



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 06:01 PM UTC

PDB ID : 2RJH / pdb_00002rjh
Title : Crystal structure of biosynthetic alaine racemase in D-cycloserine-bound form from Escherichia coli
Authors : Wu, D.; Hu, T.; Zhang, L.; Jiang, H.; Shen, X.
Deposited on : 2007-10-15
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
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
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

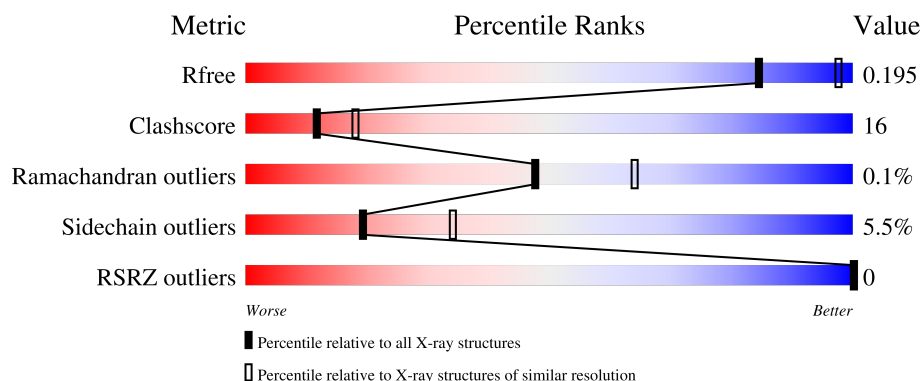
1 Overall quality at a glance

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	379	
1	B	379	
1	C	379	
1	D	379	

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	DCS	A	362	X	-	-	-
3	DCS	B	362	X	X	-	-
3	DCS	C	362	X	-	-	-
3	DCS	D	360	X	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11633 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 Alanine racemase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	359	Total	C	N	O	S	0	0	0
			2754	1734	498	507	15			
1	B	359	Total	C	N	O	S	0	0	0
			2754	1734	498	507	15			
1	C	353	Total	C	N	O	S	0	0	0
			2704	1704	488	498	14			
1	D	357	Total	C	N	O	S	0	0	0
			2737	1724	495	504	14			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP P0A6B4
A	-18	GLY	-	expression tag	UNP P0A6B4
A	-17	SER	-	expression tag	UNP P0A6B4
A	-16	SER	-	expression tag	UNP P0A6B4
A	-15	HIS	-	expression tag	UNP P0A6B4
A	-14	HIS	-	expression tag	UNP P0A6B4
A	-13	HIS	-	expression tag	UNP P0A6B4
A	-12	HIS	-	expression tag	UNP P0A6B4
A	-11	HIS	-	expression tag	UNP P0A6B4
A	-10	HIS	-	expression tag	UNP P0A6B4
A	-9	SER	-	expression tag	UNP P0A6B4
A	-8	SER	-	expression tag	UNP P0A6B4
A	-7	GLY	-	expression tag	UNP P0A6B4
A	-6	LEU	-	expression tag	UNP P0A6B4
A	-5	VAL	-	expression tag	UNP P0A6B4
A	-4	PRO	-	expression tag	UNP P0A6B4
A	-3	ARG	-	expression tag	UNP P0A6B4
A	-2	GLY	-	expression tag	UNP P0A6B4
A	-1	SER	-	expression tag	UNP P0A6B4
A	0	HIS	-	expression tag	UNP P0A6B4
B	-19	MET	-	expression tag	UNP P0A6B4

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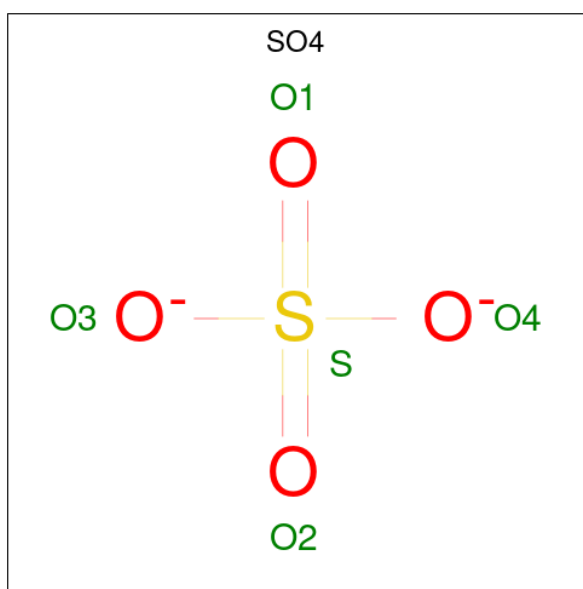
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B	-18	GLY	-	expression tag	UNP P0A6B4
B	-17	SER	-	expression tag	UNP P0A6B4
B	-16	SER	-	expression tag	UNP P0A6B4
B	-15	HIS	-	expression tag	UNP P0A6B4
B	-14	HIS	-	expression tag	UNP P0A6B4
B	-13	HIS	-	expression tag	UNP P0A6B4
B	-12	HIS	-	expression tag	UNP P0A6B4
B	-11	HIS	-	expression tag	UNP P0A6B4
B	-10	HIS	-	expression tag	UNP P0A6B4
B	-9	SER	-	expression tag	UNP P0A6B4
B	-8	SER	-	expression tag	UNP P0A6B4
B	-7	GLY	-	expression tag	UNP P0A6B4
B	-6	LEU	-	expression tag	UNP P0A6B4
B	-5	VAL	-	expression tag	UNP P0A6B4
B	-4	PRO	-	expression tag	UNP P0A6B4
B	-3	ARG	-	expression tag	UNP P0A6B4
B	-2	GLY	-	expression tag	UNP P0A6B4
B	-1	SER	-	expression tag	UNP P0A6B4
B	0	HIS	-	expression tag	UNP P0A6B4
C	-19	MET	-	expression tag	UNP P0A6B4
C	-18	GLY	-	expression tag	UNP P0A6B4
C	-17	SER	-	expression tag	UNP P0A6B4
C	-16	SER	-	expression tag	UNP P0A6B4
C	-15	HIS	-	expression tag	UNP P0A6B4
C	-14	HIS	-	expression tag	UNP P0A6B4
C	-13	HIS	-	expression tag	UNP P0A6B4
C	-12	HIS	-	expression tag	UNP P0A6B4
C	-11	HIS	-	expression tag	UNP P0A6B4
C	-10	HIS	-	expression tag	UNP P0A6B4
C	-9	SER	-	expression tag	UNP P0A6B4
C	-8	SER	-	expression tag	UNP P0A6B4
C	-7	GLY	-	expression tag	UNP P0A6B4
C	-6	LEU	-	expression tag	UNP P0A6B4
C	-5	VAL	-	expression tag	UNP P0A6B4
C	-4	PRO	-	expression tag	UNP P0A6B4
C	-3	ARG	-	expression tag	UNP P0A6B4
C	-2	GLY	-	expression tag	UNP P0A6B4
C	-1	SER	-	expression tag	UNP P0A6B4
C	0	HIS	-	expression tag	UNP P0A6B4
D	-19	MET	-	expression tag	UNP P0A6B4
D	-18	GLY	-	expression tag	UNP P0A6B4
D	-17	SER	-	expression tag	UNP P0A6B4

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-16	SER	-	expression tag	UNP P0A6B4
D	-15	HIS	-	expression tag	UNP P0A6B4
D	-14	HIS	-	expression tag	UNP P0A6B4
D	-13	HIS	-	expression tag	UNP P0A6B4
D	-12	HIS	-	expression tag	UNP P0A6B4
D	-11	HIS	-	expression tag	UNP P0A6B4
D	-10	HIS	-	expression tag	UNP P0A6B4
D	-9	SER	-	expression tag	UNP P0A6B4
D	-8	SER	-	expression tag	UNP P0A6B4
D	-7	GLY	-	expression tag	UNP P0A6B4
D	-6	LEU	-	expression tag	UNP P0A6B4
D	-5	VAL	-	expression tag	UNP P0A6B4
D	-4	PRO	-	expression tag	UNP P0A6B4
D	-3	ARG	-	expression tag	UNP P0A6B4
D	-2	GLY	-	expression tag	UNP P0A6B4
D	-1	SER	-	expression tag	UNP P0A6B4
D	0	HIS	-	expression tag	UNP P0A6B4

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



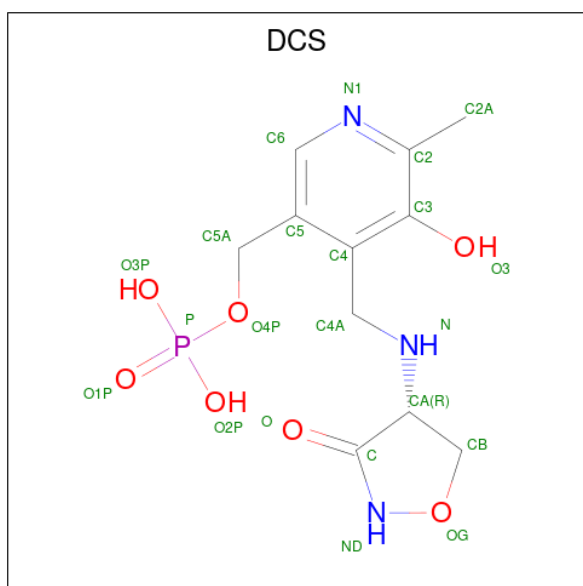
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

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

- Molecule 3 is D-[3-HYDROXY-2-METHYL-5-PHOSPHONOOXYMETHYL-PYRIDIN-4-YLMETHYL]-N,O-CYCLOSERYLAMIDE (CCD ID: DCS) (formula: C₁₁H₁₆N₃O₇P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			22	11	3	7	1		
3	B	1	Total	C	N	O	P	0	0
			22	11	3	7	1		
3	C	1	Total	C	N	O	P	0	0
			22	11	3	7	1		
3	D	1	Total	C	N	O	P	0	0
			22	11	3	7	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	185	Total	O	0	0
			185	185		
4	B	190	Total	O	0	0
			190	190		

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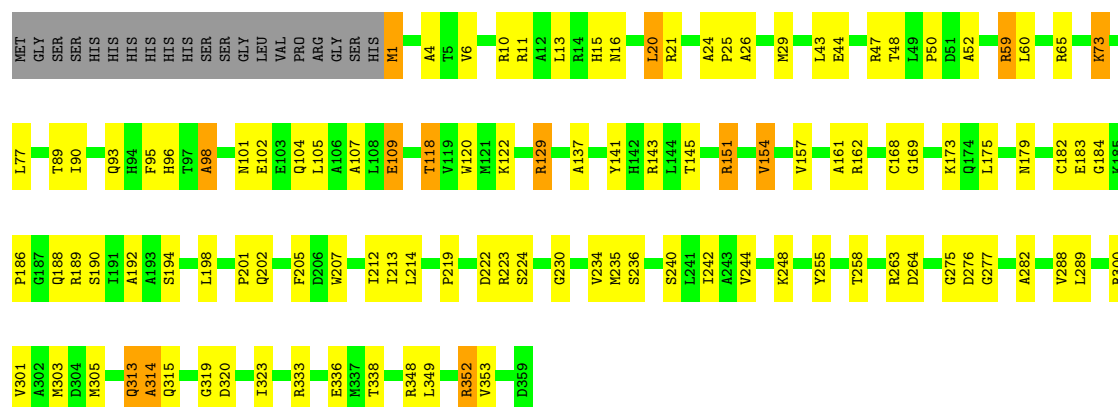
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	83	Total 83	O 83	0	0
4	D	108	Total 108	O 108	0	0

3 Residue-property plots

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.

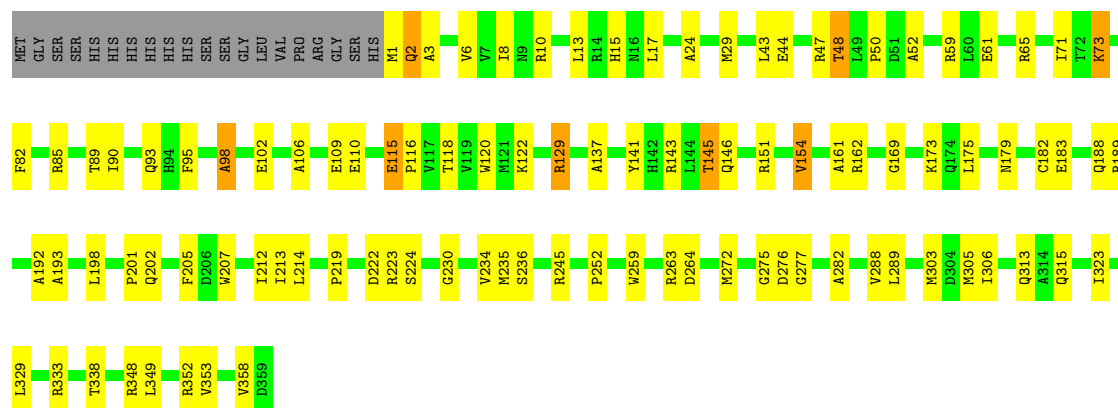
• Molecule 1: Alanine racemase

Chain A: 



• Molecule 1: Alanine racemase

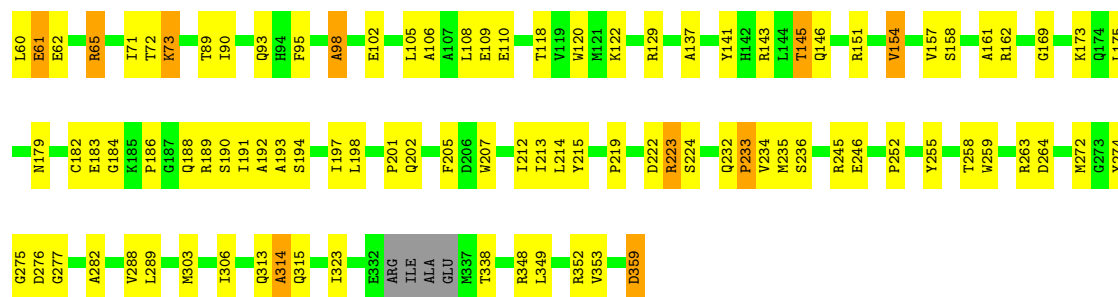
Chain B: 



• Molecule 1: Alanine racemase

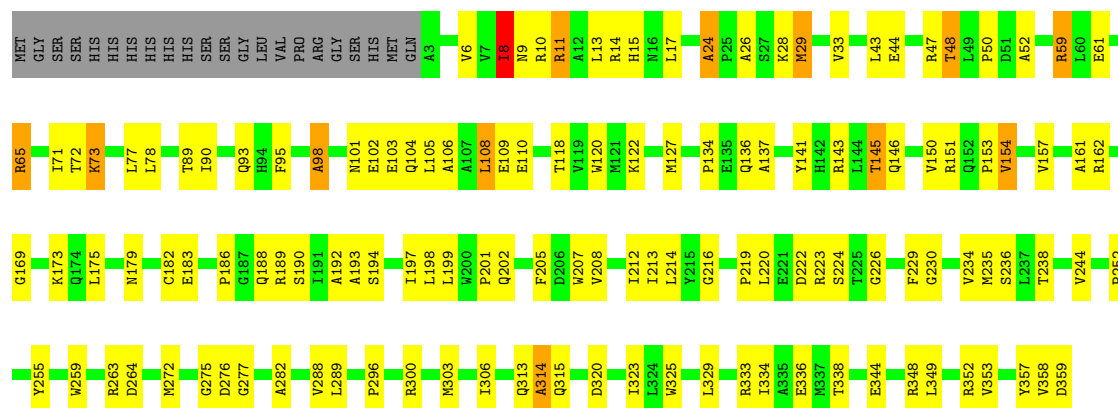
Chain C: 





• Molecule 1: Alanine racemase

Chain D: 59% 32% 6%



4 Data and refinement statistics

Property	Value	Source
Space group	P 6	Depositor
Cell constants a, b, c, α , β , γ	147.71Å 147.71Å 163.72Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	15.00 – 2.40 15.00 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.9 (15.00-2.40) 99.9 (15.00-2.40)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.71 (at 2.39Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.199 , 0.228 (Not available) , 0.195	Depositor DCC
R_{free} test set	7605 reflections (9.63%)	wwPDB-VP
Wilson B-factor (Å ²)	15.9	Xtriage
Anisotropy	0.403	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 10.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.487 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11633	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.86% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: DCS, SO4, KCX

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	0.81	9/2799 (0.3%)	1.21	28/3800 (0.7%)
1	B	0.75	4/2799 (0.1%)	1.14	24/3800 (0.6%)
1	C	0.60	2/2748 (0.1%)	1.15	24/3731 (0.6%)
1	D	0.59	0/2782	1.13	25/3778 (0.7%)
All	All	0.69	15/11128 (0.1%)	1.16	101/15109 (0.7%)

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	D	0	1

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	89	THR	C-O	-6.99	1.16	1.24
1	A	109	GLU	C-O	-6.48	1.16	1.24
1	A	352	ARG	C-O	-6.06	1.16	1.24
1	A	59	ARG	CZ-NH2	-5.88	1.25	1.33
1	C	233	PRO	C-O	-5.64	1.17	1.23
1	B	115	GLU	C-O	-5.62	1.18	1.24
1	A	60	LEU	C-O	-5.58	1.17	1.24
1	A	313	GLN	C-O	-5.42	1.15	1.23
1	A	48	THR	C-O	-5.42	1.17	1.24
1	A	59	ARG	C-O	-5.28	1.17	1.23
1	A	118	THR	CB-CG2	-5.26	1.35	1.52
1	A	107	ALA	CA-CB	-5.22	1.45	1.53
1	B	71	ILE	C-O	-5.16	1.18	1.24
1	B	61	GLU	C-O	-5.13	1.18	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	15	HIS	C-O	-5.08	1.18	1.24

All (101) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	333	ARG	NE-CZ-NH2	11.37	129.43	119.20
1	A	333	ARG	NE-CZ-NH1	-10.97	110.53	121.50
1	B	85	ARG	NE-CZ-NH2	10.80	128.92	119.20
1	C	151	ARG	NE-CZ-NH2	-10.51	109.74	119.20
1	C	151	ARG	CD-NE-CZ	10.40	138.96	124.40
1	C	223	ARG	NE-CZ-NH2	-10.38	109.86	119.20
1	B	85	ARG	NE-CZ-NH1	-10.37	111.13	121.50
1	D	8	ILE	N-CA-C	10.09	122.72	107.99
1	C	223	ARG	CD-NE-CZ	9.68	137.96	124.40
1	C	151	ARG	NE-CZ-NH1	9.31	130.81	121.50
1	C	223	ARG	NE-CZ-NH1	9.13	130.63	121.50
1	A	65	ARG	NE-CZ-NH2	9.12	127.41	119.20
1	D	65	ARG	NE-CZ-NH2	9.02	127.32	119.20
1	C	11	ARG	NE-CZ-NH2	8.91	127.22	119.20
1	A	65	ARG	NE-CZ-NH1	-8.86	112.64	121.50
1	C	11	ARG	NE-CZ-NH1	-8.52	112.98	121.50
1	D	65	ARG	NE-CZ-NH1	-8.48	113.02	121.50
1	A	11	ARG	NE-CZ-NH2	8.22	126.60	119.20
1	D	314	ALA	N-CA-C	8.08	124.70	111.37
1	A	333	ARG	CD-NE-CZ	8.01	135.62	124.40
1	A	129	ARG	NE-CZ-NH1	-7.94	113.56	121.50
1	A	11	ARG	NE-CZ-NH1	-7.85	113.65	121.50
1	C	314	ALA	N-CA-C	7.67	124.02	111.37
1	A	129	ARG	NE-CZ-NH2	7.65	126.08	119.20
1	B	129	ARG	NE-CZ-NH2	7.61	126.05	119.20
1	D	11	ARG	CD-NE-CZ	7.58	135.02	124.40
1	D	11	ARG	NE-CZ-NH2	-7.50	112.45	119.20
1	B	129	ARG	NE-CZ-NH1	-7.42	114.08	121.50
1	D	65	ARG	CD-NE-CZ	7.34	134.67	124.40
1	A	65	ARG	CD-NE-CZ	7.31	134.64	124.40
1	B	85	ARG	CD-NE-CZ	7.12	134.37	124.40
1	B	65	ARG	CD-NE-CZ	6.92	134.09	124.40
1	D	8	ILE	N-CA-CB	6.84	119.21	111.21
1	A	11	ARG	CD-NE-CZ	6.76	133.86	124.40
1	C	65	ARG	CD-NE-CZ	6.75	133.85	124.40
1	A	352	ARG	N-CA-C	6.74	120.77	112.54
1	C	11	ARG	CD-NE-CZ	6.69	133.76	124.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	315	GLN	N-CA-C	6.62	120.82	113.21
1	A	314	ALA	N-CA-C	6.52	124.70	110.80
1	C	65	ARG	NE-CZ-NH2	-6.49	113.36	119.20
1	A	98	ALA	N-CA-C	-6.42	101.88	110.55
1	D	11	ARG	NE-CZ-NH1	6.38	127.88	121.50
1	B	137	ALA	N-CA-C	6.36	118.21	111.28
1	A	198	LEU	N-CA-C	6.29	117.93	111.14
1	B	65	ARG	NE-CZ-NH2	-6.26	113.57	119.20
1	D	137	ALA	N-CA-C	6.10	117.93	111.28
1	A	137	ALA	N-CA-C	6.08	117.91	111.28
1	C	137	ALA	N-CA-C	6.00	117.82	111.28
1	B	98	ALA	N-CA-C	-5.97	102.48	110.55
1	C	198	LEU	N-CA-C	5.93	117.82	111.36
1	B	198	LEU	N-CA-C	5.89	117.50	111.14
1	C	129	ARG	NE-CZ-NH2	-5.88	113.91	119.20
1	D	198	LEU	N-CA-C	5.88	117.77	111.36
1	C	65	ARG	NE-CZ-NH1	5.87	127.37	121.50
1	C	98	ALA	N-CA-C	-5.86	102.64	110.55
1	A	275	GLY	N-CA-C	-5.81	108.28	114.67
1	A	264	ASP	N-CA-C	-5.80	101.70	110.28
1	D	98	ALA	N-CA-C	-5.79	102.73	110.55
1	C	24	ALA	CA-C-N	5.79	125.49	119.82
1	C	24	ALA	C-N-CA	5.79	125.49	119.82
1	B	65	ARG	NE-CZ-NH1	5.67	127.17	121.50
1	B	151	ARG	NE-CZ-NH2	5.63	124.27	119.20
1	C	275	GLY	N-CA-C	-5.60	108.52	114.67
1	D	29	MET	N-CA-C	5.57	118.03	109.07
1	C	29	MET	N-CA-C	5.55	118.01	109.07
1	B	358	VAL	N-CA-C	5.54	117.63	108.99
1	D	224	SER	N-CA-C	-5.54	103.07	110.55
1	C	192	ALA	N-CA-C	5.51	119.21	109.96
1	D	264	ASP	N-CA-C	-5.49	101.72	109.96
1	D	275	GLY	N-CA-C	-5.47	108.16	115.32
1	B	151	ARG	NE-CZ-NH1	-5.44	116.06	121.50
1	A	192	ALA	N-CA-C	5.43	119.08	109.96
1	B	275	GLY	N-CA-C	-5.42	108.71	114.67
1	B	192	ALA	N-CA-C	5.40	119.04	109.96
1	D	108	LEU	N-CA-CB	-5.38	103.45	110.88
1	D	244	VAL	N-CA-C	-5.38	99.81	107.77
1	D	151	ARG	NE-CZ-NH2	5.36	124.02	119.20
1	B	29	MET	N-CA-C	5.34	117.67	109.07
1	C	224	SER	N-CA-C	-5.30	103.40	110.55

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	329	LEU	CA-C-N	5.29	125.21	119.76
1	B	329	LEU	C-N-CA	5.29	125.21	119.76
1	D	24	ALA	CA-C-N	5.28	125.00	119.82
1	D	24	ALA	C-N-CA	5.28	125.00	119.82
1	C	264	ASP	N-CA-C	-5.27	102.05	109.96
1	D	192	ALA	N-CA-C	5.27	118.81	109.96
1	A	29	MET	N-CA-C	5.23	117.50	109.07
1	B	224	SER	N-CA-C	-5.23	103.48	110.55
1	B	264	ASP	N-CA-C	-5.22	102.14	109.96
1	D	151	ARG	NE-CZ-NH1	-5.20	116.30	121.50
1	A	224	SER	N-CA-C	-5.16	103.59	110.55
1	D	320	ASP	CA-C-N	5.14	124.90	119.76
1	D	320	ASP	C-N-CA	5.14	124.90	119.76
1	A	24	ALA	CA-C-N	5.14	124.80	119.56
1	A	24	ALA	C-N-CA	5.14	124.80	119.56
1	A	244	VAL	N-CA-C	-5.13	100.17	107.77
1	A	301	VAL	N-CA-C	-5.08	102.18	108.89
1	A	320	ASP	CA-C-N	5.07	124.83	119.76
1	A	320	ASP	C-N-CA	5.07	124.83	119.76
1	B	24	ALA	CA-C-N	5.04	124.76	119.82
1	B	24	ALA	C-N-CA	5.04	124.76	119.82
1	A	107	ALA	N-CA-C	5.00	117.11	111.11

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	358	VAL	Mainchain

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	2754	0	2760	81	0
1	B	2754	0	2760	68	0
1	C	2704	0	2704	94	0
1	D	2737	0	2740	115	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	10	0	0	0	0
2	B	10	0	0	0	0
2	C	10	0	0	0	0
3	A	22	0	13	2	0
3	B	22	0	13	6	0
3	C	22	0	13	5	0
3	D	22	0	13	7	0
4	A	185	0	0	11	0
4	B	190	0	0	3	0
4	C	83	0	0	11	0
4	D	108	0	0	26	0
All	All	11633	0	11016	352	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 16.

All (352) 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:B:362:DCS:P	3:B:362:DCS:O3P	1.15	1.54
1:C:6:VAL:HG23	1:C:353:VAL:HG11	1.42	1.01
3:B:362:DCS:O3P	3:B:362:DCS:O4P	1.78	1.00
1:C:52:ALA:O	1:C:73:LYS:HD3	1.61	1.00
1:C:222:ASP:OD2	4:C:498:HOH:O	1.79	0.99
1:A:1:MET:HE1	1:A:242:ILE:HG22	1.44	0.99
1:A:52:ALA:O	1:A:73:LYS:HD3	1.62	0.98
1:D:6:VAL:HG23	1:D:353:VAL:HG11	1.45	0.97
1:C:11:ARG:NH1	1:C:359:ASP:O	1.97	0.97
1:B:52:ALA:O	1:B:73:LYS:HD3	1.64	0.96
1:D:52:ALA:O	1:D:73:LYS:HD3	1.65	0.95
3:D:360:DCS:N	3:D:360:DCS:H5A1	1.82	0.94
1:D:134:PRO:HG2	4:D:449:HOH:O	1.69	0.93
4:C:437:HOH:O	1:D:348:ARG:HB2	1.69	0.91
1:D:29:MET:HE1	4:D:439:HOH:O	1.73	0.89
1:A:73:LYS:HZ2	1:A:73:LYS:HB2	1.36	0.89
1:D:44:GLU:O	1:D:48:THR:HG23	1.73	0.88
1:C:44:GLU:O	1:C:48:THR:HG23	1.73	0.88
1:B:44:GLU:O	1:B:48:THR:HG23	1.75	0.86
3:C:362:DCS:HND	1:D:303:MET:H	1.22	0.86
3:B:362:DCS:O3P	3:B:362:DCS:O2P	1.92	0.86
1:C:73:LYS:NZ	1:C:73:LYS:HB2	1.90	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:73:LYS:HB2	1:D:73:LYS:NZ	1.90	0.85
1:C:222:ASP:O	1:C:223:ARG:HG2	1.76	0.85
1:A:73:LYS:HB2	1:A:73:LYS:NZ	1.91	0.84
1:D:229:PHE:HZ	4:D:459:HOH:O	1.60	0.83
1:B:73:LYS:HB2	1:B:73:LYS:NZ	1.92	0.83
1:D:229:PHE:CZ	4:D:459:HOH:O	2.32	0.82
1:A:151:ARG:HG2	1:A:151:ARG:HH21	1.44	0.82
1:D:24:ALA:HB3	4:D:439:HOH:O	1.82	0.80
1:A:1:MET:HE1	1:A:242:ILE:CG2	2.11	0.80
3:A:362:DCS:HND	1:B:303:MET:H	1.31	0.78
1:A:73:LYS:NZ	4:A:458:HOH:O	2.16	0.78
1:C:215:TYR:HE1	4:C:451:HOH:O	1.67	0.77
1:A:1:MET:HE3	1:A:240:SER:HB2	1.66	0.76
1:A:276:ASP:OD1	1:A:352:ARG:HD3	1.84	0.76
1:A:189:ARG:HB2	1:A:189:ARG:HH11	1.52	0.75
1:A:258:THR:HG21	4:A:405:HOH:O	1.86	0.74
1:D:189:ARG:HH11	1:D:189:ARG:HB2	1.53	0.74
1:B:189:ARG:HH11	1:B:189:ARG:HB2	1.53	0.73
3:D:360:DCS:H5A1	3:D:360:DCS:HN	1.50	0.73
1:C:145:THR:HG22	1:C:146:GLN:HE21	1.55	0.72
1:C:258:THR:HB	4:C:475:HOH:O	1.90	0.72
1:C:189:ARG:HH11	1:C:189:ARG:HB2	1.55	0.72
1:B:145:THR:HG22	1:B:146:GLN:HE21	1.55	0.72
1:D:220:LEU:HD21	4:D:386:HOH:O	1.89	0.71
1:A:1:MET:SD	1:B:82:PHE:HB3	2.30	0.71
1:A:168:CYS:O	4:A:516:HOH:O	2.10	0.70
1:D:78:LEU:HD22	3:D:360:DCS:H2A3	1.74	0.69
3:B:362:DCS:O3P	3:B:362:DCS:C5A	2.40	0.69
1:C:105:LEU:O	1:C:109:GLU:HG3	1.93	0.68
1:D:145:THR:HG22	1:D:146:GLN:HE21	1.57	0.68
1:C:6:VAL:HG23	1:C:353:VAL:CG1	2.21	0.67
1:D:220:LEU:HD12	4:D:459:HOH:O	1.95	0.67
1:A:282:ALA:HA	1:A:338:THR:HG22	1.77	0.67
1:D:6:VAL:HG23	1:D:353:VAL:CG1	2.23	0.67
1:A:21:ARG:NH2	4:A:434:HOH:O	2.28	0.66
1:A:47:ARG:NH1	4:A:382:HOH:O	2.28	0.66
1:B:145:THR:HG22	1:B:146:GLN:NE2	2.11	0.66
1:C:145:THR:HG22	1:C:146:GLN:NE2	2.10	0.66
1:C:90:ILE:HG23	1:C:95:PHE:HB2	1.76	0.65
1:C:212:ILE:HG21	1:C:219:PRO:HB3	1.78	0.65
4:C:437:HOH:O	1:D:344:GLU:HG3	1.96	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:73:LYS:HB2	1:C:73:LYS:HZ3	1.61	0.65
1:C:282:ALA:HA	1:C:338:THR:HG22	1.79	0.65
1:D:145:THR:HG22	1:D:146:GLN:NE2	2.12	0.65
3:D:360:DCS:N	3:D:360:DCS:C5A	2.57	0.65
1:A:151:ARG:HG2	1:A:151:ARG:NH2	2.08	0.65
1:C:303:MET:H	3:D:360:DCS:HND	1.45	0.65
1:A:1:MET:N	1:B:82:PHE:HB2	2.11	0.64
1:D:282:ALA:HA	1:D:338:THR:HG22	1.80	0.64
1:D:296:PRO:HB3	4:D:463:HOH:O	1.96	0.64
1:C:274:TYR:HB2	4:C:437:HOH:O	1.96	0.64
1:D:90:ILE:HG23	1:D:95:PHE:HB2	1.78	0.64
1:A:25:PRO:HG3	1:C:184:GLY:O	1.97	0.64
1:B:276:ASP:OD1	1:B:352:ARG:HD3	1.98	0.64
1:C:65:ARG:NH2	4:C:496:HOH:O	2.27	0.63
1:D:212:ILE:HG21	1:D:219:PRO:HB3	1.79	0.63
1:A:90:ILE:HG23	1:A:95:PHE:HB2	1.80	0.63
1:C:246:GLU:HG2	4:C:430:HOH:O	1.98	0.63
1:A:186:PRO:HB3	1:C:25:PRO:HB2	1.80	0.63
1:C:276:ASP:OD1	1:C:352:ARG:HD3	1.98	0.63
1:D:73:LYS:HB2	1:D:73:LYS:HZ3	1.61	0.63
1:D:208:VAL:HG11	4:D:440:HOH:O	1.99	0.63
3:D:360:DCS:HA	4:D:437:HOH:O	1.98	0.63
1:D:230:GLY:HA2	4:D:389:HOH:O	1.98	0.62
1:C:233:PRO:HB3	4:C:439:HOH:O	2.00	0.62
1:B:282:ALA:HA	1:B:338:THR:HG22	1.82	0.62
1:D:276:ASP:OD1	1:D:352:ARG:HD3	2.00	0.62
1:B:90:ILE:HG23	1:B:95:PHE:HB2	1.81	0.61
1:D:105:LEU:O	1:D:109:GLU:HG3	2.00	0.61
1:A:212:ILE:HG21	1:A:219:PRO:HB3	1.82	0.61
1:A:186:PRO:HA	1:C:26:ALA:HB2	1.82	0.61
1:D:43:LEU:HD21	1:D:47:ARG:NH2	2.15	0.61
1:A:21:ARG:NE	4:A:434:HOH:O	2.33	0.61
1:B:212:ILE:HG21	1:B:219:PRO:HB3	1.83	0.60
1:C:50:PRO:O	1:C:73:LYS:HE2	2.01	0.60
1:C:222:ASP:C	1:C:223:ARG:HG2	2.26	0.60
1:A:21:ARG:CZ	4:A:434:HOH:O	2.49	0.60
1:A:43:LEU:HD21	1:A:47:ARG:NH2	2.17	0.60
1:A:50:PRO:O	1:A:73:LYS:HE2	2.02	0.60
1:B:73:LYS:HB2	1:B:73:LYS:HZ3	1.66	0.60
1:D:8:ILE:HG12	1:D:9:ASN:N	2.17	0.60
1:B:102:GLU:OE2	1:B:143:ARG:NH2	2.34	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:360:DCS:HN	3:D:360:DCS:C5A	2.13	0.59
1:D:50:PRO:O	1:D:73:LYS:HE2	2.02	0.59
1:D:188:GLN:HE21	1:D:207:TRP:HE1	1.50	0.59
1:D:102:GLU:OE2	1:D:143:ARG:NH2	2.35	0.59
1:B:73:LYS:HB2	1:B:73:LYS:HZ2	1.67	0.59
1:D:8:ILE:HG12	1:D:234:VAL:CG1	2.32	0.59
1:A:188:GLN:HE21	1:A:207:TRP:HE1	1.48	0.59
1:C:352:ARG:HD3	1:D:59:ARG:NH2	2.17	0.59
1:A:26:ALA:HB2	1:C:186:PRO:HA	1.85	0.58
1:B:50:PRO:O	1:B:73:LYS:HE2	2.03	0.58
1:B:188:GLN:HE21	1:B:207:TRP:HE1	1.48	0.58
1:A:189:ARG:HB2	1:A:189:ARG:NH1	2.17	0.58
1:B:43:LEU:HD21	1:B:47:ARG:NH2	2.18	0.58
1:C:43:LEU:HD21	1:C:47:ARG:NH2	2.18	0.58
1:D:136:GLN:HB3	4:D:373:HOH:O	2.03	0.58
1:C:188:GLN:HE21	1:C:207:TRP:HE1	1.49	0.58
1:A:77:LEU:HB3	4:A:369:HOH:O	2.03	0.58
1:A:222:ASP:O	1:A:223:ARG:CG	2.52	0.57
1:B:222:ASP:O	1:B:223:ARG:HB2	2.02	0.57
1:C:15:HIS:CD2	1:C:232:GLN:HG3	2.40	0.57
1:D:201:PRO:HG2	1:D:202:GLN:HE21	1.68	0.57
1:B:189:ARG:HB2	1:B:189:ARG:NH1	2.19	0.56
1:C:102:GLU:OE2	1:C:143:ARG:NH2	2.38	0.56
1:D:189:ARG:HB2	1:D:189:ARG:NH1	2.18	0.56
1:D:103:GLU:HG2	4:D:448:HOH:O	2.05	0.56
1:D:179:ASN:O	1:D:183:GLU:HG3	2.06	0.56
1:C:189:ARG:HB2	1:C:189:ARG:NH1	2.20	0.56
1:C:61:GLU:HG2	1:C:62:GLU:N	2.20	0.55
1:C:169:GLY:O	1:C:173:LYS:HG3	2.07	0.55
1:D:222:ASP:O	1:D:223:ARG:CG	2.54	0.55
1:C:212:ILE:HG23	1:C:213:ILE:N	2.22	0.55
1:D:8:ILE:CG1	1:D:234:VAL:CG1	2.84	0.55
1:B:15:HIS:HD2	4:B:647:HOH:O	1.88	0.55
1:D:329:LEU:HG	1:D:334:ILE:HD11	1.87	0.55
1:B:179:ASN:O	1:B:183:GLU:HG3	2.05	0.55
1:D:24:ALA:HB2	4:D:440:HOH:O	2.06	0.55
1:C:222:ASP:O	1:C:223:ARG:CG	2.51	0.54
1:B:102:GLU:CD	1:B:143:ARG:HH22	2.16	0.54
1:D:13:LEU:HD13	1:D:234:VAL:HG21	1.89	0.54
1:D:26:ALA:HA	4:D:444:HOH:O	2.07	0.54
1:C:179:ASN:O	1:C:183:GLU:HG3	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:212:ILE:HG23	1:D:213:ILE:N	2.23	0.54
1:A:102:GLU:OE2	1:A:143:ARG:NH2	2.37	0.54
1:D:73:LYS:HB2	1:D:73:LYS:HZ2	1.69	0.54
1:D:108:LEU:HD21	1:D:150:VAL:CG2	2.37	0.54
1:C:73:LYS:HB2	1:C:73:LYS:HZ2	1.70	0.53
1:C:245:ARG:HD2	4:D:410:HOH:O	2.07	0.53
1:C:235:MET:HG2	1:C:236:SER:N	2.23	0.53
1:D:169:GLY:O	1:D:173:LYS:HG3	2.07	0.53
1:D:199:LEU:HG	4:D:386:HOH:O	2.08	0.53
1:C:13:LEU:HD13	1:C:234:VAL:HG21	1.90	0.53
1:B:188:GLN:HE21	1:B:207:TRP:NE1	2.06	0.53
1:C:6:VAL:HG13	1:C:235:MET:HG3	1.91	0.53
1:A:179:ASN:O	1:A:183:GLU:HG3	2.08	0.53
1:B:193:ALA:HB1	3:B:362:DCS:H5A2	1.90	0.53
1:D:8:ILE:HG23	1:D:357:TYR:CE1	2.44	0.53
1:A:235:MET:HG2	1:A:236:SER:N	2.24	0.53
1:C:314:ALA:O	1:C:315:GLN:HB2	2.08	0.53
1:A:98:ALA:HA	1:A:120:TRP:O	2.08	0.52
1:A:188:GLN:HE21	1:A:207:TRP:NE1	2.06	0.52
1:D:102:GLU:CD	1:D:143:ARG:HH22	2.16	0.52
1:A:303:MET:H	3:B:362:DCS:HND	1.56	0.52
1:C:102:GLU:CD	1:C:143:ARG:HH22	2.17	0.52
1:B:13:LEU:HD13	1:B:234:VAL:HG21	1.91	0.52
1:A:13:LEU:HD13	1:A:234:VAL:HG21	1.91	0.52
1:A:102:GLU:CD	1:A:143:ARG:HH22	2.18	0.52
1:C:93:GLN:HE21	1:C:93:GLN:HA	1.75	0.52
1:D:106:ALA:O	1:D:110:GLU:HG2	2.10	0.52
1:A:129:ARG:HD3	1:B:305:MET:SD	2.49	0.52
1:C:188:GLN:HE21	1:C:207:TRP:NE1	2.08	0.52
1:B:245:ARG:HD2	4:B:547:HOH:O	2.09	0.52
1:A:212:ILE:HG23	1:A:213:ILE:N	2.25	0.51
1:B:169:GLY:O	1:B:173:LYS:HG3	2.10	0.51
1:D:220:LEU:HA	4:D:400:HOH:O	2.09	0.51
1:A:182:CYS:HB3	1:A:189:ARG:HD3	1.92	0.51
1:B:1:MET:HG3	1:B:2:GLN:N	2.26	0.51
1:B:189:ARG:HH11	1:B:189:ARG:CB	2.23	0.51
1:D:188:GLN:HE21	1:D:207:TRP:NE1	2.09	0.51
1:D:235:MET:HG2	1:D:236:SER:N	2.25	0.51
1:B:98:ALA:HA	1:B:120:TRP:O	2.10	0.51
1:A:348:ARG:HG2	1:B:348:ARG:HG2	1.93	0.51
1:B:235:MET:HG2	1:B:236:SER:N	2.25	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:43:LEU:HD12	1:D:71:ILE:HD13	1.93	0.51
1:A:6:VAL:HG13	1:A:235:MET:HG3	1.93	0.51
1:A:25:PRO:HB2	1:C:186:PRO:HB3	1.91	0.51
1:D:189:ARG:HH11	1:D:189:ARG:CB	2.23	0.51
1:D:28:LYS:HE3	4:D:406:HOH:O	2.11	0.51
1:B:201:PRO:HG2	1:B:202:GLN:HE21	1.75	0.50
1:C:109:GLU:OE2	1:C:143:ARG:HD2	2.11	0.50
1:A:314:ALA:O	1:A:315:GLN:HB2	2.10	0.50
1:C:43:LEU:HD12	1:C:71:ILE:HD13	1.93	0.50
1:D:6:VAL:HG13	1:D:235:MET:HG3	1.92	0.50
1:B:182:CYS:HB3	1:B:189:ARG:HD3	1.93	0.50
1:D:93:GLN:HE21	1:D:93:GLN:HA	1.76	0.50
1:A:151:ARG:HH21	1:A:151:ARG:CG	2.21	0.50
1:A:169:GLY:O	1:A:173:LYS:HG3	2.12	0.50
1:C:43:LEU:CD1	1:C:71:ILE:HD13	2.42	0.50
1:D:10:ARG:HE	1:D:48:THR:HG22	1.76	0.50
1:B:10:ARG:HH21	1:B:48:THR:HG22	1.77	0.50
1:B:212:ILE:HG23	1:B:213:ILE:N	2.27	0.50
1:A:258:THR:HB	4:A:417:HOH:O	2.10	0.50
1:C:98:ALA:HA	1:C:120:TRP:O	2.11	0.49
1:C:182:CYS:HB3	1:C:189:ARG:HD3	1.94	0.49
1:D:182:CYS:HB3	1:D:189:ARG:HD3	1.94	0.49
1:B:6:VAL:HG13	1:B:235:MET:HG3	1.94	0.49
1:C:289:LEU:HB2	1:C:323:ILE:HB	1.95	0.49
1:D:8:ILE:HG12	1:D:234:VAL:HG13	1.93	0.49
1:D:8:ILE:HG23	1:D:357:TYR:CD1	2.47	0.49
1:D:238:THR:HB	4:D:376:HOH:O	2.12	0.49
1:C:276:ASP:OD1	1:C:352:ARG:CD	2.60	0.49
1:C:348:ARG:HG2	1:D:348:ARG:HG2	1.93	0.49
1:B:93:GLN:HA	1:B:93:GLN:HE21	1.77	0.49
1:C:141:TYR:CE2	1:C:154:VAL:HG22	2.46	0.49
1:D:314:ALA:O	1:D:315:GLN:HB2	2.11	0.49
1:D:325:TRP:HB3	1:D:334:ILE:HD13	1.95	0.49
1:C:201:PRO:HG2	1:C:202:GLN:HE21	1.75	0.49
1:D:186:PRO:HG2	4:D:425:HOH:O	2.13	0.49
1:D:61:GLU:O	1:D:65:ARG:HG3	2.12	0.48
1:A:222:ASP:O	1:A:223:ARG:HG3	2.13	0.48
1:D:43:LEU:CD1	1:D:71:ILE:HD13	2.43	0.48
1:A:305:MET:SD	1:B:129:ARG:HD3	2.53	0.48
1:D:73:LYS:HE3	4:D:428:HOH:O	2.13	0.48
1:D:276:ASP:OD1	1:D:352:ARG:CD	2.61	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:201:PRO:HG2	1:A:202:GLN:HE21	1.77	0.48
1:A:129:ARG:NH1	3:A:362:DCS:O	2.46	0.48
1:D:141:TYR:CE2	1:D:154:VAL:HG22	2.49	0.48
1:C:10:ARG:HE	1:C:48:THR:HG22	1.79	0.48
1:D:43:LEU:HD21	1:D:47:ARG:HH21	1.78	0.48
1:D:109:GLU:OE2	1:D:143:ARG:HD2	2.14	0.48
1:A:105:LEU:O	1:A:109:GLU:HG3	2.13	0.47
1:B:141:TYR:CE2	1:B:154:VAL:HG22	2.49	0.47
1:C:32:VAL:O	4:C:451:HOH:O	2.20	0.47
1:A:1:MET:HG3	1:A:319:GLY:HA2	1.95	0.47
1:B:289:LEU:HB2	1:B:323:ILE:HB	1.96	0.47
1:D:98:ALA:HA	1:D:120:TRP:O	2.14	0.47
1:D:222:ASP:O	1:D:223:ARG:HG3	2.15	0.47
1:A:1:MET:N	1:B:82:PHE:CB	2.76	0.47
1:A:96:HIS:ND1	4:A:365:HOH:O	2.33	0.47
1:C:10:ARG:HE	1:C:48:THR:CG2	2.27	0.47
1:A:15:HIS:HE1	1:A:230:GLY:O	1.97	0.47
1:B:276:ASP:OD1	1:B:352:ARG:CD	2.61	0.47
1:D:8:ILE:CG1	1:D:234:VAL:HG12	2.44	0.47
1:D:10:ARG:HE	1:D:48:THR:CG2	2.27	0.47
1:D:223:ARG:NH1	1:D:336:GLU:OE1	2.48	0.47
1:D:272:MET:O	1:D:306:ILE:HG13	2.15	0.47
1:A:10:ARG:NH2	1:A:44:GLU:OE1	2.40	0.46
1:A:141:TYR:CE2	1:A:154:VAL:HG22	2.49	0.46
1:C:10:ARG:HH21	1:C:48:THR:HG22	1.80	0.46
1:C:193:ALA:O	1:C:197:ILE:HG13	2.15	0.46
1:A:1:MET:H1	1:B:82:PHE:HB2	1.77	0.46
1:D:11:ARG:NH1	1:D:359:ASP:O	2.48	0.46
1:D:15:HIS:HE1	1:D:230:GLY:O	1.98	0.46
1:B:15:HIS:HE1	1:B:230:GLY:O	1.97	0.46
1:A:222:ASP:O	1:A:223:ARG:HG2	2.15	0.46
1:B:109:GLU:OE2	1:B:143:ARG:HD2	2.14	0.46
1:C:43:LEU:HD21	1:C:47:ARG:HH21	1.80	0.46
1:D:277:GLY:HA3	1:D:349:LEU:HA	1.98	0.46
1:D:289:LEU:HB2	1:D:323:ILE:HB	1.98	0.46
1:A:189:ARG:HH11	1:A:189:ARG:CB	2.23	0.46
1:C:93:GLN:HA	1:C:93:GLN:NE2	2.31	0.46
1:C:189:ARG:NH1	1:C:205:PHE:CD2	2.84	0.46
1:A:93:GLN:HA	1:A:93:GLN:HE21	1.80	0.46
1:B:115:GLU:HB2	1:B:116:PRO:HD2	1.98	0.46
1:C:277:GLY:HA3	1:C:349:LEU:HA	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:277:GLY:HA3	1:B:349:LEU:HA	1.99	0.45
1:C:189:ARG:HH11	1:C:189:ARG:CB	2.25	0.45
1:D:193:ALA:O	1:D:197:ILE:HG13	2.17	0.45
1:A:184:GLY:O	1:C:25:PRO:HG3	2.15	0.45
1:A:277:GLY:HA3	1:A:349:LEU:HA	1.99	0.45
1:D:157:VAL:HA	1:D:190:SER:O	2.17	0.45
1:D:8:ILE:HD13	1:D:8:ILE:C	2.41	0.45
1:A:189:ARG:NH1	1:A:205:PHE:CD2	2.85	0.45
1:D:101:ASN:OD1	1:D:104:GLN:HG3	2.17	0.45
1:A:300:ARG:NH1	4:A:477:HOH:O	2.49	0.45
1:D:153:PRO:HB2	4:D:425:HOH:O	2.17	0.45
1:A:43:LEU:HD21	1:A:47:ARG:HH21	1.81	0.45
1:C:276:ASP:OD1	1:D:59:ARG:NH2	2.49	0.44
1:D:8:ILE:HG13	1:D:234:VAL:CG1	2.47	0.44
1:A:101:ASN:OD1	1:A:104:GLN:HG3	2.18	0.44
1:B:93:GLN:HA	1:B:93:GLN:NE2	2.32	0.44
1:D:10:ARG:HH21	1:D:48:THR:HG22	1.81	0.44
1:D:93:GLN:HA	1:D:93:GLN:NE2	2.31	0.44
1:B:189:ARG:NH1	1:B:205:PHE:CD2	2.85	0.44
1:A:16:ASN:O	1:A:20:LEU:HD22	2.18	0.44
1:C:93:GLN:HE21	1:C:93:GLN:CA	2.31	0.44
1:D:189:ARG:NH1	1:D:205:PHE:CD2	2.86	0.44
1:C:108:LEU:HA	1:C:108:LEU:HD23	1.49	0.44
1:A:289:LEU:HB2	1:A:323:ILE:HB	2.00	0.44
1:B:10:ARG:HE	1:B:48:THR:HG22	1.82	0.43
1:C:141:TYR:CE2	1:C:154:VAL:CG2	3.01	0.43
1:D:325:TRP:HB3	1:D:334:ILE:CD1	2.48	0.43
1:C:272:MET:O	1:C:306:ILE:HG13	2.18	0.43
1:D:161:ALA:C	1:D:162:ARG:HG3	2.44	0.43
1:D:194:SER:OG	1:D:213:ILE:HG22	2.18	0.43
1:C:252:PRO:HA	1:C:259:TRP:O	2.18	0.43
3:C:362:DCS:O4P	3:C:362:DCS:C4A	2.67	0.43
1:D:216:GLY:O	1:D:226:GLY:HA3	2.19	0.43
1:D:222:ASP:O	1:D:223:ARG:HG2	2.19	0.43
1:D:255:TYR:HB2	1:D:300:ARG:HD2	2.01	0.43
1:A:223:ARG:NH1	1:A:336:GLU:OE1	2.52	0.43
1:C:8:ILE:HG23	1:C:234:VAL:HG12	2.01	0.43
1:D:252:PRO:HA	1:D:259:TRP:O	2.18	0.43
1:B:115:GLU:HG2	4:B:564:HOH:O	2.18	0.43
3:C:362:DCS:O3	3:C:362:DCS:N	2.52	0.42
1:B:10:ARG:HE	1:B:48:THR:CG2	2.32	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:194:SER:OG	1:A:213:ILE:HG22	2.19	0.42
1:A:93:GLN:HA	1:A:93:GLN:NE2	2.35	0.42
1:B:93:GLN:HE21	1:B:93:GLN:CA	2.31	0.42
1:C:212:ILE:CG2	1:C:219:PRO:HB3	2.48	0.42
1:A:255:TYR:HB2	1:A:300:ARG:HD2	2.02	0.42
1:B:8:ILE:HG23	1:B:234:VAL:HG12	2.01	0.42
1:B:106:ALA:O	1:B:110:GLU:HG2	2.19	0.42
1:C:193:ALA:HB1	3:C:362:DCS:H5A2	2.02	0.41
1:A:161:ALA:C	1:A:162:ARG:HG3	2.45	0.41
1:B:43:LEU:HD21	1:B:47:ARG:HH21	1.84	0.41
1:B:43:LEU:O	1:B:47:ARG:HG3	2.19	0.41
1:D:93:GLN:HE21	1:D:93:GLN:CA	2.32	0.41
1:A:353:VAL:O	1:A:353:VAL:HG23	2.20	0.41
1:B:161:ALA:C	1:B:162:ARG:HG3	2.45	0.41
1:C:10:ARG:NH2	1:C:44:GLU:OE1	2.40	0.41
3:C:362:DCS:HA	1:D:255:TYR:OH	2.20	0.41
1:B:252:PRO:HA	1:B:259:TRP:O	2.20	0.41
1:B:1:MET:C	1:B:3:ALA:H	2.28	0.41
1:C:233:PRO:CG	4:C:439:HOH:O	2.69	0.41
1:B:175:LEU:HD23	1:B:175:LEU:HA	1.94	0.41
1:B:272:MET:O	1:B:306:ILE:HG13	2.20	0.41
1:C:106:ALA:O	1:C:110:GLU:HG2	2.21	0.41
1:C:161:ALA:C	1:C:162:ARG:HG3	2.45	0.41
1:C:194:SER:OG	1:C:213:ILE:HG22	2.21	0.41
1:D:10:ARG:NH2	1:D:44:GLU:OE1	2.41	0.41
1:D:33:VAL:HG21	4:D:392:HOH:O	2.20	0.41
1:D:77:LEU:HB3	4:D:372:HOH:O	2.21	0.41
1:C:157:VAL:HA	1:C:190:SER:O	2.21	0.41
1:C:158:SER:HB3	1:C:191:ILE:HG13	2.03	0.41
1:A:157:VAL:HA	1:A:190:SER:O	2.21	0.40
1:B:212:ILE:CG2	1:B:219:PRO:HB3	2.49	0.40
1:C:34:LYS:HB3	1:C:57:VAL:HA	2.03	0.40
1:C:61:GLU:O	1:C:65:ARG:HG3	2.20	0.40
1:C:212:ILE:CG2	1:C:213:ILE:N	2.84	0.40
1:D:14:ARG:HD2	4:D:403:HOH:O	2.20	0.40
1:D:212:ILE:CG2	1:D:219:PRO:HB3	2.48	0.40
1:C:255:TYR:CZ	1:D:127:MET:HE1	2.57	0.40
1:A:212:ILE:CG2	1:A:219:PRO:HB3	2.49	0.40
1:B:44:GLU:O	1:B:48:THR:CG2	2.59	0.40
1:C:43:LEU:O	1:C:47:ARG:HG3	2.21	0.40
1:C:232:GLN:HA	1:C:233:PRO:HD3	1.86	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:43:LEU:O	1:D:47:ARG:HG3	2.22	0.40
1:D:212:ILE:CG2	1:D:213:ILE:N	2.85	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	356/379 (94%)	341 (96%)	14 (4%)	1 (0%)	36	50
1	B	356/379 (94%)	342 (96%)	14 (4%)	0	100	100
1	C	348/379 (92%)	333 (96%)	15 (4%)	0	100	100
1	D	354/379 (93%)	340 (96%)	14 (4%)	0	100	100
All	All	1414/1516 (93%)	1356 (96%)	57 (4%)	1 (0%)	48	64

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	4	ALA

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	286/303 (94%)	271 (95%)	15 (5%)	21	36

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	286/303 (94%)	272 (95%)	14 (5%)	22	39
1	C	281/303 (93%)	264 (94%)	17 (6%)	17	31
1	D	284/303 (94%)	268 (94%)	16 (6%)	19	33
All	All	1137/1212 (94%)	1075 (94%)	62 (6%)	19	34

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	20	LEU
1	A	59	ARG
1	A	73	LYS
1	A	89	THR
1	A	118	THR
1	A	145	THR
1	A	151	ARG
1	A	154	VAL
1	A	175	LEU
1	A	214	LEU
1	A	248	LYS
1	A	263	ARG
1	A	288	VAL
1	A	313	GLN
1	B	2	GLN
1	B	17	LEU
1	B	48	THR
1	B	59	ARG
1	B	73	LYS
1	B	118	THR
1	B	145	THR
1	B	154	VAL
1	B	214	LEU
1	B	263	ARG
1	B	288	VAL
1	B	313	GLN
1	B	333	ARG
1	B	353	VAL
1	C	17	LEU
1	C	48	THR
1	C	59	ARG
1	C	60	LEU

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Mol	Chain	Res	Type
1	C	61	GLU
1	C	72	THR
1	C	73	LYS
1	C	89	THR
1	C	118	THR
1	C	145	THR
1	C	154	VAL
1	C	175	LEU
1	C	214	LEU
1	C	263	ARG
1	C	288	VAL
1	C	313	GLN
1	C	359	ASP
1	D	8	ILE
1	D	17	LEU
1	D	48	THR
1	D	59	ARG
1	D	72	THR
1	D	73	LYS
1	D	89	THR
1	D	118	THR
1	D	145	THR
1	D	154	VAL
1	D	175	LEU
1	D	214	LEU
1	D	263	ARG
1	D	288	VAL
1	D	313	GLN
1	D	333	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	2	GLN
1	A	15	HIS
1	A	93	GLN
1	A	149	ASN
1	A	188	GLN
1	A	202	GLN
1	B	15	HIS
1	B	93	GLN
1	B	100	HIS

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Mol	Chain	Res	Type
1	B	146	GLN
1	B	149	ASN
1	B	188	GLN
1	B	202	GLN
1	C	93	GLN
1	C	100	HIS
1	C	146	GLN
1	C	179	ASN
1	C	188	GLN
1	C	202	GLN
1	D	15	HIS
1	D	93	GLN
1	D	96	HIS
1	D	146	GLN
1	D	188	GLN
1	D	202	GLN
1	D	313	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	KCX	C	122	1	10,11,12	0.90	0	6,12,14	2.15	2 (33%)
1	KCX	A	122	1	10,11,12	0.93	1 (10%)	6,12,14	2.18	2 (33%)
1	KCX	D	122	1	10,11,12	0.93	1 (10%)	6,12,14	2.15	2 (33%)
1	KCX	B	122	1	10,11,12	1.92	3 (30%)	6,12,14	0.60	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	KCX	C	122	1	-	0/9/10/12	-
1	KCX	A	122	1	-	0/9/10/12	-
1	KCX	D	122	1	-	0/9/10/12	-
1	KCX	B	122	1	-	1/9/10/12	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	122	KCX	CX-NZ	-4.40	1.27	1.35
1	B	122	KCX	OQ1-CX	-2.90	1.16	1.21
1	B	122	KCX	OQ2-CX	-2.27	1.14	1.36
1	D	122	KCX	OQ1-CX	2.04	1.25	1.21
1	A	122	KCX	OQ1-CX	2.02	1.25	1.21

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	122	KCX	OQ1-CX-NZ	-4.78	117.65	124.92
1	C	122	KCX	OQ1-CX-NZ	-4.73	117.73	124.92
1	D	122	KCX	OQ1-CX-NZ	-4.64	117.87	124.92
1	D	122	KCX	CD-CE-NZ	2.28	118.59	112.20
1	A	122	KCX	CD-CE-NZ	2.12	118.15	112.20
1	C	122	KCX	CD-CE-NZ	2.03	117.90	112.20

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	122	KCX	CG-CD-CE-NZ

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry

10 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	B	360	-	4,4,4	0.90	0	6,6,6	0.70	0
3	DCS	C	362	-	22,23,23	0.93	0	27,33,33	1.25	2 (7%)
2	SO4	B	361	-	4,4,4	0.81	0	6,6,6	0.74	0
2	SO4	A	360	-	4,4,4	1.11	0	6,6,6	0.96	0
3	DCS	B	362	-	22,23,23	4.94	13 (59%)	27,33,33	3.09	15 (55%)
2	SO4	A	361	-	4,4,4	0.57	0	6,6,6	0.54	0
3	DCS	A	362	-	22,23,23	3.81	13 (59%)	27,33,33	2.83	14 (51%)
2	SO4	C	361	-	4,4,4	0.64	0	6,6,6	0.67	0
3	DCS	D	360	-	22,23,23	0.94	0	27,33,33	1.30	3 (11%)
2	SO4	C	360	-	4,4,4	0.94	0	6,6,6	0.66	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	DCS	B	362	-	1/1/3/5	7/11/21/21	0/2/2/2
3	DCS	C	362	-	1/1/3/5	4/11/21/21	0/2/2/2
3	DCS	D	360	-	1/1/3/5	4/11/21/21	0/2/2/2
3	DCS	A	362	-	1/1/3/5	2/11/21/21	0/2/2/2

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	362	DCS	CA-C	-12.51	1.43	1.52
3	B	362	DCS	P-O3P	-10.52	1.15	1.54
3	B	362	DCS	P-O4P	-9.85	1.29	1.60
3	A	362	DCS	P-O4P	-8.07	1.34	1.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	362	DCS	P-O2P	-7.32	1.27	1.54
3	B	362	DCS	P-O2P	-7.11	1.28	1.54
3	A	362	DCS	C-ND	-6.39	1.27	1.34
3	A	362	DCS	P-O3P	-5.69	1.33	1.54
3	A	362	DCS	C3-C2	-5.27	1.35	1.41
3	A	362	DCS	P-O1P	-5.17	1.34	1.50
3	A	362	DCS	CA-C	-4.66	1.49	1.52
3	B	362	DCS	C-ND	-4.65	1.29	1.34
3	B	362	DCS	P-O1P	-4.62	1.36	1.50
3	B	362	DCS	C6-N1	-3.75	1.26	1.34
3	B	362	DCS	C3-C2	-3.66	1.37	1.41
3	A	362	DCS	OG-CB	-3.43	1.36	1.44
3	B	362	DCS	C4A-C4	-3.42	1.46	1.52
3	A	362	DCS	C6-N1	-3.00	1.28	1.34
3	B	362	DCS	OG-CB	-2.91	1.37	1.44
3	A	362	DCS	O3-C3	-2.60	1.30	1.36
3	B	362	DCS	O4P-C5A	-2.52	1.35	1.44
3	A	362	DCS	CB-CA	-2.46	1.47	1.53
3	A	362	DCS	C3-C4	-2.32	1.36	1.40
3	B	362	DCS	O3-C3	-2.32	1.31	1.36
3	A	362	DCS	O-C	-2.26	1.19	1.23
3	B	362	DCS	CB-CA	-2.14	1.48	1.53

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	362	DCS	C4A-N-CA	-7.58	99.63	113.84
3	A	362	DCS	C4A-N-CA	-6.29	102.04	113.84
3	A	362	DCS	C2A-C2-C3	-6.01	113.76	120.80
3	B	362	DCS	O4P-P-O1P	5.81	122.16	106.44
3	B	362	DCS	O3P-P-O4P	-4.99	93.66	106.67
3	B	362	DCS	OG-CB-CA	-4.76	98.01	104.80
3	A	362	DCS	O-C-CA	4.66	129.53	126.19
3	A	362	DCS	C4A-C4-C3	4.55	126.03	119.98
3	B	362	DCS	C5A-C5-C6	-4.50	112.02	119.36
3	B	362	DCS	C4A-C4-C3	4.36	125.78	119.98
3	B	362	DCS	C2A-C2-C3	-3.76	116.40	120.80
3	C	362	DCS	C4A-N-CA	-3.70	106.91	113.84
3	A	362	DCS	CB-OG-ND	3.41	111.86	105.92
3	A	362	DCS	C2A-C2-N1	3.33	123.92	117.64
3	D	360	DCS	O4P-P-O1P	3.11	114.84	106.44
3	A	362	DCS	O2P-P-O1P	3.06	122.74	110.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	362	DCS	O-C-ND	-3.05	121.98	125.73
3	B	362	DCS	O4P-C5A-C5	3.05	115.07	109.36
3	D	360	DCS	C4A-N-CA	-3.04	108.14	113.84
3	B	362	DCS	C5A-C5-C4	2.99	128.96	122.67
3	B	362	DCS	C3-C4-C5	-2.97	116.03	118.73
3	A	362	DCS	C3-C4-C5	-2.96	116.04	118.73
3	A	362	DCS	O4P-C5A-C5	2.83	114.65	109.36
3	A	362	DCS	C6-C5-C4	2.73	120.13	118.06
3	C	362	DCS	O4P-P-O1P	2.50	113.20	106.44
3	A	362	DCS	O3P-P-O4P	-2.45	100.28	106.67
3	B	362	DCS	O2P-P-O4P	-2.42	100.37	106.67
3	D	360	DCS	C4A-C4-C5	2.40	122.37	119.75
3	B	362	DCS	O-C-ND	2.33	128.59	125.73
3	B	362	DCS	O3-C3-C4	2.19	124.54	118.18
3	A	362	DCS	O2P-P-O4P	-2.18	100.99	106.67
3	A	362	DCS	O3-C3-C4	2.10	124.27	118.18
3	B	362	DCS	CA-C-ND	-2.08	106.35	107.63
3	B	362	DCS	O2P-P-O1P	2.02	118.70	110.83

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	362	DCS	CA
3	B	362	DCS	CA
3	C	362	DCS	CA
3	D	360	DCS	CA

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	362	DCS	C4-C5-C5A-O4P
3	B	362	DCS	C6-C5-C5A-O4P
3	B	362	DCS	C-CA-N-C4A
3	B	362	DCS	CB-CA-N-C4A
3	C	362	DCS	C4-C5-C5A-O4P
3	C	362	DCS	C6-C5-C5A-O4P
3	D	360	DCS	C3-C4-C4A-N
3	D	360	DCS	C5-C4-C4A-N
3	D	360	DCS	C5A-O4P-P-O2P
3	A	362	DCS	C5-C4-C4A-N
3	B	362	DCS	C5-C4-C4A-N
3	C	362	DCS	C5-C4-C4A-N

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Mol	Chain	Res	Type	Atoms
3	B	362	DCS	C3-C4-C4A-N
3	A	362	DCS	C3-C4-C4A-N
3	C	362	DCS	C3-C4-C4A-N
3	D	360	DCS	C5A-O4P-P-O3P
3	B	362	DCS	C5A-O4P-P-O2P

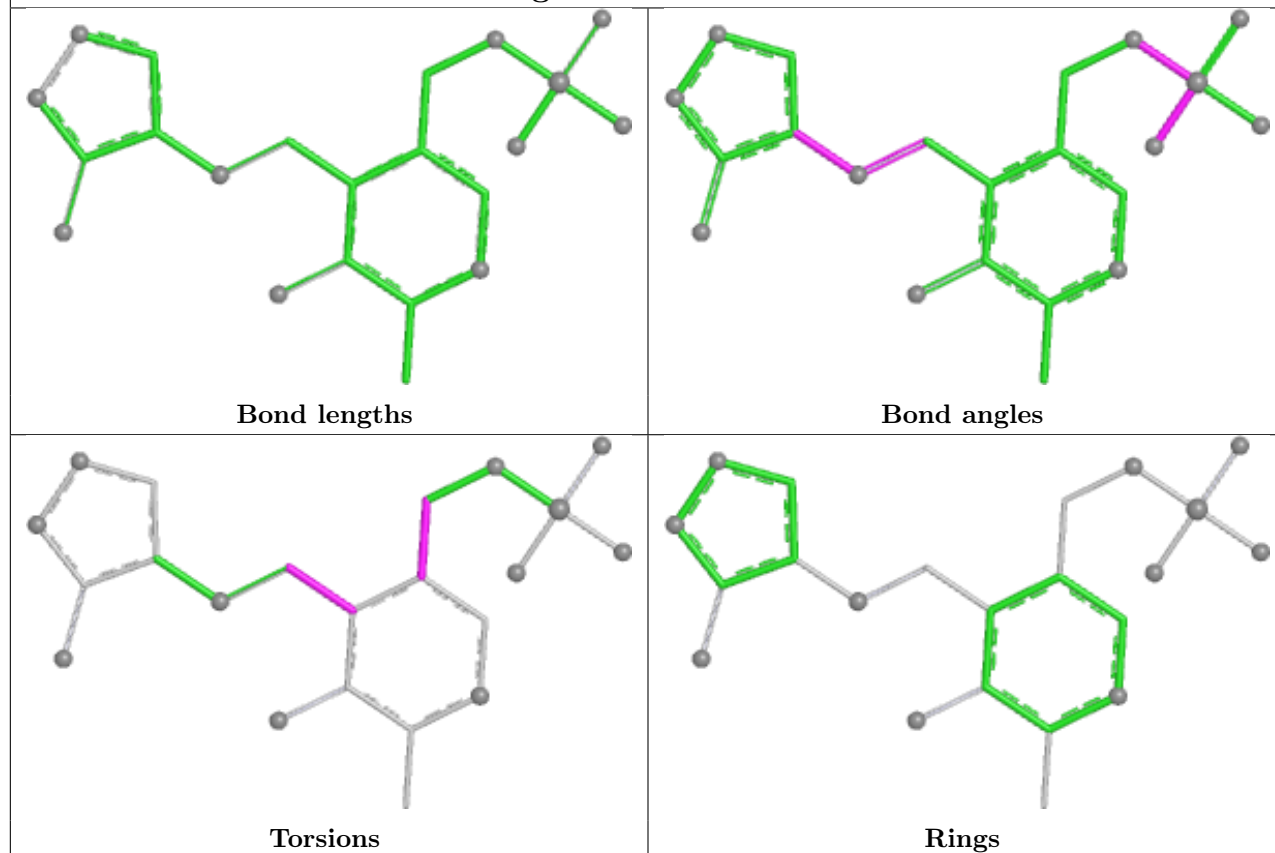
There are no ring outliers.

4 monomers are involved in 20 short contacts:

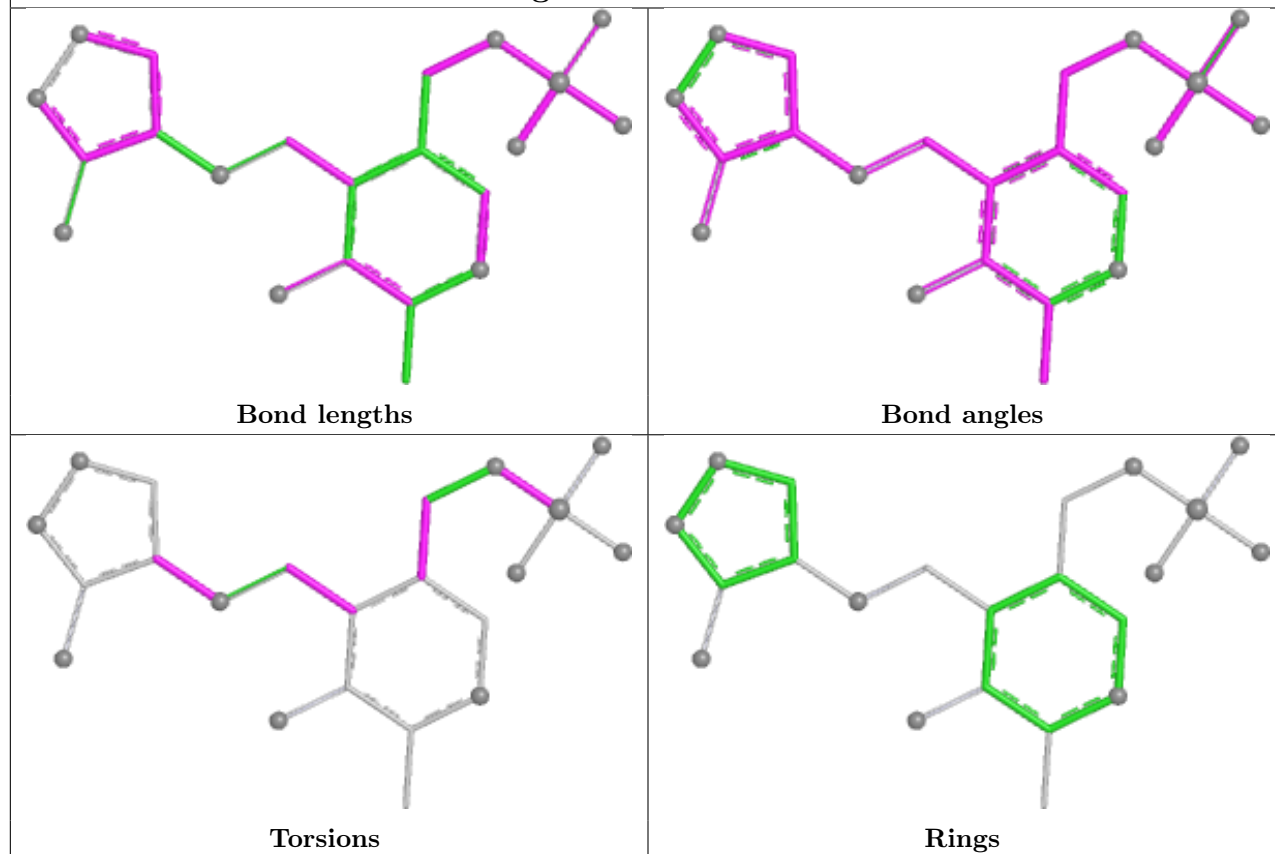
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	362	DCS	5	0
3	B	362	DCS	6	0
3	A	362	DCS	2	0
3	D	360	DCS	7	0

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 than 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.

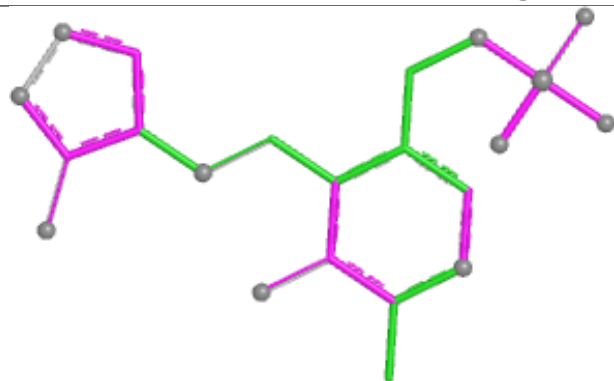
Ligand DCS C 362



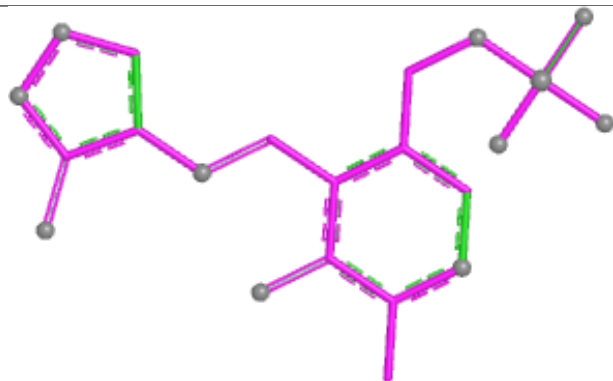
Ligand DCS B 362



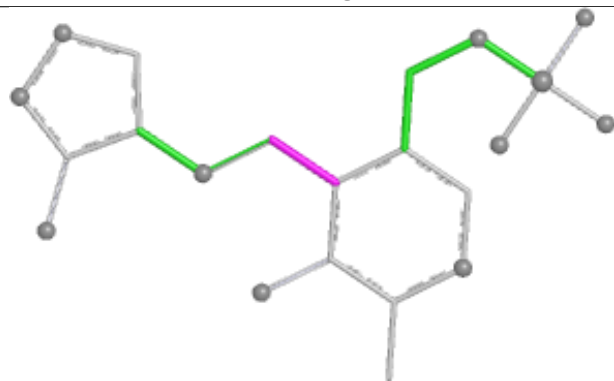
Ligand DCS A 362



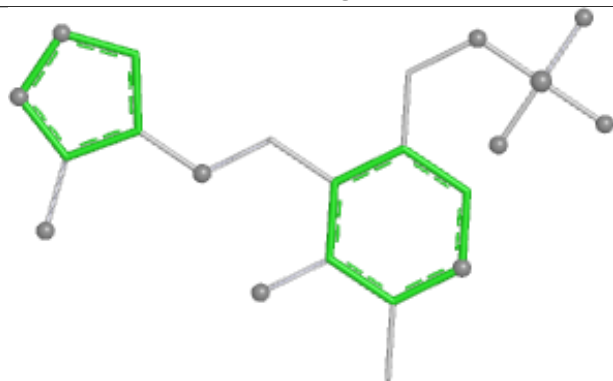
Bond lengths



Bond angles

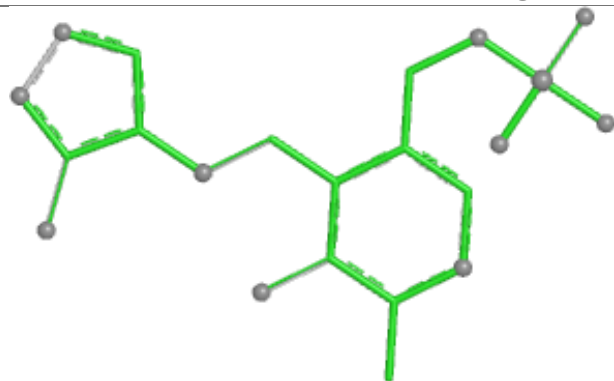


Torsions

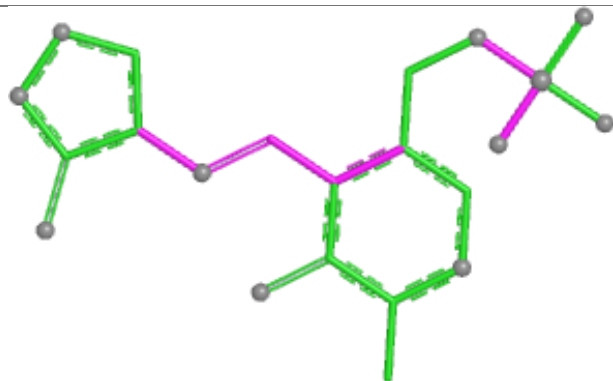


Rings

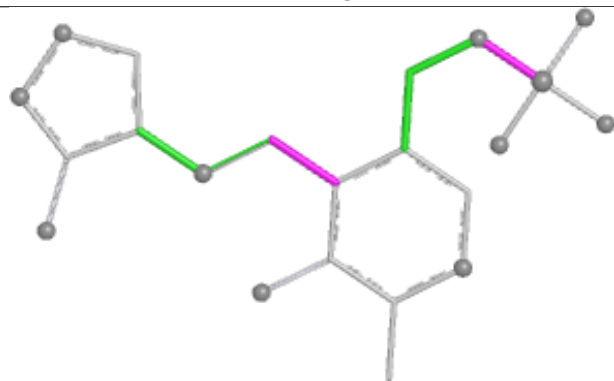
Ligand DCS D 360



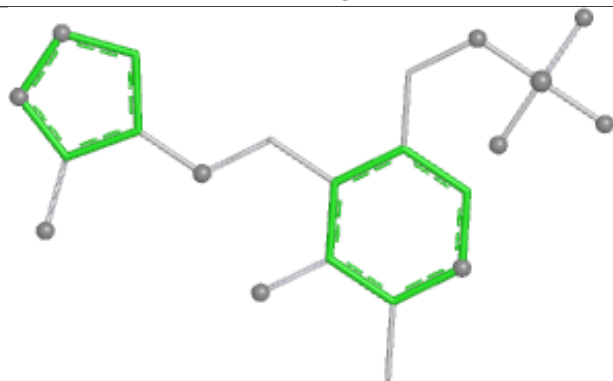
Bond lengths



Bond angles



Torsions



Rings

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	358/379 (94%)	-1.68	0 100 100	2, 10, 24, 43	0
1	B	358/379 (94%)	-1.66	0 100 100	2, 11, 26, 53	0
1	C	352/379 (92%)	-1.42	0 100 100	13, 27, 41, 52	0
1	D	356/379 (93%)	-1.34	0 100 100	10, 30, 47, 53	0
All	All	1424/1516 (93%)	-1.52	0 100 100	2, 20, 41, 53	0

There are no RSRZ outliers to report.

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
1	KCX	A	122	12/13	0.99	0.04	12,16,21,21	0
1	KCX	C	122	12/13	0.99	0.04	17,23,28,28	0
1	KCX	D	122	12/13	0.99	0.04	28,31,38,38	0
1	KCX	B	122	12/13	1.00	0.03	8,10,14,16	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands

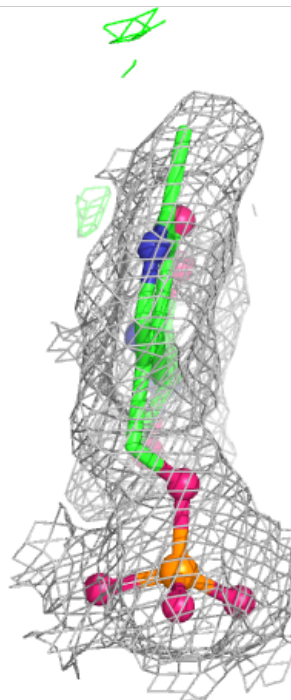
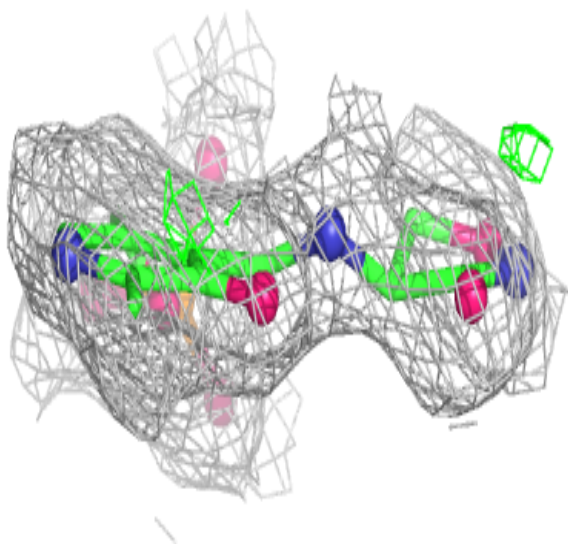
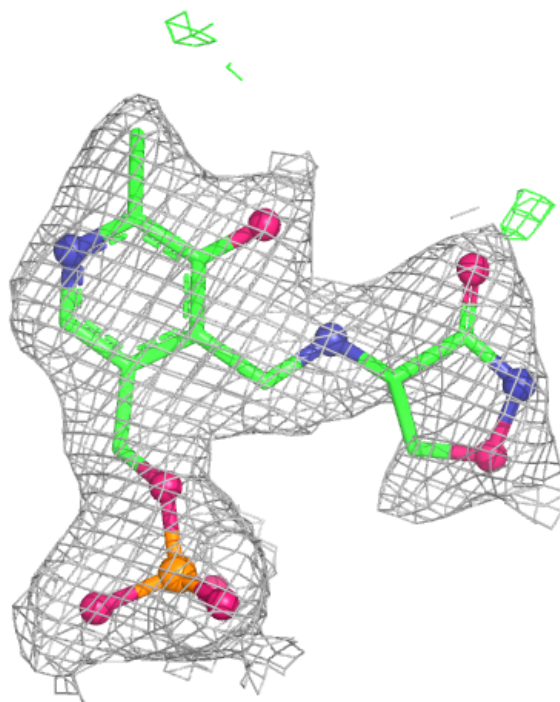
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	SO4	A	361	5/5	0.98	0.10	83,83,84,84	0
2	SO4	C	361	5/5	0.98	0.06	69,69,70,70	0
2	SO4	B	360	5/5	0.99	0.04	56,56,58,58	0
2	SO4	B	361	5/5	0.99	0.05	73,74,74,75	0
2	SO4	C	360	5/5	0.99	0.08	63,64,64,65	0
2	SO4	A	360	5/5	0.99	0.05	53,53,54,55	0
3	DCS	C	362	22/22	0.99	0.03	24,29,31,31	0
3	DCS	D	360	22/22	0.99	0.03	20,28,31,31	0
3	DCS	A	362	22/22	1.00	0.02	6,15,18,19	0
3	DCS	B	362	22/22	1.00	0.02	9,13,16,16	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

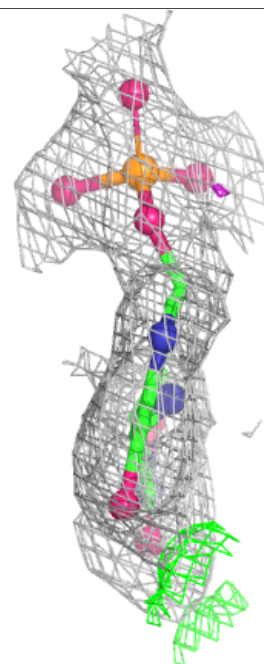
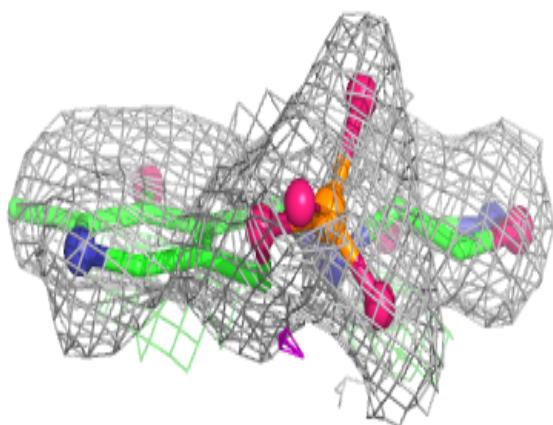
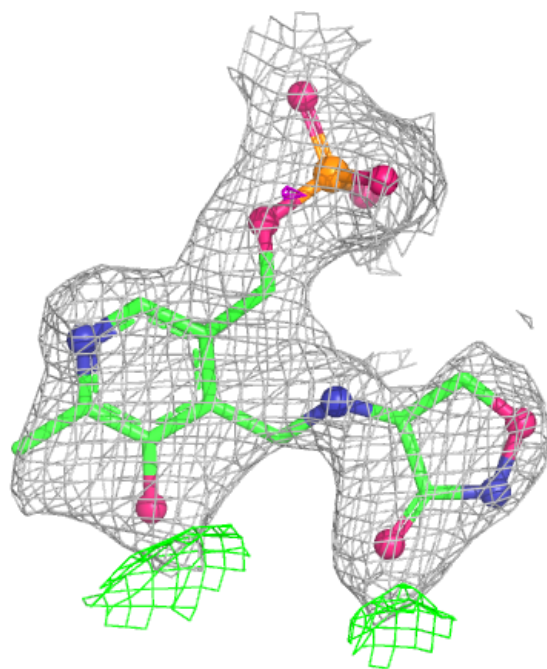
Electron density around DCS C 362:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



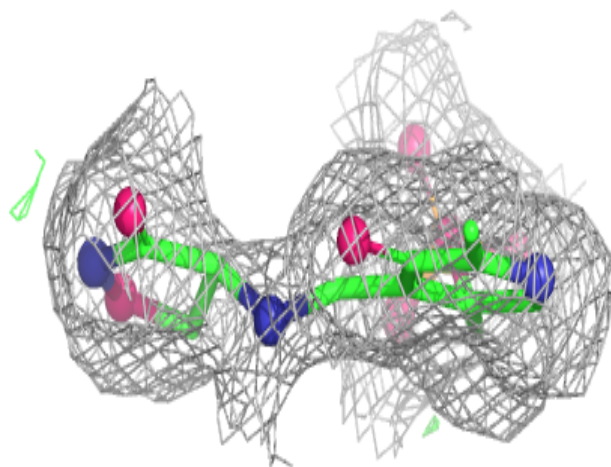
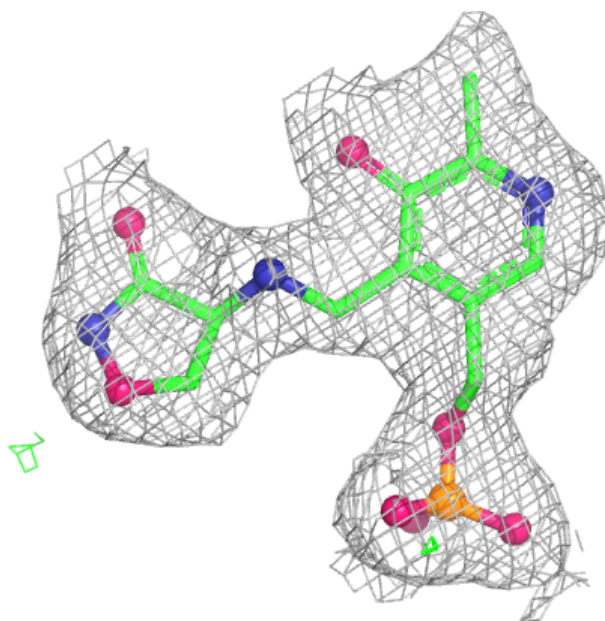
Electron density around DCS D 360:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



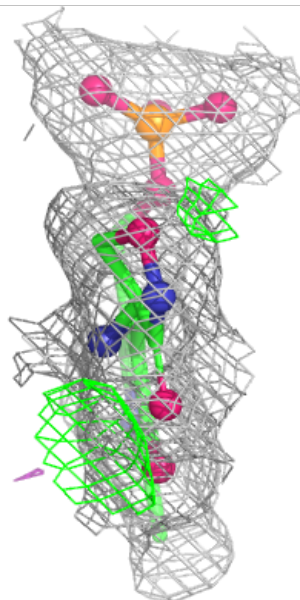
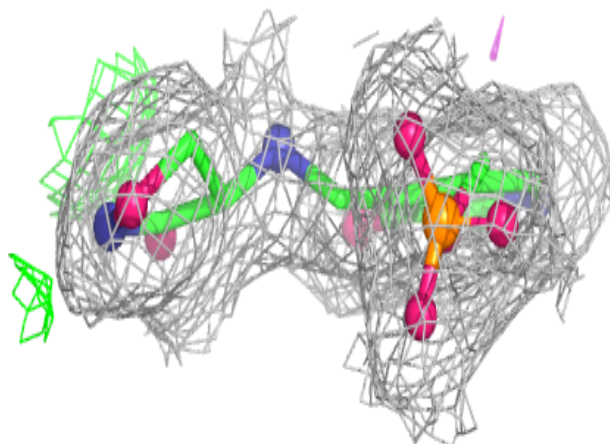
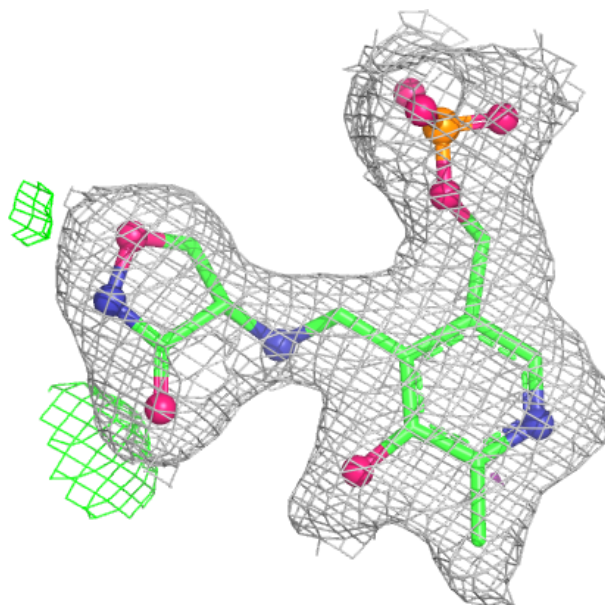
Electron density around DCS A 362:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around DCS B 362:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.