



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:40 AM UTC

PDB ID : 3D6K / pdb_00003d6k
Title : The crystal structure of a putative aminotransferase from *Corynebacterium diphtheriae*
Authors : Tan, K.; Zhang, R.; Duggan, E.; Clancy, S.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2008-05-19
Resolution : 2.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

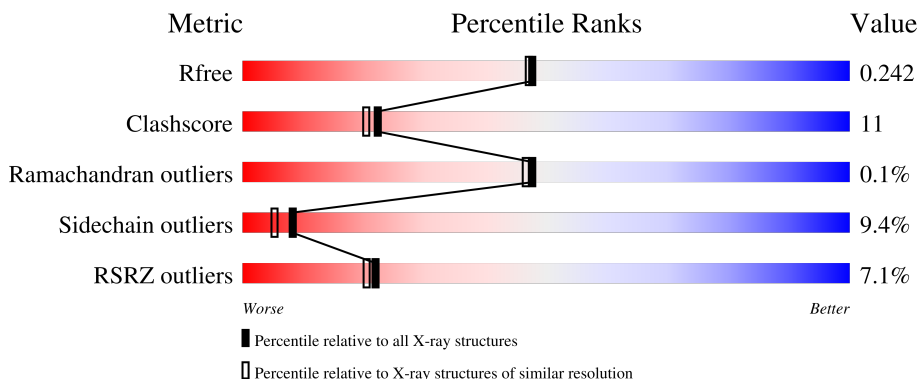
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	422	 3% 72% 21% 5% . .
1	B	422	 12% 70% 22% 5% . .
1	C	422	 10% 76% 17% 5% .
1	D	422	 2% 76% 18% 5% . .

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 13332 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.

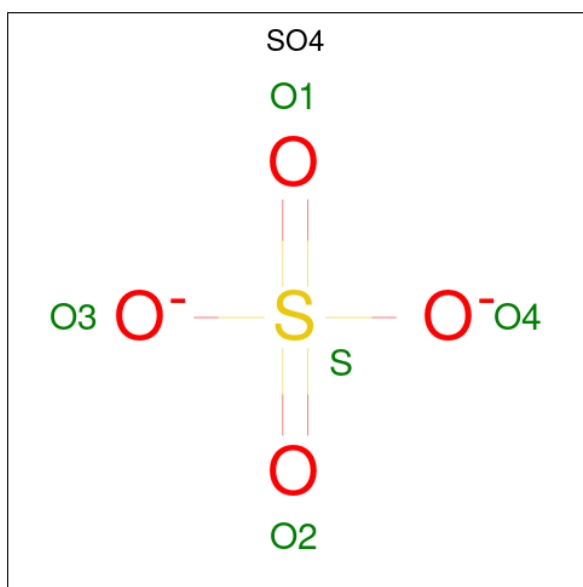
- Molecule 1 is a protein called Putative aminotransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	405	Total	C	N	O	S	Se	0	0	0
			3118	1983	527	594	4	10			
1	B	409	Total	C	N	O	S	Se	0	0	0
			3147	2002	532	599	4	10			
1	C	413	Total	C	N	O	S	Se	0	0	0
			3176	2018	539	605	4	10			
1	D	412	Total	C	N	O	S	Se	0	0	0
			3168	2014	538	602	4	10			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP Q6NJY4
A	-1	ASN	-	expression tag	UNP Q6NJY4
A	0	ALA	-	expression tag	UNP Q6NJY4
B	-2	SER	-	expression tag	UNP Q6NJY4
B	-1	ASN	-	expression tag	UNP Q6NJY4
B	0	ALA	-	expression tag	UNP Q6NJY4
C	-2	SER	-	expression tag	UNP Q6NJY4
C	-1	ASN	-	expression tag	UNP Q6NJY4
C	0	ALA	-	expression tag	UNP Q6NJY4
D	-2	SER	-	expression tag	UNP Q6NJY4
D	-1	ASN	-	expression tag	UNP Q6NJY4
D	0	ALA	-	expression tag	UNP Q6NJY4

- 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	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Cl	0	0
			1	1		
3	B	1	Total	Cl	0	0
			1	1		
3	C	1	Total	Cl	0	0
			1	1		
3	D	1	Total	Cl	0	0
			1	1		

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		
4	C	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	0
			4	2	2		

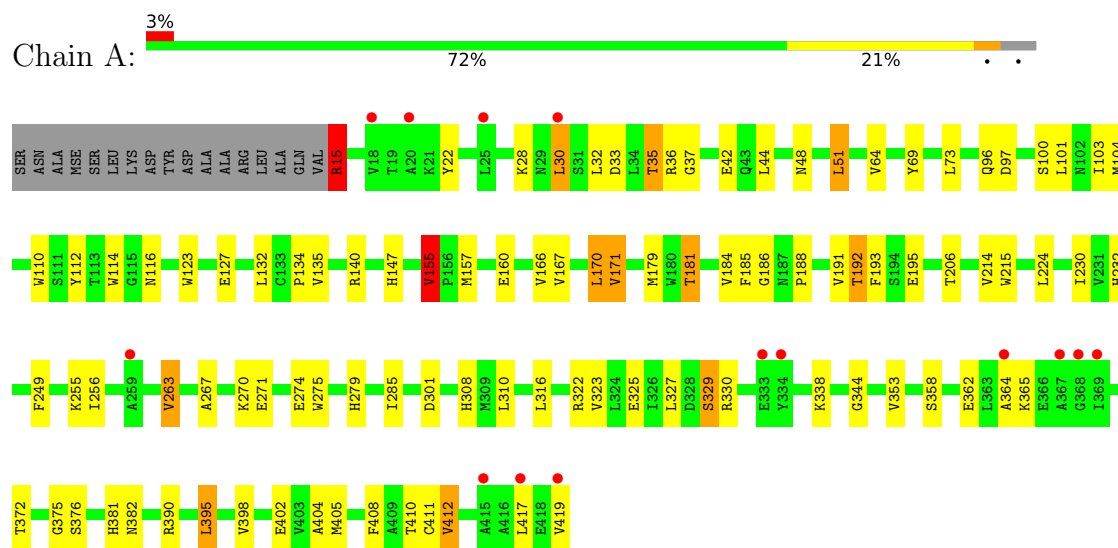
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	211	Total	O	0	0
			211	211		
5	B	182	Total	O	0	0
			182	182		
5	C	164	Total	O	0	0
			164	164		
5	D	126	Total	O	0	0
			126	126		

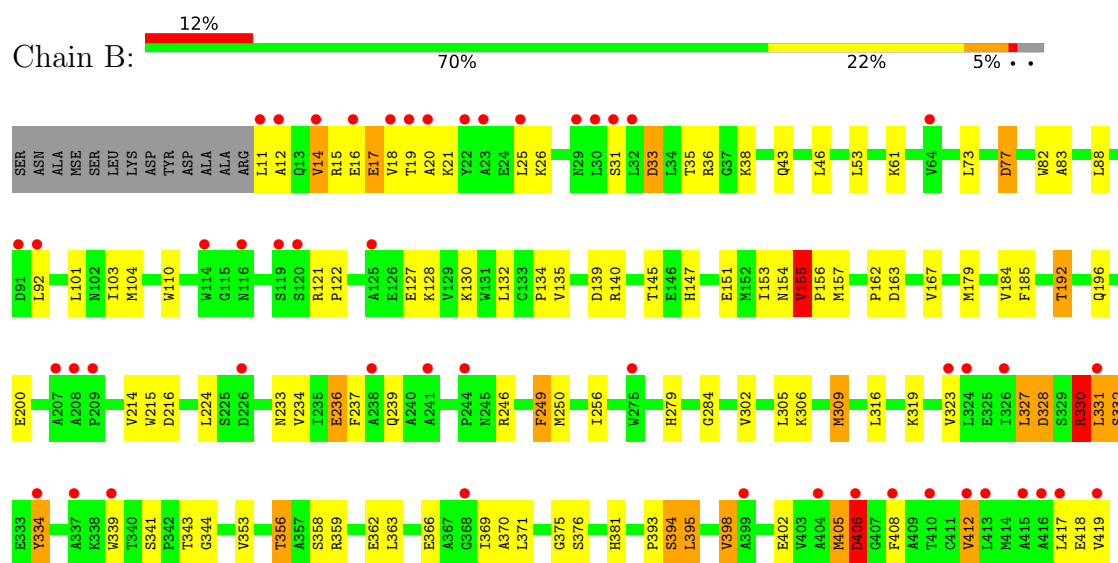
3 Residue-property plots

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.

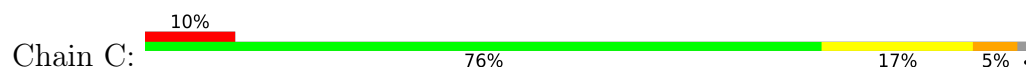
• Molecule 1: Putative aminotransferase

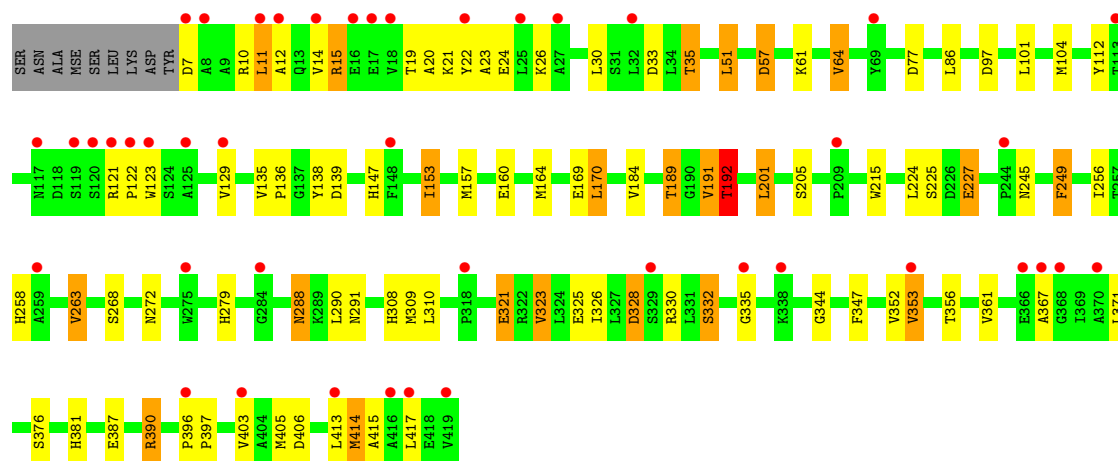


• Molecule 1: Putative aminotransferase

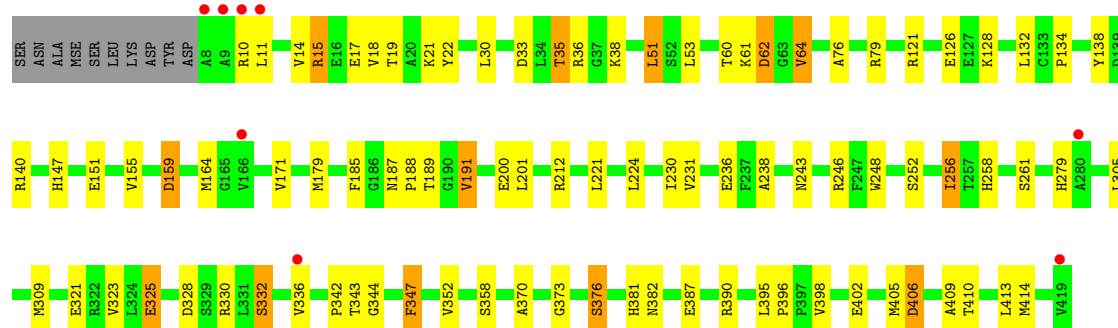
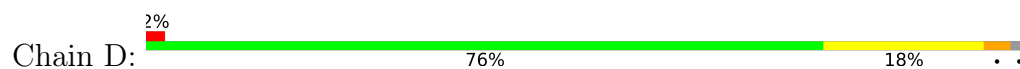


• Molecule 1: Putative aminotransferase





● Molecule 1: Putative aminotransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	89.25Å 56.08Å 173.59Å 90.00° 102.47° 90.00°	Depositor
Resolution (Å)	42.68 – 2.00 42.68 – 2.00	Depositor EDS
% Data completeness (in resolution range)	90.4 (42.68-2.00) 90.4 (42.68-2.00)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.03 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.190 , 0.241 0.191 , 0.242	Depositor DCC
R_{free} test set	5228 reflections (4.59%)	wwPDB-VP
Wilson B-factor (Å ²)	32.3	Xtriage
Anisotropy	0.197	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 33.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.028 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	13332	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.28% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.85	0/3188	1.04	5/4322 (0.1%)
1	B	0.96	7/3217 (0.2%)	1.07	12/4362 (0.3%)
1	C	0.94	4/3246 (0.1%)	1.03	7/4401 (0.2%)
1	D	0.77	0/3238	0.96	4/4390 (0.1%)
All	All	0.88	11/12889 (0.1%)	1.03	28/17475 (0.2%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	406	ASP	C-N	9.08	1.46	1.33
1	C	415	ALA	C-O	8.34	1.34	1.24
1	B	20	ALA	CA-C	8.23	1.63	1.52
1	B	12	ALA	C-N	8.00	1.44	1.33
1	B	16	GLU	C-O	7.98	1.33	1.24
1	C	328	ASP	CG-OD1	6.12	1.36	1.25
1	B	155	VAL	CA-CB	6.06	1.59	1.54
1	C	263	VAL	CA-CB	5.83	1.63	1.54
1	B	319	LYS	C-N	5.73	1.42	1.33
1	B	406	ASP	C-O	-5.17	1.18	1.24
1	C	367	ALA	C-N	5.16	1.41	1.33

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	22	TYR	O-C-N	8.28	131.03	122.09
1	C	22	TYR	CA-C-O	-7.61	112.86	120.70
1	D	256	ILE	N-CA-C	-7.04	106.20	111.90
1	C	192	THR	CB-CA-C	6.63	121.19	109.72
1	B	20	ALA	O-C-N	6.48	128.99	122.12
1	B	341	SER	CA-C-N	6.22	126.73	120.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	341	SER	C-N-CA	6.22	126.73	120.14
1	A	186	GLY	N-CA-C	6.20	118.62	111.36
1	B	154	ASN	CA-C-N	-6.20	117.94	123.33
1	B	154	ASN	C-N-CA	-6.20	117.94	123.33
1	A	155	VAL	CB-CA-C	6.02	115.97	110.13
1	C	57	ASP	CB-CA-C	-5.87	103.31	112.12
1	B	246	ARG	N-CA-C	5.87	117.68	111.28
1	B	12	ALA	CA-C-O	5.75	125.74	119.24
1	D	347	PHE	N-CA-C	5.69	117.83	109.24
1	B	330	ARG	N-CA-C	5.62	120.29	113.38
1	B	394	SER	N-CA-C	5.43	117.20	111.28
1	B	405	MSE	N-CA-C	5.39	117.23	111.36
1	D	396	PRO	O-C-N	5.37	123.67	121.15
1	C	263	VAL	CB-CA-C	5.23	118.56	110.50
1	B	334	TYR	N-CA-C	-5.20	106.94	113.28
1	D	246	ARG	N-CA-C	5.16	117.80	111.82
1	C	64	VAL	N-CA-CB	5.15	116.33	110.72
1	A	15	ARG	CA-C-N	-5.06	111.72	121.58
1	A	15	ARG	C-N-CA	-5.06	111.72	121.58
1	B	33	ASP	N-CA-C	5.04	116.72	108.76
1	A	263	VAL	CB-CA-C	5.02	118.24	110.50
1	C	15	ARG	N-CA-C	5.02	116.60	111.03

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3118	0	3026	59	0
1	B	3147	0	3059	90	0
1	C	3176	0	3086	60	0
1	D	3168	0	3082	66	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	5	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	4	0	6	0	0
4	B	4	0	6	0	0
4	C	4	0	6	1	0
4	D	4	0	6	1	0
5	A	211	0	0	3	0
5	B	182	0	0	5	0
5	C	164	0	0	1	0
5	D	126	0	0	0	0
All	All	13332	0	12277	264	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 11.

All (264) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:104:MSE:HE2	1:B:214:VAL:CG1	1.50	1.39
1:B:104:MSE:HE3	1:B:250:MSE:HG3	1.21	1.18
1:A:15:ARG:HB3	1:A:15:ARG:NH2	1.59	1.15
1:B:104:MSE:HE2	1:B:214:VAL:HG11	1.16	1.12
1:D:22:TYR:HB2	1:D:414:MSE:HE3	1.30	1.12
1:D:256:ILE:HG23	1:D:309:MSE:HE1	1.23	1.09
1:B:104:MSE:HE1	1:B:216:ASP:HB2	1.17	1.09
1:B:33:ASP:HB3	1:B:370:ALA:HB3	1.37	1.07
5:B:511:HOH:O	1:D:382:ASN:HB2	1.61	0.99
1:B:104:MSE:CE	1:B:214:VAL:CG1	2.40	0.99
1:B:185:PHE:H	1:B:192:THR:HG22	1.29	0.97
1:B:309:MSE:HE2	1:B:309:MSE:HA	1.45	0.96
1:D:33:ASP:HB3	1:D:370:ALA:HB3	1.45	0.96
1:B:134:PRO:HB3	1:B:157:MSE:HE3	1.45	0.96
1:D:22:TYR:CB	1:D:414:MSE:HE3	1.96	0.96
1:B:33:ASP:CB	1:B:370:ALA:HB3	1.96	0.95
1:D:11:LEU:O	1:D:15:ARG:HG2	1.68	0.93
1:C:189:THR:HG23	1:C:191:VAL:HG13	1.51	0.93
1:C:256:ILE:HG23	1:C:309:MSE:HE1	1.51	0.92
1:A:184:VAL:HA	1:A:192:THR:HB	1.54	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:11:LEU:O	1:C:14:VAL:HG12	1.71	0.88
1:C:256:ILE:HG23	1:C:309:MSE:CE	2.03	0.88
1:A:147:HIS:NE2	1:B:279:HIS:HD2	1.72	0.87
1:A:15:ARG:HB3	1:A:15:ARG:HH21	1.40	0.87
1:C:326:ILE:O	1:C:330:ARG:HB3	1.74	0.86
1:B:104:MSE:HE1	1:B:216:ASP:CB	2.05	0.86
1:C:279:HIS:HD2	1:D:147:HIS:NE2	1.74	0.85
1:B:104:MSE:HE3	1:B:250:MSE:CG	2.06	0.85
1:B:157:MSE:CE	1:B:162:PRO:HA	2.07	0.84
1:B:104:MSE:HE2	1:B:214:VAL:HG12	1.58	0.83
1:B:104:MSE:CE	1:B:216:ASP:HB2	2.07	0.83
1:D:256:ILE:HG23	1:D:309:MSE:CE	2.07	0.83
1:C:330:ARG:NH1	1:C:406:ASP:OD2	2.11	0.82
1:D:22:TYR:HB2	1:D:414:MSE:CE	2.10	0.81
1:D:330:ARG:HG2	1:D:409:ALA:CB	2.11	0.81
1:A:33:ASP:OD1	1:A:35:THR:HB	1.79	0.81
1:A:279:HIS:HD2	1:B:147:HIS:NE2	1.79	0.80
1:B:132:LEU:HB2	1:B:179:MSE:HG3	1.66	0.78
1:C:33:ASP:OD1	1:C:35:THR:HB	1.83	0.77
1:D:33:ASP:CB	1:D:370:ALA:HB3	2.14	0.77
1:B:104:MSE:CE	1:B:250:MSE:HG3	2.10	0.76
1:B:157:MSE:HE2	1:B:162:PRO:HA	1.67	0.76
1:C:153:ILE:HD11	1:C:170:LEU:HD21	1.69	0.75
1:A:15:ARG:HB3	1:A:15:ARG:CZ	2.17	0.75
1:B:157:MSE:HE1	1:B:162:PRO:CA	2.16	0.75
1:B:309:MSE:HA	1:B:309:MSE:CE	2.16	0.74
1:D:347:PHE:CD1	1:D:390:ARG:HD3	2.23	0.74
1:B:406:ASP:HB2	5:B:489:HOH:O	1.88	0.73
1:B:353:VAL:O	1:B:356:THR:HB	1.88	0.73
1:A:15:ARG:NH2	1:A:15:ARG:CB	2.45	0.73
1:A:15:ARG:CZ	1:A:15:ARG:CB	2.66	0.73
1:B:359:ARG:NH2	1:B:362:GLU:OE1	2.23	0.72
1:C:184:VAL:HA	1:C:192:THR:HB	1.70	0.72
1:D:330:ARG:HG2	1:D:409:ALA:HB1	1.70	0.72
1:B:157:MSE:HE2	1:B:157:MSE:HA	1.71	0.72
1:D:321:GLU:O	1:D:325:GLU:HG2	1.90	0.71
1:C:147:HIS:NE2	1:D:279:HIS:HD2	1.90	0.69
1:C:157:MSE:HE3	1:C:189:THR:HG21	1.74	0.69
1:B:104:MSE:CE	1:B:214:VAL:HG12	2.19	0.67
1:A:147:HIS:NE2	1:B:279:HIS:CD2	2.60	0.67
1:B:184:VAL:HA	1:B:192:THR:HB	1.77	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:330:ARG:O	1:C:330:ARG:HG3	1.97	0.65
1:B:134:PRO:HB3	1:B:157:MSE:CE	2.23	0.65
1:D:159:ASP:OD2	1:D:159:ASP:N	2.29	0.65
1:A:28:LYS:HB3	1:A:30:LEU:HD13	1.79	0.64
1:B:17:GLU:O	1:B:21:LYS:HB2	1.97	0.64
1:B:139:ASP:HB3	1:B:376:SER:O	1.98	0.64
1:D:330:ARG:CG	1:D:409:ALA:HB1	2.28	0.63
1:A:181:THR:HG22	1:A:215:TRP:HA	1.81	0.63
1:A:255:LYS:HE2	5:B:555:HOH:O	1.97	0.63
1:D:330:ARG:HH12	1:D:406:ASP:HA	1.62	0.63
1:A:322:ARG:HD3	1:A:402:GLU:OE2	1.98	0.62
1:B:104:MSE:CE	1:B:214:VAL:HG11	2.10	0.62
1:C:279:HIS:CD2	1:D:147:HIS:NE2	2.63	0.62
1:D:238:ALA:HB1	1:D:243:ASN:O	1.99	0.62
1:A:192:THR:HG23	5:A:581:HOH:O	1.98	0.62
1:D:230:ILE:HD12	1:D:231:VAL:N	2.14	0.62
1:D:330:ARG:HH12	1:D:406:ASP:CA	2.12	0.62
1:D:330:ARG:NH2	1:D:410:THR:OG1	2.33	0.62
1:D:230:ILE:HD12	1:D:231:VAL:H	1.66	0.61
1:C:15:ARG:HD2	1:C:417:LEU:HD22	1.82	0.61
1:D:19:THR:HA	1:D:414:MSE:HE1	1.83	0.61
1:C:347:PHE:CD1	1:C:390:ARG:HD3	2.35	0.61
1:B:331:LEU:HD21	1:B:412:VAL:HG11	1.83	0.60
1:A:166:VAL:HG12	1:A:170:LEU:HD22	1.82	0.60
1:A:69:TYR:HB2	1:B:38:LYS:HD2	1.83	0.60
1:A:112:TYR:HA	1:A:123:TRP:HB2	1.84	0.60
1:A:157:MSE:HE1	1:A:191:VAL:HG11	1.82	0.60
1:B:33:ASP:HB2	1:B:370:ALA:HB3	1.82	0.60
1:A:375:GLY:O	1:A:381:HIS:HA	2.02	0.59
1:A:37:GLY:O	1:A:255:LYS:NZ	2.35	0.59
1:A:271:GLU:HG3	5:A:588:HOH:O	2.02	0.59
1:C:20:ALA:O	1:C:21:LYS:C	2.45	0.59
1:A:181:THR:CG2	1:A:215:TRP:HA	2.33	0.59
1:C:12:ALA:O	1:C:15:ARG:HB2	2.03	0.59
1:C:164:MSE:HE1	1:C:201:LEU:HD13	1.84	0.58
1:B:192:THR:HG23	5:B:599:HOH:O	2.04	0.58
1:B:224:LEU:HG	1:B:344:GLY:HA2	1.86	0.58
1:A:270:LYS:O	1:A:274:GLU:HG3	2.03	0.58
1:C:224:LEU:HG	1:C:344:GLY:HA2	1.85	0.57
1:B:21:LYS:HE2	1:B:330:ARG:NH1	2.20	0.57
1:A:181:THR:HG22	1:A:214:VAL:O	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:212:ARG:HG3	1:D:248:TRP:CZ2	2.40	0.57
1:B:157:MSE:HE1	1:B:162:PRO:HA	1.79	0.57
1:D:330:ARG:HG2	1:D:409:ALA:HB3	1.86	0.57
1:A:185:PHE:H	1:A:192:THR:HG22	1.70	0.56
1:B:157:MSE:CE	1:B:162:PRO:CA	2.75	0.56
1:B:14:VAL:O	1:B:18:VAL:HG13	2.06	0.56
1:B:128:LYS:HD3	1:B:151:GLU:OE1	2.05	0.56
1:C:164:MSE:CE	1:C:201:LEU:HD13	2.36	0.56
1:C:288:ASN:C	1:C:288:ASN:HD22	2.14	0.56
1:C:328:ASP:O	1:C:332:SER:HB2	2.06	0.56
1:C:309:MSE:HA	1:C:309:MSE:HE2	1.89	0.55
1:A:188:PRO:HB3	1:A:390:ARG:HG3	1.89	0.55
1:C:189:THR:HG23	1:C:191:VAL:H	1.71	0.55
1:C:21:LYS:HD2	1:C:330:ARG:NH2	2.21	0.54
1:C:353:VAL:HG23	1:C:356:THR:OG1	2.08	0.54
1:B:121:ARG:HB2	1:B:122:PRO:HD2	1.90	0.54
1:D:132:LEU:HB2	1:D:179:MSE:HG3	1.90	0.54
1:D:330:ARG:NH1	1:D:406:ASP:HA	2.22	0.54
1:B:15:ARG:HG2	1:B:417:LEU:HD22	1.90	0.53
1:A:167:VAL:O	1:A:171:VAL:HB	2.09	0.53
1:B:323:VAL:HG22	1:B:405:MSE:HG3	1.91	0.53
1:D:61:LYS:HG3	1:D:62:ASP:OD1	2.10	0.52
1:B:323:VAL:HG21	1:B:393:PRO:HB3	1.92	0.52
1:D:35:THR:HG22	1:D:36:ARG:HG3	1.92	0.52
1:B:25:LEU:HD11	1:B:369:ILE:HD11	1.92	0.52
1:B:398:VAL:O	1:B:402:GLU:HG3	2.09	0.51
1:C:11:LEU:O	1:C:14:VAL:CG1	2.53	0.51
1:C:288:ASN:ND2	1:C:291:ASN:H	2.07	0.51
1:D:323:VAL:HG22	1:D:405:MSE:HG3	1.91	0.51
1:D:134:PRO:HA	1:D:155:VAL:O	2.09	0.51
1:B:104:MSE:HE2	1:B:214:VAL:HG13	1.74	0.51
1:D:18:VAL:O	1:D:414:MSE:HE1	2.11	0.50
1:B:302:VAL:HG22	1:B:306:LYS:HD2	1.93	0.50
1:C:414:MSE:HE2	1:C:414:MSE:HA	1.93	0.50
1:B:157:MSE:HE1	1:B:162:PRO:CG	2.41	0.50
1:B:302:VAL:HG12	5:B:545:HOH:O	2.11	0.50
1:A:101:LEU:HD23	1:B:284:GLY:HA2	1.93	0.50
1:A:325:GLU:O	1:A:329:SER:HB3	2.12	0.49
1:A:48:ASN:HA	1:A:51:LEU:HD22	1.94	0.49
1:B:216:ASP:HA	1:B:250:MSE:HB2	1.93	0.49
1:B:157:MSE:HE1	1:B:162:PRO:HG3	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:308:HIS:HD2	1:C:309:MSE:CE	2.26	0.49
1:A:181:THR:HG21	1:A:215:TRP:CE3	2.47	0.49
1:A:256:ILE:O	1:A:308:HIS:HE1	1.96	0.49
1:C:169:GLU:OE1	1:C:169:GLU:HA	2.13	0.49
1:C:215:TRP:HB3	1:C:249:PHE:HA	1.94	0.49
1:D:398:VAL:O	1:D:402:GLU:HG3	2.13	0.48
1:C:7:ASP:OD2	1:C:335:GLY:HA3	2.13	0.48
1:A:279:HIS:CD2	1:B:147:HIS:NE2	2.71	0.48
1:A:224:LEU:HG	1:A:344:GLY:HA2	1.94	0.48
1:C:376:SER:HA	1:C:381:HIS:CD2	2.49	0.48
1:D:342:PRO:HG3	1:D:347:PHE:O	2.14	0.48
1:C:101:LEU:HA	1:C:104:MSE:HE3	1.95	0.48
1:C:225:SER:OG	1:C:227:GLU:HG3	2.14	0.48
1:D:328:ASP:O	1:D:332:SER:HB2	2.13	0.47
1:D:164:MSE:HE2	1:D:200:GLU:HB3	1.96	0.47
1:A:132:LEU:HB2	1:A:179:MSE:HB2	1.96	0.47
1:B:305:LEU:O	1:B:309:MSE:HG2	2.14	0.47
1:B:334:TYR:N	1:B:334:TYR:CD1	2.83	0.47
1:C:135:VAL:HA	1:C:136:PRO:C	2.40	0.47
1:B:83:ALA:HB1	1:B:88:LEU:O	2.15	0.47
1:C:139:ASP:HB3	1:C:376:SER:O	2.15	0.47
1:D:376:SER:HA	1:D:381:HIS:CD2	2.51	0.46
1:A:15:ARG:HH21	1:A:15:ARG:CB	2.19	0.46
1:C:112:TYR:HA	1:C:123:TRP:HB2	1.98	0.46
1:B:77:ASP:OD1	1:B:77:ASP:N	2.48	0.46
1:C:121:ARG:HB2	1:C:122:PRO:HD2	1.96	0.46
1:D:11:LEU:HA	1:D:14:VAL:HG12	1.98	0.46
1:B:38:LYS:HB2	1:B:395:LEU:HD12	1.97	0.46
1:B:157:MSE:HE1	1:B:162:PRO:N	2.30	0.46
1:C:290:LEU:HD12	1:D:261:SER:HB2	1.97	0.46
1:A:408:PHE:O	1:A:412:VAL:HG12	2.16	0.46
1:D:18:VAL:HG12	1:D:414:MSE:HE2	1.97	0.46
1:D:347:PHE:CE1	1:D:390:ARG:HD3	2.51	0.46
1:B:418:GLU:O	1:B:419:VAL:C	2.58	0.46
1:C:352:VAL:HG22	1:C:387:GLU:O	2.16	0.46
1:D:138:TYR:CZ	1:D:140:ARG:HB2	2.51	0.46
1:A:116:ASN:HA	1:A:275:TRP:CE2	2.50	0.45
1:B:328:ASP:O	1:B:332:SER:HB2	2.16	0.45
1:D:185:PHE:HE1	1:D:343:THR:CG2	2.28	0.45
1:A:110:TRP:CD1	1:A:114:TRP:HZ3	2.34	0.45
1:A:134:PRO:HA	1:A:155:VAL:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:417:LEU:C	1:B:419:VAL:H	2.25	0.45
1:C:330:ARG:O	1:C:330:ARG:CG	2.64	0.45
1:A:327:LEU:HD21	1:A:405:MSE:HE1	1.98	0.45
1:C:157:MSE:HE1	5:C:498:HOH:O	2.15	0.45
1:A:323:VAL:HG22	1:A:405:MSE:HG3	1.98	0.45
1:A:330:ARG:NH1	1:A:410:THR:OG1	2.50	0.44
1:B:26:LYS:NZ	1:B:366:GLU:O	2.51	0.44
1:C:164:MSE:SE	1:C:201:LEU:HD13	2.68	0.44
1:D:14:VAL:HA	1:D:17:GLU:HG2	1.98	0.44
1:D:138:TYR:CE1	4:D:422:EDO:H21	2.52	0.44
1:A:101:LEU:HD21	1:A:140:ARG:HB3	2.00	0.44
1:B:130:LYS:HD3	1:B:153:ILE:HD11	1.98	0.44
1:D:121:ARG:NH2	1:D:126:GLU:OE1	2.51	0.44
1:D:164:MSE:HE1	1:D:201:LEU:HD23	1.98	0.44
1:B:234:VAL:HA	1:B:237:PHE:HB2	1.99	0.44
1:D:224:LEU:HG	1:D:344:GLY:HA2	1.99	0.44
1:C:245:ASN:HB3	1:C:272:ASN:OD1	2.18	0.44
1:A:32:LEU:HD13	1:A:404:ALA:HB2	2.00	0.43
1:D:10:ARG:O	1:D:14:VAL:HG12	2.18	0.43
1:C:77:ASP:OD1	1:C:77:ASP:N	2.51	0.43
1:D:352:VAL:HG22	1:D:387:GLU:O	2.19	0.43
1:D:51:LEU:HD13	1:D:258:HIS:NE2	2.32	0.43
1:B:233:ASN:ND2	1:B:236:GLU:HB3	2.33	0.43
1:C:189:THR:CG2	1:C:191:VAL:HG13	2.36	0.43
1:C:361:VAL:HA	1:C:371:LEU:HD12	2.00	0.43
1:A:36:ARG:O	5:A:607:HOH:O	2.21	0.43
1:B:110:TRP:CE3	1:B:279:HIS:HB2	2.53	0.43
1:C:321:GLU:O	1:C:325:GLU:HG2	2.19	0.43
1:A:195:GLU:HB2	1:A:232:HIS:CE1	2.53	0.43
1:B:196:GLN:O	1:B:200:GLU:HG3	2.18	0.43
1:C:309:MSE:HE2	1:C:309:MSE:CA	2.49	0.43
1:D:188:PRO:HB3	1:D:390:ARG:HG3	2.00	0.43
1:B:82:TRP:HH2	1:B:256:ILE:CD1	2.32	0.42
1:B:256:ILE:HG23	1:B:309:MSE:CE	2.49	0.42
1:A:193:PHE:N	1:A:193:PHE:CD2	2.86	0.42
1:A:395:LEU:CD1	1:A:395:LEU:C	2.92	0.42
1:D:256:ILE:HG12	1:D:309:MSE:SE	2.69	0.42
1:A:157:MSE:CE	1:A:191:VAL:HG11	2.48	0.42
1:B:43:GLN:O	1:B:46:LEU:HB2	2.19	0.42
1:C:138:TYR:CE1	4:C:422:EDO:H21	2.54	0.42
1:D:60:THR:OG1	1:D:64:VAL:HG13	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:128:LYS:HD3	1:D:151:GLU:OE1	2.19	0.42
1:D:62:ASP:OD1	1:D:62:ASP:N	2.33	0.42
1:B:101:LEU:HD21	1:B:140:ARG:HB3	2.02	0.42
1:B:128:LYS:HB2	1:B:128:LYS:HE3	1.73	0.42
1:C:330:ARG:HH12	1:C:406:ASP:CG	2.26	0.42
1:D:76:ALA:HA	1:D:79:ARG:HH11	1.84	0.42
1:D:187:ASN:HA	1:D:188:PRO:HA	1.78	0.42
1:C:51:LEU:HD13	1:C:258:HIS:NE2	2.35	0.42
1:A:100:SER:O	1:A:104:MSE:HG3	2.19	0.42
1:A:364:ALA:HA	1:A:411:CYS:SG	2.60	0.42
1:A:249:PHE:HB3	1:A:267:ALA:HB3	2.02	0.42
1:B:327:LEU:HB3	1:B:339:TRP:CZ2	2.55	0.42
1:D:138:TYR:CE1	1:D:140:ARG:HB2	2.54	0.42
1:C:308:HIS:CD2	1:C:309:MSE:CE	3.03	0.41
1:D:309:MSE:HE2	1:D:309:MSE:N	2.35	0.41
1:A:22:TYR:HA	1:A:410:THR:HG21	2.03	0.41
1:B:36:ARG:HD2	1:B:395:LEU:HD13	2.01	0.41
1:C:323:VAL:HB	1:C:405:MSE:HG3	2.01	0.41
1:D:305:LEU:HD12	1:D:309:MSE:HE3	2.02	0.41
1:D:189:THR:OG1	1:D:191:VAL:HG13	2.20	0.41
1:B:43:GLN:HA	1:B:46:LEU:HD13	2.03	0.41
1:B:155:VAL:HA	1:B:156:PRO:HD3	1.87	0.41
1:B:408:PHE:O	1:B:412:VAL:HG12	2.20	0.41
1:C:20:ALA:O	1:C:23:ALA:HB3	2.19	0.41
1:B:375:GLY:O	1:B:381:HIS:HA	2.21	0.41
1:B:134:PRO:HA	1:B:155:VAL:O	2.20	0.41
1:C:20:ALA:O	1:C:24:GLU:N	2.49	0.41
1:A:285:ILE:HG12	1:B:140:ARG:CZ	2.51	0.41
1:B:163:ASP:O	1:B:167:VAL:HG23	2.21	0.41
1:B:327:LEU:HD13	1:B:405:MSE:HE2	2.02	0.41
1:C:396:PRO:HA	1:C:397:PRO:HD3	1.91	0.41
1:B:215:TRP:HB3	1:B:249:PHE:HA	2.04	0.40
1:A:96:GLN:HB2	1:A:103:ILE:HD11	2.04	0.40
1:A:171:VAL:HG13	1:A:206:THR:HA	2.02	0.40
1:A:376:SER:HA	1:A:381:HIS:CD2	2.57	0.40
1:D:18:VAL:HG12	1:D:414:MSE:CE	2.52	0.40
1:B:38:LYS:O	1:B:394:SER:OG	2.34	0.40
1:B:103:ILE:HG21	1:B:250:MSE:HE1	2.02	0.40
1:D:373:GLY:O	1:D:376:SER:HB3	2.22	0.40

There are no symmetry-related clashes.

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	403/422 (96%)	389 (96%)	13 (3%)	1 (0%)	43	42
1	B	407/422 (96%)	388 (95%)	19 (5%)	0	100	100
1	C	411/422 (97%)	395 (96%)	15 (4%)	1 (0%)	43	42
1	D	410/422 (97%)	394 (96%)	16 (4%)	0	100	100
All	All	1631/1688 (97%)	1566 (96%)	63 (4%)	2 (0%)	48	46

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	97	ASP
1	C	97	ASP

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	332/334 (99%)	298 (90%)	34 (10%)	7	4
1	B	335/334 (100%)	300 (90%)	35 (10%)	7	4
1	C	337/334 (101%)	303 (90%)	34 (10%)	7	4
1	D	336/334 (101%)	313 (93%)	23 (7%)	14	11
All	All	1340/1336 (100%)	1214 (91%)	126 (9%)	8	5

All (126) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	ARG
1	A	30	LEU
1	A	35	THR
1	A	42	GLU
1	A	44	LEU
1	A	51	LEU
1	A	64	VAL
1	A	73	LEU
1	A	127	GLU
1	A	135	VAL
1	A	155	VAL
1	A	160	GLU
1	A	170	LEU
1	A	171	VAL
1	A	181	THR
1	A	192	THR
1	A	230	ILE
1	A	263	VAL
1	A	301	ASP
1	A	310	LEU
1	A	316	LEU
1	A	329	SER
1	A	338	LYS
1	A	353	VAL
1	A	358	SER
1	A	362	GLU
1	A	365	LYS
1	A	372	THR
1	A	382	ASN
1	A	395	LEU
1	A	398	VAL
1	A	412	VAL
1	A	417	LEU
1	A	419	VAL
1	B	11	LEU
1	B	14	VAL
1	B	17	GLU
1	B	19	THR
1	B	31	SER
1	B	35	THR
1	B	53	LEU
1	B	61	LYS
1	B	73	LEU

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Mol	Chain	Res	Type
1	B	77	ASP
1	B	92	LEU
1	B	127	GLU
1	B	135	VAL
1	B	145	THR
1	B	155	VAL
1	B	192	THR
1	B	236	GLU
1	B	239	GLN
1	B	249	PHE
1	B	309	MSE
1	B	316	LEU
1	B	327	LEU
1	B	328	ASP
1	B	330	ARG
1	B	331	LEU
1	B	332	SER
1	B	343	THR
1	B	356	THR
1	B	358	SER
1	B	363	LEU
1	B	371	LEU
1	B	395	LEU
1	B	398	VAL
1	B	406	ASP
1	B	412	VAL
1	C	10	ARG
1	C	11	LEU
1	C	19	THR
1	C	26	LYS
1	C	30	LEU
1	C	35	THR
1	C	51	LEU
1	C	57	ASP
1	C	61	LYS
1	C	64	VAL
1	C	86	LEU
1	C	129	VAL
1	C	153	ILE
1	C	160	GLU
1	C	170	LEU
1	C	189	THR

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Mol	Chain	Res	Type
1	C	191	VAL
1	C	192	THR
1	C	201	LEU
1	C	205	SER
1	C	227	GLU
1	C	249	PHE
1	C	263	VAL
1	C	268	SER
1	C	288	ASN
1	C	310	LEU
1	C	321	GLU
1	C	323	VAL
1	C	332	SER
1	C	353	VAL
1	C	390	ARG
1	C	403	VAL
1	C	413	LEU
1	C	414	MSE
1	D	15	ARG
1	D	21	LYS
1	D	30	LEU
1	D	35	THR
1	D	38	LYS
1	D	51	LEU
1	D	53	LEU
1	D	62	ASP
1	D	64	VAL
1	D	159	ASP
1	D	171	VAL
1	D	191	VAL
1	D	221	LEU
1	D	236	GLU
1	D	252	SER
1	D	325	GLU
1	D	332	SER
1	D	336	VAL
1	D	358	SER
1	D	376	SER
1	D	395	LEU
1	D	406	ASP
1	D	413	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	GLN
1	A	96	GLN
1	A	232	HIS
1	A	245	ASN
1	A	279	HIS
1	A	308	HIS
1	A	381	HIS
1	B	102	ASN
1	B	233	ASN
1	B	245	ASN
1	B	258	HIS
1	B	279	HIS
1	B	381	HIS
1	C	43	GLN
1	C	102	ASN
1	C	239	GLN
1	C	279	HIS
1	C	288	ASN
1	C	308	HIS
1	C	381	HIS
1	D	96	GLN
1	D	102	ASN
1	D	175	GLN
1	D	239	GLN
1	D	279	HIS
1	D	281	ASN
1	D	381	HIS

5.3.3 RNA ⓘ

There are no RNA molecules 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 oligosaccharides in this entry.

5.6 Ligand geometry

Of 12 ligands modelled in this entry, 4 are monoatomic - 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	$\# Z > 2$
4	EDO	C	422	-	3,3,3	0.60	0	2,2,2	0.26	0
2	SO4	A	420	-	4,4,4	0.40	0	6,6,6	0.35	0
2	SO4	C	420	-	4,4,4	0.47	0	6,6,6	0.56	0
4	EDO	A	422	-	3,3,3	0.75	0	2,2,2	0.36	0
2	SO4	D	420	-	4,4,4	0.54	0	6,6,6	0.22	0
4	EDO	B	422	-	3,3,3	0.59	0	2,2,2	0.33	0
2	SO4	B	420	-	4,4,4	0.31	0	6,6,6	0.30	0
4	EDO	D	422	-	3,3,3	0.77	0	2,2,2	0.46	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	B	422	-	-	1/1/1/1	-
4	EDO	C	422	-	-	0/1/1/1	-
4	EDO	A	422	-	-	0/1/1/1	-
4	EDO	D	422	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	422	EDO	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	422	EDO	1	0
4	D	422	EDO	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	395/422 (93%)	0.51	14 (3%) 47 46	27, 34, 49, 54	0
1	B	399/422 (94%)	1.03	49 (12%) 8 7	27, 35, 51, 55	0
1	C	403/422 (95%)	0.94	43 (10%) 11 10	28, 35, 46, 55	0
1	D	402/422 (95%)	0.51	8 (1%) 65 64	28, 35, 42, 62	0
All	All	1599/1688 (94%)	0.75	114 (7%) 22 20	27, 35, 48, 62	0

All (114) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	404	ALA	5.4
1	B	415	ALA	5.4
1	B	32	LEU	4.9
1	B	20	ALA	4.8
1	C	417	LEU	4.8
1	B	419	VAL	4.8
1	B	11	LEU	4.7
1	A	30	LEU	4.4
1	D	8	ALA	4.3
1	B	417	LEU	4.2
1	D	419	VAL	4.2
1	A	419	VAL	4.1
1	B	326	ILE	3.7
1	B	337	ALA	3.6
1	B	410	THR	3.5
1	B	412	VAL	3.5
1	A	368	GLY	3.4
1	B	406	ASP	3.4
1	B	25	LEU	3.4
1	A	364	ALA	3.4
1	A	367	ALA	3.4

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Mol	Chain	Res	Type	RSRZ
1	D	9	ALA	3.4
1	B	18	VAL	3.4
1	B	22	TYR	3.3
1	B	416	ALA	3.2
1	C	416	ALA	3.2
1	B	23	ALA	3.2
1	C	11	LEU	3.1
1	C	27	ALA	3.0
1	A	334	TYR	3.0
1	C	275	TRP	3.0
1	A	18	VAL	2.9
1	C	12	ALA	2.9
1	A	333	GLU	2.8
1	C	259	ALA	2.8
1	B	275	TRP	2.8
1	C	413	LEU	2.8
1	C	125	ALA	2.7
1	A	417	LEU	2.7
1	C	335	GLY	2.7
1	A	25	LEU	2.7
1	B	19	THR	2.7
1	B	334	TYR	2.7
1	C	284	GLY	2.6
1	B	241	ALA	2.6
1	C	370	ALA	2.6
1	C	14	VAL	2.5
1	C	353	VAL	2.5
1	B	208	ALA	2.5
1	C	32	LEU	2.5
1	C	122	PRO	2.5
1	C	367	ALA	2.5
1	C	396	PRO	2.5
1	B	30	LEU	2.5
1	D	11	LEU	2.4
1	B	16	GLU	2.4
1	C	419	VAL	2.4
1	C	119	SER	2.4
1	C	244	PRO	2.4
1	B	91	ASP	2.4
1	B	238	ALA	2.3
1	C	368	GLY	2.3
1	C	148	PHE	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	129	VAL	2.3
1	B	399	ALA	2.3
1	B	92	LEU	2.3
1	B	119	SER	2.3
1	C	366	GLU	2.3
1	C	403	VAL	2.3
1	C	7	ASP	2.3
1	A	369	ILE	2.3
1	C	69	TYR	2.3
1	B	339	TRP	2.3
1	B	244	PRO	2.2
1	C	113	THR	2.2
1	D	336	VAL	2.2
1	B	29	ASN	2.2
1	D	10	ARG	2.2
1	B	209	PRO	2.2
1	C	117	ASN	2.2
1	C	209	PRO	2.2
1	B	31	SER	2.2
1	B	120	SER	2.2
1	B	226	ASP	2.2
1	B	64	VAL	2.2
1	A	20	ALA	2.2
1	B	12	ALA	2.2
1	B	125	ALA	2.2
1	C	121	ARG	2.2
1	C	123	TRP	2.2
1	D	166	VAL	2.1
1	A	415	ALA	2.1
1	C	338	LYS	2.1
1	B	114	TRP	2.1
1	C	18	VAL	2.1
1	B	207	ALA	2.1
1	B	324	LEU	2.1
1	B	331	LEU	2.1
1	B	413	LEU	2.1
1	C	120	SER	2.1
1	C	329	SER	2.1
1	C	17	GLU	2.1
1	B	368	GLY	2.1
1	B	14	VAL	2.1
1	C	8	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	25	LEU	2.1
1	C	318	PRO	2.1
1	B	323	VAL	2.1
1	A	259	ALA	2.1
1	C	16	GLU	2.0
1	D	280	ALA	2.1
1	B	116	ASN	2.0
1	B	408	PHE	2.0
1	C	22	TYR	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	SO4	A	420	5/5	0.80	0.16	25,30,31,36	5
4	EDO	D	422	4/4	0.84	0.13	27,28,29,32	0
2	SO4	D	420	5/5	0.87	0.15	37,39,39,41	5
4	EDO	C	422	4/4	0.88	0.14	38,42,43,44	0
4	EDO	A	422	4/4	0.89	0.18	25,28,29,29	0
2	SO4	B	420	5/5	0.90	0.15	41,42,45,47	5
2	SO4	C	420	5/5	0.91	0.14	24,27,29,31	5
4	EDO	B	422	4/4	0.93	0.12	37,38,38,39	0
3	CL	A	421	1/1	0.96	0.13	36,36,36,36	0
3	CL	D	421	1/1	0.96	0.17	42,42,42,42	0
3	CL	C	421	1/1	0.97	0.14	39,39,39,39	0
3	CL	B	421	1/1	0.97	0.12	43,43,43,43	0

6.5 Other polymers [i](#)

There are no such residues in this entry.