2:A:760[A]:HDD:CBB	4:A:754:H2S:S	2.61	0.89
1:D:274:CYS:SG	2:D:760[A]:HDD:CMB	2.62	0.87

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:C:761[B]:HDE:CBC	3:C:761[B]:HDE:CMC	2.44	0.86
3:B:761[B]:HDE:CMC	3:B:761[B]:HDE:CBC	2.33	0.84
1:B:73:LYS:CE	5:D:3517:HOH:O	2.25	0.83
3:B:761[B]:HDE:HMB	3:B:761[B]:HDE:HBBB	1.61	0.81
3:D:761[B]:HDE:CMC	3:D:761[B]:HDE:HBCB	2.11	0.80
1:C:274:CYS:SG	2:C:760[A]:HDD:CMB	2.70	0.80
1:D:731:GLU:OE2	5:D:3028:HOH:O	2.02	0.78
3:B:761[B]:HDE:CBB	3:B:761[B]:HDE:HMB	2.16	0.76
1:A:716:GLU:CG	5:A:3090:HOH:O	2.36	0.74
3:D:761[B]:HDE:HMB	3:D:761[B]:HDE:CBB	2.17	0.74
1:A:451:MET:SD	5:A:3609:HOH:O	2.47	0.72
1:C:725:ASP:O	5:C:2403:HOH:O	2.08	0.72
3:D:761[B]:HDE:HMC	4:D:754:H2S:S	2.31	0.70
1:A:532:SER:OG	5:A:2410:HOH:O	2.09	0.69
1:C:59:ASP:OD2	5:C:2529:HOH:O	2.11	0.69
1:C:267:ARG:CG	5:C:2916:HOH:O	2.42	0.68
1:A:377:ARG:NH1	5:A:1001:HOH:O	2.27	0.67
1:C:578:ASP:CB	5:C:2919:HOH:O	2.42	0.67
3:D:761[B]:HDE:HMB	3:D:761[B]:HDE:HBBA	1.78	0.66
1:A:751:ILE:O	1:A:751:ILE:CD1	2.45	0.65
3:A:761[B]:HDE:HBBB	3:A:761[B]:HDE:HMB	1.77	0.65
3:C:761[B]:HDE:HMB	3:C:761[B]:HDE:CBB	2.28	0.63
3:B:761[B]:HDE:HBCB	3:B:761[B]:HDE:HMC	1.74	0.63
1:C:59:ASP:OD2	5:C:3406:HOH:O	2.15	0.63
1:B:416:THR:CG2	5:D:2464:HOH:O	2.47	0.63
3:B:761[B]:HDE:CMC	3:B:761[B]:HDE:HBCB	2.28	0.62
1:B:533[A]:LYS:CE	5:B:3100:HOH:O	2.47	0.61
3:C:761[B]:HDE:HMC	4:C:754:H2S:S	2.41	0.61
3:B:761[B]:HDE:HMC	4:B:754:H2S:S	2.39	0.61
1:B:274:CYS:SG	2:B:760[A]:HDD:CMB	2.89	0.61
1:A:610:GLU:OE1	5:A:2394:HOH:O	2.16	0.61
1:C:368:GLN:NE2	5:C:1787:HOH:O	2.33	0.60
1:D:556:GLN:NE2	5:D:2773:HOH:O	2.33	0.60
1:A:610:GLU:CG	5:A:3555:HOH:O	2.49	0.60
1:D:158:LYS:NZ	5:D:2950:HOH:O	2.35	0.59
3:C:761[B]:HDE:CMC	3:C:761[B]:HDE:HBCA	2.31	0.59
3:D:761[B]:HDE:HBBB	3:D:761[B]:HDE:HMB	1.84	0.58
1:A:509:ARG:NE	5:A:3173:HOH:O	2.36	0.58
1:B:73:LYS:CD	5:D:3517:HOH:O	2.49	0.57
1:B:611:VAL:N	5:B:1760:HOH:O	2.36	0.57
1:B:201:ASN:ND2	2:B:760[A]:HDD:CMB	2.67	0.57
2:C:760[A]:HDD:CBB	2:C:760[A]:HDD:CMB	2.84	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:201:ASN:CG	2:A:760[A]:HDD:CMB	2.74	0.56
1:C:533:LYS:NZ	5:C:2231:HOH:O	2.40	0.55
1:A:541:GLU:OE2	5:A:2550:HOH:O	2.17	0.55
1:C:748:ILE:O	1:C:751:ILE:CG2	2.55	0.55
1:B:201:ASN:CG	2:B:760[A]:HDD:CMB	2.75	0.54
1:B:115:THR:CG2	5:B:3420:HOH:O	2.56	0.54
1:D:368[B]:GLN:NE2	5:D:1808:HOH:O	2.40	0.54
1:C:274:CYS:CB	2:C:760[A]:HDD:CBB	2.87	0.53
1:C:488:ARG:NH1	5:C:2379:HOH:O	2.40	0.53
3:D:761[B]:HDE:HBCB	4:D:754:H2S:S	2.49	0.53
2:A:760[A]:HDD:CAB	4:A:754:H2S:S	2.97	0.53
3:A:761[B]:HDE:CBB	3:A:761[B]:HDE:HMB	2.37	0.53
3:C:761[B]:HDE:HMB	3:C:761[B]:HDE:HBBB	1.90	0.53
1:D:29:LEU:N	5:D:2467:HOH:O	2.42	0.53
1:A:273:GLY:C	1:A:274:CYS:SG	2.87	0.53
1:D:201:ASN:CG	2:D:760[A]:HDD:CMB	2.78	0.52
3:C:761[B]:HDE:HMB	3:C:761[B]:HDE:HBBA	1.92	0.51
1:C:440:TYR:CD1	5:C:3454:HOH:O	2.55	0.51
1:A:201:ASN:ND2	2:A:760[A]:HDD:CMB	2.74	0.50
2:D:760[A]:HDD:CAB	4:D:754:H2S:S	2.95	0.50
1:A:416[A]:THR:CG2	5:C:3313:HOH:O	2.59	0.50
1:B:273:GLY:C	1:B:274:CYS:SG	2.90	0.49
1:C:201:ASN:CG	2:C:760[A]:HDD:CMB	2.80	0.49
3:B:761[B]:HDE:CMC	4:B:754:H2S:S	3.01	0.49
1:C:29:LEU:N	5:C:2891:HOH:O	2.46	0.49
1:C:201:ASN:ND2	2:C:760[A]:HDD:CMB	2.76	0.49
3:A:761[B]:HDE:CBC	3:A:761[B]:HDE:HMC	2.35	0.48
1:B:724:ALA:O	1:B:725:ASP:O	2.31	0.48
1:D:440:TYR:CZ	5:D:3517:HOH:O	2.55	0.48
3:D:761[B]:HDE:CMC	3:D:761[B]:HDE:CBC	2.85	0.48
1:B:449[B]:HIS:CE1	5:D:1789:HOH:O	2.67	0.47
1:D:629:HIS:CD2	5:D:1554:HOH:O	2.66	0.47
3:B:761[B]:HDE:HMAA	3:B:761[B]:HDE:HAAA	1.75	0.47
1:D:490:GLU:OE1	5:D:2308:HOH:O	2.20	0.47
3:D:761[B]:HDE:HMC	3:D:761[B]:HDE:CBC	2.22	0.46
3:C:761[B]:HDE:HAAA	3:C:761[B]:HDE:HMAA	1.76	0.46
3:D:761[B]:HDE:HMAA	3:D:761[B]:HDE:HAAA	1.70	0.46
1:A:451:MET:CE	1:C:451:MET:CE	2.94	0.45
1:A:751:ILE:CD1	1:A:751:ILE:C	2.85	0.45
1:A:274:CYS:CA	4:A:754:H2S:S	3.00	0.45
1:B:274:CYS:CB	2:B:760[A]:HDD:CBB	2.95	0.45
1:B:612:ARG:NH2	1:B:722:ASP:O	2.50	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:29:LEU:CD2	5:C:2405:HOH:O	2.65	0.44
1:A:479:ARG:NH2	5:A:2607:HOH:O	2.50	0.44
1:B:488:ARG:NH2	5:B:3477:HOH:O	2.51	0.44
1:C:308:GLN:NE2	5:C:3278:HOH:O	2.50	0.44
1:C:273:GLY:C	1:C:274:CYS:SG	2.96	0.44
1:B:157:ASN:C	1:B:157:ASN:ND2	2.71	0.44
1:B:699:GLY:O	1:B:702:ARG:NE	2.51	0.43
3:C:761[B]:HDE:HBCA	3:C:761[B]:HDE:HMC	1.82	0.43
1:B:73:LYS:NZ	5:B:1110:HOH:O	2.51	0.43
1:A:155:ASP:CG	5:A:3378:HOH:O	2.56	0.43
1:B:330:ASP:OD1	1:B:629:HIS:CE1	2.72	0.43
2:B:760[A]:HDD:CMB	2:B:760[A]:HDD:CBB	2.97	0.43
1:B:202:ASN:ND2	5:B:3407:HOH:O	2.52	0.43
3:B:761[B]:HDE:CBC	4:B:754:H2S:S	3.07	0.42
1:D:267:ARG:NH1	1:D:321:GLU:OE2	2.51	0.42
1:B:461:GLU:OE1	1:D:91:ASP:OD1	2.37	0.42
1:B:725:ASP:C	1:B:725:ASP:OD2	2.57	0.42
1:A:368:GLN:NE2	5:A:1785:HOH:O	2.53	0.42
1:B:91:ASP:OD1	1:D:461:GLU:OE1	2.39	0.41
1:B:345:ASP:OD1	1:B:348:LYS:NZ	2.54	0.41
1:B:206:PHE:CG	1:B:207:PHE:N	2.89	0.41
1:A:201:ASN:CG	3:A:761[B]:HDE:HAC	2.40	0.41
1:B:214:PHE:CD2	2:B:760[A]:HDD:CMC	3.04	0.40
1:C:377:ARG:NH1	5:C:1076:HOH:O	2.54	0.40
1:D:377:ARG:NH1	5:D:1500:HOH:O	2.53	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
5:A:3348:HOH:O	5:B:2822:HOH:O[2_545]	2.13	0.07
1:C:584:LYS:NZ	5:A:1624:HOH:O[1_554]	2.19	0.01

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

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	729/753 (97%)	707 (97%)	21 (3%)	1 (0%)	59	41
1	B	729/753 (97%)	708 (97%)	19 (3%)	2 (0%)	50	31
1	C	726/753 (96%)	700 (96%)	25 (3%)	1 (0%)	59	41
1	D	733/753 (97%)	710 (97%)	22 (3%)	1 (0%)	59	41
All	All	2917/3012 (97%)	2825 (97%)	87 (3%)	5 (0%)	56	38

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	725	ASP
1	A	75	SER
1	C	75	SER
1	B	75	SER
1	D	75	SER

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	614/634 (97%)	595 (97%)	19 (3%)	52	33
1	B	614/634 (97%)	590 (96%)	24 (4%)	43	23
1	C	611/634 (96%)	585 (96%)	26 (4%)	40	19
1	D	618/634 (98%)	590 (96%)	28 (4%)	38	17
All	All	2457/2536 (97%)	2360 (96%)	97 (4%)	43	23

All (97) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	SER
1	A	29	LEU
1	A	32	GLU
1	A	42	PRO
1	A	185	PHE
1	A	205	ILE
1	A	227	TRP

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Mol	Chain	Res	Type
1	A	252	ASN
1	A	274	CYS
1	A	369	ARG
1	A	377	ARG
1	A	398	PRO
1	A	432	PRO
1	A	440	TYR
1	A	459	ASN
1	A	552	LEU
1	A	713	GLN
1	A	732	LEU
1	A	751	ILE
1	B	157	ASN
1	B	158	LYS
1	B	185	PHE
1	B	205	ILE
1	B	227	TRP
1	B	252	ASN
1	B	274	CYS
1	B	432	PRO
1	B	440	TYR
1	B	459	ASN
1	B	552	LEU
1	B	562	LEU
1	B	565	GLU
1	B	568	ASP
1	B	571	LEU
1	B	583	LYS
1	B	616	LEU
1	B	633	LEU
1	B	703	LYS
1	B	713	GLN
1	B	725	ASP
1	B	732	LEU
1	B	750	LYS
1	B	751	ILE
1	C	32	GLU
1	C	61	ARG
1	C	159	ILE
1	C	185	PHE
1	C	205	ILE
1	C	227	TRP

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Mol	Chain	Res	Type
1	C	237	ASP
1	C	252	ASN
1	C	285	LYS
1	C	377	ARG
1	C	440	TYR
1	C	459	ASN
1	C	521	ARG
1	C	552	LEU
1	C	571	LEU
1	C	606	LEU
1	C	616	LEU
1	C	617	LEU
1	C	626	LYS
1	C	633	LEU
1	C	636	ARG
1	C	648	LEU
1	C	725	ASP
1	C	733	LEU
1	C	750	LYS
1	C	751	ILE
1	D	32	GLU
1	D	38	PRO
1	D	48	GLN
1	D	73[A]	LYS
1	D	73[B]	LYS
1	D	185	PHE
1	D	191	THR
1	D	205	ILE
1	D	227	TRP
1	D	237	ASP
1	D	252	ASN
1	D	274	CYS
1	D	369	ARG
1	D	416[A]	THR
1	D	416[B]	THR
1	D	432	PRO
1	D	440	TYR
1	D	459	ASN
1	D	490	GLU
1	D	552	LEU
1	D	574	THR
1	D	582	LEU

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Mol	Chain	Res	Type
1	D	616	LEU
1	D	648	LEU
1	D	700	ASP
1	D	713	GLN
1	D	747	LYS
1	D	750	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 12 ligands modelled in this entry, 4 are modelled with single atom - leaving 8 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	HDD	A	760[A]	1,5	52,52,52	2.55	23 (44%)	70,89,89	2.68	24 (34%)
3	HDE	A	761[B]	1,5	52,52,52	3.01	22 (42%)	75,89,89	2.72	35 (46%)
2	HDD	B	760[A]	1,5	52,52,52	2.27	19 (36%)	70,89,89	1.96	18 (25%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	HDE	B	761[B]	1,5	52,52,52	3.13	22 (42%)	75,89,89	2.85	35 (46%)
2	HDD	C	760[A]	1,5	52,52,52	2.38	21 (40%)	70,89,89	2.33	30 (42%)
3	HDE	C	761[B]	1,5	52,52,52	2.93	22 (42%)	75,89,89	2.84	35 (46%)
2	HDD	D	760[A]	1,5	52,52,52	2.64	17 (32%)	70,89,89	2.34	27 (38%)
3	HDE	D	761[B]	1,5	52,52,52	2.83	20 (38%)	75,89,89	2.67	35 (46%)

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	HDD	A	760[A]	1,5	-	0/5/89/89	0/1/9/9
3	HDE	A	761[B]	1,5	-	1/9/89/89	0/1/9/9
2	HDD	B	760[A]	1,5	-	0/5/89/89	0/1/9/9
3	HDE	B	761[B]	1,5	-	1/9/89/89	0/1/9/9
2	HDD	C	760[A]	1,5	-	0/5/89/89	0/1/9/9
3	HDE	C	761[B]	1,5	-	2/9/89/89	0/1/9/9
2	HDD	D	760[A]	1,5	-	0/5/89/89	0/1/9/9
3	HDE	D	761[B]	1,5	-	1/9/89/89	0/1/9/9

All (166) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	761[B]	HDE	C4C-C3C	8.79	1.49	1.37
3	B	761[B]	HDE	CHC-C4B	-8.05	1.44	1.51
3	A	761[B]	HDE	C4C-C3C	7.92	1.48	1.37
3	B	761[B]	HDE	C1C-NC	7.57	1.44	1.35
3	D	761[B]	HDE	CHC-C4B	-7.54	1.44	1.51
3	C	761[B]	HDE	C4C-C3C	7.47	1.47	1.37
3	A	761[B]	HDE	CHC-C4B	-7.34	1.44	1.51
3	A	761[B]	HDE	CHC-C1C	-7.09	1.44	1.51
2	D	760[A]	HDD	C3C-CAC	6.79	1.55	1.49
3	C	761[B]	HDE	CHD-C4C	-6.76	1.45	1.51
3	D	761[B]	HDE	C4C-C3C	6.68	1.46	1.37
3	D	761[B]	HDE	CHD-C1D	-6.67	1.45	1.51
2	D	760[A]	HDD	FE-ND	6.65	2.23	1.97
3	B	761[B]	HDE	CHD-C1D	-6.44	1.45	1.51
3	C	761[B]	HDE	CHC-C4B	-6.43	1.45	1.51
2	C	760[A]	HDD	FE-ND	6.29	2.21	1.97
3	C	761[B]	HDE	C1C-NC	6.22	1.43	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	761[B]	HDE	CHD-C4C	-6.22	1.45	1.51
3	C	761[B]	HDE	C4C-NC	6.18	1.43	1.35
3	A	761[B]	HDE	C4C-NC	6.01	1.42	1.35
3	A	761[B]	HDE	C1C-NC	6.00	1.42	1.35
3	D	761[B]	HDE	CHC-C1C	-5.95	1.46	1.51
2	A	760[A]	HDD	O1D-C3D	-5.90	1.36	1.46
2	B	760[A]	HDD	C3C-CAC	5.82	1.54	1.49
2	A	760[A]	HDD	C1C-C2C	5.77	1.47	1.40
2	D	760[A]	HDD	O1D-C3D	-5.15	1.38	1.46
3	B	761[B]	HDE	O1A-CGA	5.11	1.44	1.35
2	A	760[A]	HDD	C2D-C1D	-5.10	1.43	1.52
2	C	760[A]	HDD	C3C-C4C	5.10	1.48	1.40
2	A	760[A]	HDD	FE-ND	5.07	2.17	1.97
3	A	761[B]	HDE	O1A-CGA	5.02	1.44	1.35
3	D	761[B]	HDE	C4C-NC	5.00	1.41	1.35
2	D	760[A]	HDD	C4A-C3A	5.00	1.46	1.40
3	B	761[B]	HDE	CHD-C4C	-4.89	1.46	1.51
3	B	761[B]	HDE	C4C-NC	4.89	1.41	1.35
3	C	761[B]	HDE	CHD-C1D	-4.89	1.46	1.51
3	C	761[B]	HDE	C4B-NB	4.85	1.41	1.35
2	A	760[A]	HDD	C3C-CAC	4.84	1.53	1.49
2	B	760[A]	HDD	C4A-C3A	4.80	1.46	1.40
3	B	761[B]	HDE	C4B-NB	4.80	1.41	1.35
3	C	761[B]	HDE	O1A-CGA	4.77	1.44	1.35
2	A	760[A]	HDD	C3B-C2B	-4.72	1.33	1.41
3	C	761[B]	HDE	CHC-C1C	-4.72	1.47	1.51
2	D	760[A]	HDD	C1C-C2C	4.64	1.46	1.40
2	B	760[A]	HDD	C4B-C3B	4.60	1.47	1.41
2	C	760[A]	HDD	C3B-C2B	-4.48	1.33	1.41
2	A	760[A]	HDD	C4B-C3B	4.47	1.47	1.41
2	D	760[A]	HDD	C3B-C2B	-4.41	1.33	1.41
2	A	760[A]	HDD	C3C-C2C	-4.39	1.33	1.41
2	C	760[A]	HDD	C4B-C3B	4.33	1.47	1.41
3	B	761[B]	HDE	FE-ND	4.27	2.10	1.92
2	B	760[A]	HDD	C3C-C2C	-4.26	1.33	1.41
2	D	760[A]	HDD	FE-NB	4.26	2.10	1.92
3	D	761[B]	HDE	C1C-NC	4.26	1.40	1.35
3	B	761[B]	HDE	C4A-NA	4.22	1.44	1.38
3	A	761[B]	HDE	CHD-C1D	-4.20	1.47	1.51
2	B	760[A]	HDD	C2D-C1D	-4.18	1.44	1.52
3	C	761[B]	HDE	FE-ND	4.15	2.10	1.92
3	A	761[B]	HDE	CHD-C4C	-4.14	1.47	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	761[B]	HDE	FE-ND	4.14	2.10	1.92
2	D	760[A]	HDD	C3C-C4C	4.13	1.46	1.40
3	A	761[B]	HDE	C1A-NA	4.10	1.44	1.38
2	C	760[A]	HDD	C4A-C3A	4.09	1.45	1.40
3	B	761[B]	HDE	CHC-C1C	-4.08	1.47	1.51
3	D	761[B]	HDE	O1A-CGA	4.01	1.42	1.35
3	C	761[B]	HDE	C3B-C2B	4.00	1.49	1.37
2	D	760[A]	HDD	C2D-C1D	-3.99	1.45	1.52
3	A	761[B]	HDE	C4B-NB	3.98	1.40	1.35
3	D	761[B]	HDE	C3B-C2B	3.96	1.49	1.37
2	D	760[A]	HDD	C3B-CAB	3.91	1.55	1.48
3	A	761[B]	HDE	C4D-C3D	3.91	1.50	1.43
2	A	760[A]	HDD	C3D-C4D	-3.88	1.45	1.51
2	C	760[A]	HDD	O1D-C3D	-3.86	1.40	1.46
3	B	761[B]	HDE	C1A-NA	3.82	1.44	1.38
2	C	760[A]	HDD	C3C-CAC	3.81	1.52	1.49
3	A	761[B]	HDE	C3B-C2B	3.80	1.48	1.37
3	A	761[B]	HDE	C3C-C2C	3.79	1.48	1.37
2	B	760[A]	HDD	C3C-C4C	3.74	1.46	1.40
3	C	761[B]	HDE	C3D-C2D	3.73	1.48	1.37
2	A	760[A]	HDD	C3C-C4C	3.72	1.46	1.40
2	D	760[A]	HDD	C4B-C3B	3.70	1.46	1.41
3	A	761[B]	HDE	FE-ND	3.70	2.08	1.92
3	B	761[B]	HDE	C3C-C2C	3.68	1.48	1.37
3	A	761[B]	HDE	C3D-C2D	3.66	1.48	1.37
3	B	761[B]	HDE	C3B-C2B	3.65	1.48	1.37
2	D	760[A]	HDD	C3C-C2C	-3.59	1.35	1.41
2	C	760[A]	HDD	C1C-C2C	3.57	1.44	1.40
3	D	761[B]	HDE	C4B-NB	3.51	1.39	1.35
2	D	760[A]	HDD	FE-NC	3.46	2.07	1.92
2	B	760[A]	HDD	O1D-C3D	-3.46	1.40	1.46
2	D	760[A]	HDD	C3D-C4D	-3.46	1.45	1.51
3	D	761[B]	HDE	C1A-NA	3.46	1.43	1.38
2	A	760[A]	HDD	FE-NC	3.44	2.07	1.92
2	B	760[A]	HDD	C3B-C2B	-3.42	1.35	1.41
2	D	760[A]	HDD	CHD-C1D	3.41	1.42	1.36
3	C	761[B]	HDE	C1A-NA	3.41	1.43	1.38
2	C	760[A]	HDD	C3C-C2C	-3.38	1.35	1.41
2	C	760[A]	HDD	C3D-C4D	-3.37	1.46	1.51
3	D	761[B]	HDE	C4D-CHA	3.33	1.49	1.39
2	C	760[A]	HDD	FE-NC	3.32	2.06	1.92
2	A	760[A]	HDD	C1A-C2A	3.28	1.49	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	761[B]	HDE	C3C-C2C	3.27	1.47	1.37
2	B	760[A]	HDD	C1D-ND	3.25	1.43	1.38
3	C	761[B]	HDE	C4D-C3D	3.22	1.49	1.43
3	C	761[B]	HDE	C3C-C2C	3.21	1.47	1.37
2	C	760[A]	HDD	C3B-CAB	3.20	1.54	1.48
2	B	760[A]	HDD	C3B-CAB	3.19	1.54	1.48
2	C	760[A]	HDD	FE-NB	3.18	2.06	1.92
3	A	761[B]	HDE	C1B-CHB	3.13	1.48	1.39
3	B	761[B]	HDE	C4D-CHA	3.12	1.48	1.39
2	B	760[A]	HDD	FE-NC	3.09	2.05	1.92
2	C	760[A]	HDD	CHD-C1D	3.08	1.41	1.36
3	C	761[B]	HDE	C4D-CHA	3.00	1.48	1.39
2	B	760[A]	HDD	CMD-C2D	3.00	1.57	1.53
2	A	760[A]	HDD	FE-NB	2.94	2.05	1.92
3	D	761[B]	HDE	C3D-C2D	2.92	1.46	1.37
3	B	761[B]	HDE	C1C-C2C	2.91	1.43	1.38
2	C	760[A]	HDD	CHA-C4D	2.88	1.41	1.36
3	A	761[B]	HDE	C4D-CHA	2.82	1.47	1.39
3	A	761[B]	HDE	C4A-NA	2.81	1.42	1.38
3	B	761[B]	HDE	C3D-C2D	2.79	1.45	1.37
3	B	761[B]	HDE	C1B-NB	2.77	1.42	1.36
3	D	761[B]	HDE	C4D-C3D	2.77	1.48	1.43
3	D	761[B]	HDE	C4A-NA	2.76	1.42	1.38
2	A	760[A]	HDD	C3B-CAB	2.66	1.53	1.48
2	A	760[A]	HDD	C4A-NA	2.65	1.41	1.36
3	B	761[B]	HDE	FE-NC	2.65	2.03	1.92
2	D	760[A]	HDD	CMD-C2D	2.61	1.57	1.53
2	C	760[A]	HDD	CMD-C2D	2.58	1.57	1.53
3	A	761[B]	HDE	C1B-NB	2.57	1.41	1.36
2	C	760[A]	HDD	CMC-C2C	2.55	1.57	1.51
3	B	761[B]	HDE	CHB-C4A	-2.50	1.32	1.36
3	A	761[B]	HDE	C1D-ND	2.49	1.38	1.35
2	A	760[A]	HDD	OND-C2D	2.44	1.47	1.42
3	D	761[B]	HDE	C1B-CHB	2.43	1.46	1.39
3	A	761[B]	HDE	FE-NC	2.41	2.02	1.92
2	B	760[A]	HDD	FE-NB	2.40	2.02	1.92
2	B	760[A]	HDD	C1A-NA	2.40	1.41	1.36
3	C	761[B]	HDE	C4A-NA	2.39	1.42	1.38
2	B	760[A]	HDD	C1C-C2C	2.39	1.43	1.40
2	A	760[A]	HDD	C1B-NB	2.38	1.40	1.37
3	B	761[B]	HDE	C4D-C3D	2.36	1.47	1.43
2	B	760[A]	HDD	C3D-C4D	-2.34	1.47	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	760[A]	HDD	FE-ND	2.32	2.06	1.97
2	A	760[A]	HDD	CMC-C2C	2.29	1.56	1.51
2	B	760[A]	HDD	CMA-C3A	2.28	1.56	1.51
2	D	760[A]	HDD	C1A-NA	2.26	1.41	1.36
3	D	761[B]	HDE	C1B-NB	2.26	1.41	1.36
2	C	760[A]	HDD	C1B-NB	2.26	1.40	1.37
2	A	760[A]	HDD	CAA-C2A	2.24	1.56	1.52
2	C	760[A]	HDD	C1A-C2A	2.23	1.47	1.43
3	C	761[B]	HDE	FE-NC	2.23	2.01	1.92
3	C	761[B]	HDE	FE-NB	2.22	2.01	1.92
3	C	761[B]	HDE	C1C-C2C	2.20	1.42	1.38
2	A	760[A]	HDD	C4B-NB	2.14	1.40	1.37
3	B	761[B]	HDE	C1B-CHB	2.13	1.45	1.39
2	A	760[A]	HDD	C1A-NA	2.11	1.40	1.36
3	A	761[B]	HDE	C1C-C2C	2.08	1.41	1.38
2	A	760[A]	HDD	C4A-C3A	2.08	1.42	1.40
3	C	761[B]	HDE	C1B-NB	2.05	1.40	1.36
3	D	761[B]	HDE	O1A-C2A	-2.05	1.43	1.46
2	A	760[A]	HDD	CHD-C1D	2.04	1.39	1.36
2	C	760[A]	HDD	C2D-C1D	-2.04	1.48	1.52
3	C	761[B]	HDE	C1B-CHB	2.03	1.45	1.39
2	B	760[A]	HDD	OND-C2D	2.02	1.47	1.42
2	C	760[A]	HDD	C1C-NC	2.00	1.40	1.36

All (239) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	761[B]	HDE	C4A-NA-C1A	-7.25	103.03	107.94
3	C	761[B]	HDE	C2D-C1D-ND	7.10	117.85	110.11
2	A	760[A]	HDD	O1D-CGD-CBD	-7.05	101.85	110.21
2	C	760[A]	HDD	C3C-CAC-CBC	-7.02	111.42	125.95
2	A	760[A]	HDD	CAD-C3D-C2D	7.02	123.31	116.32
3	D	761[B]	HDE	C2D-C1D-ND	6.87	117.60	110.11
3	B	761[B]	HDE	C2D-C1D-ND	6.80	117.52	110.11
3	B	761[B]	HDE	C2C-C1C-NC	6.50	117.20	110.11
2	A	760[A]	HDD	C3B-CAB-CBB	-6.47	112.55	125.95
3	A	761[B]	HDE	C2C-C1C-NC	6.37	117.06	110.11
2	D	760[A]	HDD	C3B-CAB-CBB	-6.28	112.94	125.95
2	B	760[A]	HDD	O1D-CGD-O2D	6.17	126.89	120.81
2	A	760[A]	HDD	C1B-CHB-C4A	-6.08	119.47	127.47
3	D	761[B]	HDE	C2C-C1C-NC	5.98	116.63	110.11
3	B	761[B]	HDE	C4C-CHD-C1D	5.98	124.74	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	761[B]	HDE	C2D-C1D-ND	5.92	116.56	110.11
3	C	761[B]	HDE	C2C-C1C-NC	5.91	116.55	110.11
3	D	761[B]	HDE	C2B-C1B-NB	5.73	113.74	109.41
3	B	761[B]	HDE	C2B-C1B-NB	5.72	113.73	109.41
3	C	761[B]	HDE	C2B-C1B-NB	5.57	113.62	109.41
3	C	761[B]	HDE	C4A-NA-C1A	-5.42	104.27	107.94
2	C	760[A]	HDD	C3B-CAB-CBB	-5.39	114.80	125.95
2	D	760[A]	HDD	C4B-C3B-C2B	5.34	110.60	106.87
2	D	760[A]	HDD	C4D-ND-C1D	5.29	111.52	107.94
3	B	761[B]	HDE	C3A-C4A-NA	5.26	114.90	105.53
3	A	761[B]	HDE	C2A-C1A-NA	5.25	114.68	105.75
2	C	760[A]	HDD	C4D-ND-C1D	5.23	111.48	107.94
3	A	761[B]	HDE	C4A-NA-C1A	-5.23	104.40	107.94
2	A	760[A]	HDD	C4A-C3A-C2A	5.22	110.63	107.00
3	C	761[B]	HDE	CHD-C1D-C2D	-5.22	119.17	129.63
2	A	760[A]	HDD	CAD-CBD-CGD	-5.06	95.79	104.56
3	A	761[B]	HDE	CHD-C1D-C2D	-4.88	119.86	129.63
3	D	761[B]	HDE	CHD-C1D-C2D	-4.81	119.99	129.63
3	C	761[B]	HDE	C1B-CHB-C4A	-4.79	120.09	130.06
3	A	761[B]	HDE	C4D-C3D-C2D	-4.79	101.96	106.92
3	A	761[B]	HDE	C4B-CHC-C1C	4.78	122.32	112.60
3	D	761[B]	HDE	C3B-C4B-NB	4.75	115.09	110.07
3	D	761[B]	HDE	CAA-C2A-C3A	-4.73	111.62	116.32
2	A	760[A]	HDD	C4D-ND-C1D	4.71	111.12	107.94
2	B	760[A]	HDD	C3C-CAC-CBC	-4.69	116.24	125.95
3	C	761[B]	HDE	CBC-CAC-C3C	-4.67	96.24	111.99
3	C	761[B]	HDE	C3A-C4A-NA	4.65	113.81	105.53
3	B	761[B]	HDE	CBC-CAC-C3C	-4.59	96.51	111.99
2	B	760[A]	HDD	CMC-C2C-C1C	-4.58	121.57	128.62
2	A	760[A]	HDD	C3C-CAC-CBC	-4.56	116.50	125.95
3	A	761[B]	HDE	CAA-C2A-C1A	-4.54	108.67	116.02
2	A	760[A]	HDD	C4B-C3B-C2B	4.53	110.03	106.87
3	C	761[B]	HDE	C2A-C1A-NA	4.50	113.40	105.75
3	B	761[B]	HDE	C1C-C2C-C3C	-4.46	100.08	106.07
3	C	761[B]	HDE	C4C-CHD-C1D	4.46	121.67	112.60
3	D	761[B]	HDE	C4C-CHD-C1D	4.46	121.67	112.60
3	C	761[B]	HDE	C4D-C3D-C2D	-4.43	102.34	106.92
3	A	761[B]	HDE	CHD-C4C-NC	4.42	127.49	120.95
3	C	761[B]	HDE	CMB-C2B-C3B	4.41	133.26	124.94
2	A	760[A]	HDD	O1D-C3D-CAD	-4.41	102.30	105.29
2	D	760[A]	HDD	O1D-CGD-CBD	-4.35	105.06	110.21
3	B	761[B]	HDE	C2A-C1A-NA	4.31	113.07	105.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	760[A]	HDD	CMA-C3A-C4A	-4.30	122.01	128.62
3	A	761[B]	HDE	C4C-NC-C1C	-4.28	103.11	107.92
3	A	761[B]	HDE	C3A-C4A-NA	4.26	113.12	105.53
3	D	761[B]	HDE	C3A-C4A-NA	4.24	113.08	105.53
3	B	761[B]	HDE	CHD-C1D-C2D	-4.23	121.15	129.63
3	C	761[B]	HDE	C4C-NC-C1C	-4.21	103.19	107.92
2	D	760[A]	HDD	C3A-C4A-NA	-4.20	106.24	109.41
3	A	761[B]	HDE	O1A-CGA-O2A	4.18	124.93	120.81
2	C	760[A]	HDD	OND-C2D-CMD	-4.18	102.32	109.57
3	D	761[B]	HDE	C1B-CHB-C4A	-4.16	121.41	130.06
3	D	761[B]	HDE	C1B-C2B-C3B	-4.13	104.12	107.00
3	B	761[B]	HDE	CAD-CBD-CGD	-4.13	100.19	113.47
3	C	761[B]	HDE	CMB-C2B-C1B	-4.10	122.31	128.62
3	A	761[B]	HDE	CAC-C3C-C4C	-4.09	122.88	127.14
3	B	761[B]	HDE	C3B-C4B-NB	4.09	114.39	110.07
2	C	760[A]	HDD	C4B-C3B-C2B	4.07	109.71	106.87
3	C	761[B]	HDE	CHD-C4C-NC	4.06	126.97	120.95
3	D	761[B]	HDE	CHD-C4C-NC	4.04	126.93	120.95
2	C	760[A]	HDD	O1D-CGD-O2D	4.04	124.79	120.81
3	A	761[B]	HDE	CAD-CBD-CGD	-3.97	100.69	113.47
3	C	761[B]	HDE	C4B-CHC-C1C	3.96	120.65	112.60
2	B	760[A]	HDD	C3B-CAB-CBB	-3.96	117.76	125.95
3	D	761[B]	HDE	C4D-C3D-C2D	-3.95	102.83	106.92
2	B	760[A]	HDD	O1D-CGD-CBD	-3.85	105.64	110.21
3	B	761[B]	HDE	C1B-CHB-C4A	-3.85	122.04	130.06
2	A	760[A]	HDD	CBD-CAD-C3D	-3.79	99.06	105.00
3	B	761[B]	HDE	C4D-C3D-C2D	-3.76	103.02	106.92
3	C	761[B]	HDE	C1B-C2B-C3B	-3.76	104.38	107.00
3	B	761[B]	HDE	C4C-NC-C1C	-3.75	103.71	107.92
2	D	760[A]	HDD	C3B-C4B-NB	-3.74	106.12	111.52
3	C	761[B]	HDE	CMD-C2D-C3D	3.73	131.97	124.94
3	D	761[B]	HDE	CAD-CBD-CGD	-3.72	101.52	113.47
3	B	761[B]	HDE	CMC-C2C-C1C	3.68	131.81	127.13
2	A	760[A]	HDD	C2C-C1C-NC	-3.67	106.64	109.41
3	C	761[B]	HDE	CHC-C1C-C2C	-3.67	122.28	129.63
2	D	760[A]	HDD	O1D-CGD-O2D	3.64	124.40	120.81
3	A	761[B]	HDE	CAB-C3B-C4B	-3.63	123.35	127.14
3	B	761[B]	HDE	C1B-C2B-C3B	-3.63	104.47	107.00
2	A	760[A]	HDD	C3B-C4B-NB	-3.63	106.28	111.52
3	B	761[B]	HDE	C1B-NB-C4B	-3.62	101.41	106.67
3	D	761[B]	HDE	C2A-C1A-NA	3.61	111.90	105.75
3	D	761[B]	HDE	C4A-NA-C1A	-3.59	105.51	107.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	761[B]	HDE	CHD-C4C-C3C	-3.54	122.54	129.63
2	D	760[A]	HDD	C3C-CAC-CBC	-3.53	118.64	125.95
3	C	761[B]	HDE	C3B-C4B-NB	3.51	113.78	110.07
2	D	760[A]	HDD	C2C-C1C-NC	-3.51	106.76	109.41
2	D	760[A]	HDD	O1D-C3D-C2D	3.51	116.06	109.46
2	C	760[A]	HDD	CMC-C2C-C1C	-3.50	123.24	128.62
3	A	761[B]	HDE	C3B-C4B-NB	3.48	113.75	110.07
3	B	761[B]	HDE	C4B-CHC-C1C	3.48	119.67	112.60
3	A	761[B]	HDE	CBA-CAA-C2A	3.48	110.44	105.00
2	C	760[A]	HDD	C3C-C4C-NC	-3.48	105.26	108.64
2	D	760[A]	HDD	C3C-C2C-C1C	3.45	109.02	107.00
2	A	760[A]	HDD	CHB-C4A-NA	3.45	130.34	124.58
2	D	760[A]	HDD	CHA-C1A-NA	3.44	130.33	124.58
3	A	761[B]	HDE	C4C-CHD-C1D	3.44	119.59	112.60
2	D	760[A]	HDD	CAA-CBA-CGA	-3.43	102.45	113.47
2	C	760[A]	HDD	CBD-CAD-C3D	-3.40	99.66	105.00
2	B	760[A]	HDD	OND-C2D-CMD	-3.37	103.72	109.57
2	D	760[A]	HDD	O1D-C3D-CAD	-3.35	103.02	105.29
3	C	761[B]	HDE	CAA-C2A-C3A	-3.35	112.99	116.32
3	A	761[B]	HDE	C1C-C2C-C3C	-3.35	101.58	106.07
3	D	761[B]	HDE	C4B-CHC-C1C	3.34	119.38	112.60
3	C	761[B]	HDE	CAA-C2A-C1A	-3.31	110.66	116.02
3	D	761[B]	HDE	CHC-C4B-C3B	-3.30	123.02	129.63
3	A	761[B]	HDE	CHC-C4B-C3B	-3.29	123.03	129.63
2	A	760[A]	HDD	CMD-C2D-C1D	-3.29	107.53	113.22
3	D	761[B]	HDE	C4C-NC-C1C	-3.27	104.24	107.92
3	D	761[B]	HDE	C1B-NB-C4B	-3.26	101.93	106.67
2	D	760[A]	HDD	C2B-C3B-CAB	-3.25	120.67	127.33
2	C	760[A]	HDD	C3B-C4B-NB	-3.23	106.86	111.52
3	D	761[B]	HDE	CHC-C1C-C2C	-3.21	123.20	129.63
3	A	761[B]	HDE	C1B-CHB-C4A	-3.21	123.38	130.06
2	C	760[A]	HDD	O1D-CGD-CBD	-3.21	106.41	110.21
3	D	761[B]	HDE	CMA-C3A-C4A	-3.17	107.75	113.22
3	D	761[B]	HDE	CMB-C2B-C3B	3.16	130.90	124.94
3	C	761[B]	HDE	C1C-C2C-C3C	-3.15	101.84	106.07
2	B	760[A]	HDD	CAD-C3D-C4D	-3.13	110.94	116.02
3	B	761[B]	HDE	CHC-C4B-C3B	-3.13	123.36	129.63
3	B	761[B]	HDE	CHC-C1C-C2C	-3.11	123.39	129.63
2	B	760[A]	HDD	CAA-CBA-CGA	-3.11	103.48	113.47
3	D	761[B]	HDE	C4D-CHA-C1A	-3.09	123.63	130.06
3	A	761[B]	HDE	C2B-C1B-NB	3.08	111.73	109.41
2	C	760[A]	HDD	CMA-C3A-C4A	-3.08	123.89	128.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	761[B]	HDE	CMC-C2C-C3C	3.05	130.70	124.94
3	A	761[B]	HDE	CHC-C1C-C2C	-3.05	123.51	129.63
2	D	760[A]	HDD	C3D-C4D-ND	3.04	110.92	105.75
2	B	760[A]	HDD	C3D-C4D-ND	3.04	110.91	105.75
2	A	760[A]	HDD	CHA-C1A-NA	3.03	129.64	124.58
3	C	761[B]	HDE	CHC-C4B-C3B	-3.03	123.57	129.63
3	D	761[B]	HDE	CMD-C2D-C3D	3.02	130.64	124.94
2	D	760[A]	HDD	C4A-C3A-C2A	3.00	109.08	107.00
3	C	761[B]	HDE	C2B-C1B-CHB	-3.00	120.31	126.00
3	C	761[B]	HDE	C1B-NB-C4B	-2.99	102.33	106.67
3	D	761[B]	HDE	CMC-C2C-C3C	2.98	130.56	124.94
3	D	761[B]	HDE	CBC-CAC-C3C	-2.98	101.94	111.99
3	C	761[B]	HDE	CHD-C4C-C3C	-2.97	123.68	129.63
2	C	760[A]	HDD	C3D-C4D-ND	2.97	110.79	105.75
3	B	761[B]	HDE	C2B-C1B-CHB	-2.96	120.39	126.00
3	B	761[B]	HDE	CBA-CAA-C2A	2.96	109.63	105.00
3	D	761[B]	HDE	CAC-C3C-C4C	-2.95	124.07	127.14
2	C	760[A]	HDD	CMD-C2D-C1D	2.92	118.27	113.22
2	C	760[A]	HDD	C1B-CHB-C4A	-2.92	123.63	127.47
2	D	760[A]	HDD	CMC-C2C-C1C	-2.91	124.14	128.62
3	D	761[B]	HDE	C1C-C2C-C3C	-2.89	102.19	106.07
3	C	761[B]	HDE	C4D-CHA-C1A	-2.89	124.05	130.06
3	A	761[B]	HDE	C4D-CHA-C1A	-2.88	124.06	130.06
2	D	760[A]	HDD	OND-C2D-CMD	-2.88	104.57	109.57
3	B	761[B]	HDE	C4D-CHA-C1A	-2.86	124.10	130.06
3	C	761[B]	HDE	CMC-C2C-C3C	2.82	130.26	124.94
3	C	761[B]	HDE	CAD-CBD-CGD	-2.82	104.42	113.47
3	B	761[B]	HDE	C4D-ND-C1D	-2.79	102.61	106.67
3	B	761[B]	HDE	CHD-C4C-NC	2.79	125.08	120.95
3	A	761[B]	HDE	C3D-C4D-ND	2.78	113.60	109.73
3	B	761[B]	HDE	C4C-C3C-C2C	2.76	110.18	105.77
3	D	761[B]	HDE	CHD-C4C-C3C	-2.75	124.11	129.63
3	A	761[B]	HDE	O1A-C2A-CAA	-2.73	103.44	105.29
3	B	761[B]	HDE	CAA-C2A-C3A	-2.71	113.63	116.32
2	D	760[A]	HDD	CMA-C3A-C4A	-2.70	124.47	128.62
2	B	760[A]	HDD	C1A-CHA-C4D	-2.69	124.45	130.06
3	C	761[B]	HDE	CMA-C3A-C4A	-2.68	108.59	113.22
3	C	761[B]	HDE	C3D-C4D-ND	2.67	113.45	109.73
2	A	760[A]	HDD	C3D-C4D-ND	2.67	110.29	105.75
2	C	760[A]	HDD	O1D-C3D-CAD	-2.63	103.51	105.29
3	D	761[B]	HDE	CBA-CAA-C2A	2.62	109.10	105.00
3	A	761[B]	HDE	C1B-C2B-C3B	-2.61	105.18	107.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	760[A]	HDD	CMC-C2C-C3C	2.60	129.07	124.97
2	B	760[A]	HDD	C4B-CHC-C1C	-2.60	124.05	127.47
3	B	761[B]	HDE	CMA-C3A-C4A	-2.60	108.74	113.22
2	C	760[A]	HDD	C4C-NC-C1C	2.58	110.16	106.76
3	B	761[B]	HDE	CAA-C2A-C1A	-2.57	111.85	116.02
3	B	761[B]	HDE	O1A-CGA-O2A	2.56	123.33	120.81
3	B	761[B]	HDE	CAC-C3C-C2C	-2.52	121.65	127.81
2	C	760[A]	HDD	C4C-C3C-CAC	-2.52	122.07	127.18
2	C	760[A]	HDD	C2C-C1C-NC	-2.52	107.51	109.41
3	D	761[B]	HDE	C4D-ND-C1D	-2.51	103.02	106.67
2	D	760[A]	HDD	CBD-CAD-C3D	-2.49	101.10	105.00
2	B	760[A]	HDD	CBD-CAD-C3D	-2.49	101.10	105.00
3	B	761[B]	HDE	C3D-C4D-ND	2.48	113.18	109.73
3	B	761[B]	HDE	CMB-C2B-C3B	2.48	129.62	124.94
2	A	760[A]	HDD	C3C-C2C-C1C	2.47	108.45	107.00
3	D	761[B]	HDE	C3D-C4D-ND	2.47	113.16	109.73
2	B	760[A]	HDD	C2D-C1D-ND	2.46	109.91	105.53
3	B	761[B]	HDE	ONA-C3A-C4A	2.45	116.22	111.72
2	D	760[A]	HDD	CHB-C4A-NA	2.45	128.67	124.58
3	C	761[B]	HDE	O1A-CGA-O2A	2.45	123.22	120.81
3	C	761[B]	HDE	C4D-ND-C1D	-2.42	103.14	106.67
2	A	760[A]	HDD	O1D-C3D-C2D	-2.41	104.92	109.46
3	D	761[B]	HDE	CMB-C2B-C1B	-2.39	124.94	128.62
3	A	761[B]	HDE	C4D-ND-C1D	-2.36	103.24	106.67
3	D	761[B]	HDE	O1A-CGA-O2A	2.35	123.13	120.81
2	C	760[A]	HDD	C2D-C1D-CHD	2.34	129.21	123.22
2	B	760[A]	HDD	C1A-C2A-C3A	2.33	109.33	106.92
3	A	761[B]	HDE	C3D-C4D-CHA	-2.30	121.64	126.00
3	A	761[B]	HDE	CMD-C2D-C3D	2.30	129.27	124.94
2	D	760[A]	HDD	CMB-C2B-C1B	2.29	130.68	126.16
2	C	760[A]	HDD	C4A-C3A-C2A	2.27	108.58	107.00
2	D	760[A]	HDD	C1A-CHA-C4D	-2.25	125.37	130.06
2	D	760[A]	HDD	C4B-NB-C1B	2.24	108.94	105.58
2	A	760[A]	HDD	C3A-C4A-CHB	-2.23	121.78	126.00
2	D	760[A]	HDD	C2A-C1A-NA	-2.22	106.65	109.73
3	A	761[B]	HDE	CMB-C2B-C3B	2.22	129.13	124.94
2	A	760[A]	HDD	CAA-CBA-CGA	-2.20	106.39	113.47
2	B	760[A]	HDD	CMC-C2C-C3C	2.20	128.43	124.97
2	C	760[A]	HDD	O1A-CGA-CBA	-2.18	115.51	123.03
2	B	760[A]	HDD	C3C-C4C-NC	-2.18	106.53	108.64
3	C	761[B]	HDE	C1D-C2D-C3D	-2.17	103.16	106.07
2	B	760[A]	HDD	C3B-C4B-NB	-2.16	108.40	111.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	760[A]	HDD	C3D-O1D-CGD	2.15	116.67	109.33
3	A	761[B]	HDE	C1B-NB-C4B	-2.14	103.56	106.67
2	C	760[A]	HDD	CAD-CBD-CGD	-2.14	100.85	104.56
2	C	760[A]	HDD	C2A-C1A-NA	-2.12	106.78	109.73
2	B	760[A]	HDD	CMB-C2B-C3B	2.12	128.30	124.97
2	A	760[A]	HDD	C2B-C3B-CAB	-2.11	123.01	127.33
3	A	761[B]	HDE	O1A-C2A-C3A	2.09	113.39	109.46
2	C	760[A]	HDD	CHA-C1A-NA	2.08	128.05	124.58
2	C	760[A]	HDD	CHB-C4A-NA	2.04	127.98	124.58
2	C	760[A]	HDD	C3A-C4A-NA	-2.03	107.88	109.41
2	C	760[A]	HDD	CHD-C4C-NC	2.03	127.97	124.58
2	C	760[A]	HDD	C4C-C3C-C2C	2.03	109.43	106.97
2	A	760[A]	HDD	C3A-C4A-NA	-2.02	107.88	109.41
3	D	761[B]	HDE	C1D-C2D-C3D	-2.02	103.36	106.07

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	761[B]	HDE	C2B-C3B-CAB-CBB
3	D	761[B]	HDE	C2B-C3B-CAB-CBB
3	B	761[B]	HDE	C2B-C3B-CAB-CBB
3	C	761[B]	HDE	C2C-C3C-CAC-CBC
3	A	761[B]	HDE	C2B-C3B-CAB-CBB

There are no ring outliers.

5.7 Other polymers ⓘ

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	726/753 (96%)	-0.65	3 (0%) 90 89	3, 10, 26, 44	2 (0%)
1	B	726/753 (96%)	-0.50	10 (1%) 72 68	3, 11, 32, 47	2 (0%)
1	C	726/753 (96%)	-0.53	2 (0%) 91 90	3, 11, 31, 46	2 (0%)
1	D	726/753 (96%)	-0.65	3 (0%) 90 89	3, 10, 27, 44	1 (0%)
All	All	2904/3012 (96%)	-0.58	18 (0%) 84 85	3, 11, 30, 47	7 (0%)

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	28	SER	3.8
1	D	28	SER	3.5
1	B	32	GLU	3.1
1	A	32	GLU	3.0
1	C	594	PRO	2.9
1	B	726	GLY	2.8
1	B	713	GLN	2.8
1	B	28	SER	2.6
1	B	673	ALA	2.5
1	C	750	LYS	2.5
1	D	713	GLN	2.3
1	B	712	ASP	2.3
1	B	552	LEU	2.3
1	D	749	ASP	2.3
1	A	713	GLN	2.1
1	B	568	ASP	2.1
1	B	725	ASP	2.0
1	B	583	LYS	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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, 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	H2S	D	754	1/1	0.24	22.81	23,23,23,23	1
4	H2S	C	754	1/1	0.26	12.35	20,20,20,20	1
4	H2S	A	754	1/1	0.30	10.90	28,28,28,28	1
4	H2S	B	754	1/1	0.20	7.31	26,26,26,26	1
2	HDD	C	760[A]	44/44	0.08	1.17	2,4,9,14	44
3	HDE	C	761[B]	44/44	0.07	0.97	3,11,13,15	44
2	HDD	D	760[A]	44/44	0.07	0.92	2,3,7,12	44
2	HDD	A	760[A]	44/44	0.08	0.69	2,3,6,9	44
3	HDE	B	761[B]	44/44	0.07	0.62	2,9,13,14	44
2	HDD	B	760[A]	44/44	0.07	0.48	2,3,6,7	44
3	HDE	A	761[B]	44/44	0.08	0.43	2,7,10,13	44
3	HDE	D	761[B]	44/44	0.06	0.26	2,8,12,16	44

6.5 Other polymers ⓘ

There are no such residues in this entry.