



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 02:39 am GMT

PDB ID : 1RKW
Title : CRYSTAL STRUCTURE OF THE MULTIDRUG BINDING TRANSCRIPTIONAL REPRESSOR QACR BOUND TO PENTAMADINE
Authors : Murray, D.S.; Schumacher, M.A.; Brennan, R.G.
Deposited on : 2003-11-23
Resolution : 2.62 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

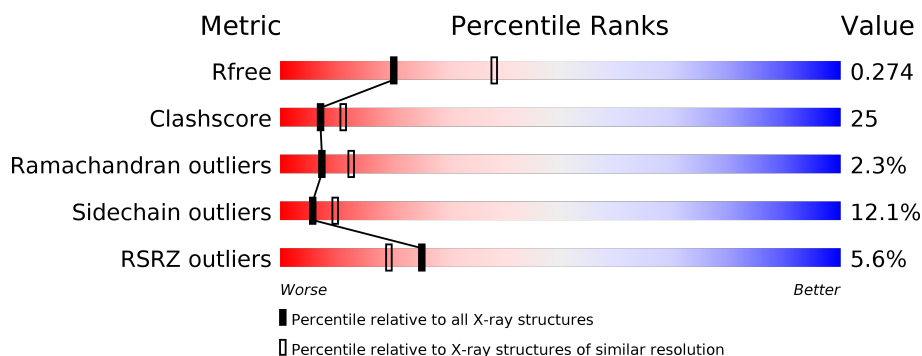
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.62 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}	100719	2983 (2.64-2.60)
Clashscore	112137	3351 (2.64-2.60)
Ramachandran outliers	110173	3298 (2.64-2.60)
Sidechain outliers	110143	3298 (2.64-2.60)
RSRZ outliers	101464	2992 (2.64-2.60)

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. 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	194	
1	B	194	
1	D	194	
1	E	194	

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	SO4	A	199	-	-	-	X
3	SO4	A	279	-	-	-	X
3	SO4	B	259	-	-	-	X
3	SO4	B	299	-	-	-	X
3	SO4	D	639	-	-	-	X
3	SO4	E	239	-	-	-	X
3	SO4	E	899	-	-	-	X
3	SO4	E	999	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6403 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 Transcriptional regulator qacR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	186	Total	C	N	O	S	0	0	0
			1547	998	252	295	2			
1	D	186	Total	C	N	O	S	0	0	0
			1547	998	252	295	2			
1	A	186	Total	C	N	O	S	0	0	0
			1547	998	252	295	2			
1	E	186	Total	C	N	O	S	0	0	0
			1547	998	252	295	2			

There are 32 discrepancies between the modelled and reference sequences:

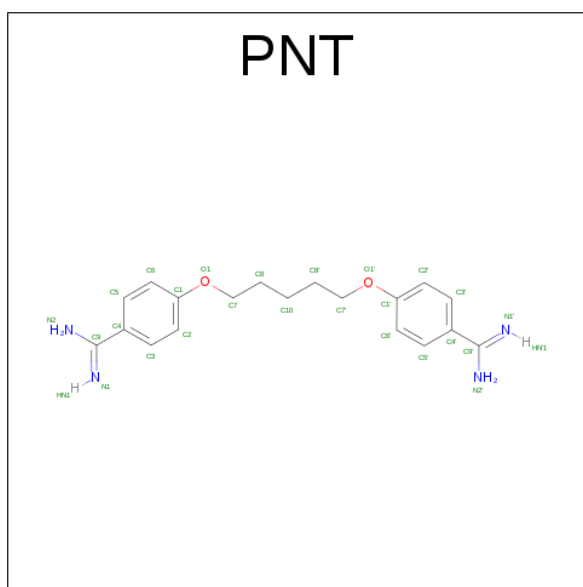
Chain	Residue	Modelled	Actual	Comment	Reference
B	72	ALA	CYS	ENGINEERED	UNP P0A0N4
B	141	SER	CYS	ENGINEERED	UNP P0A0N4
B	189	HIS	-	EXPRESSION TAG	UNP P0A0N4
B	190	HIS	-	EXPRESSION TAG	UNP P0A0N4
B	191	HIS	-	EXPRESSION TAG	UNP P0A0N4
B	192	HIS	-	EXPRESSION TAG	UNP P0A0N4
B	193	HIS	-	EXPRESSION TAG	UNP P0A0N4
B	194	HIS	-	EXPRESSION TAG	UNP P0A0N4
D	72	ALA	CYS	ENGINEERED	UNP P0A0N4
D	141	SER	CYS	ENGINEERED	UNP P0A0N4
D	189	HIS	-	EXPRESSION TAG	UNP P0A0N4
D	190	HIS	-	EXPRESSION TAG	UNP P0A0N4
D	191	HIS	-	EXPRESSION TAG	UNP P0A0N4
D	192	HIS	-	EXPRESSION TAG	UNP P0A0N4
D	193	HIS	-	EXPRESSION TAG	UNP P0A0N4
D	194	HIS	-	EXPRESSION TAG	UNP P0A0N4
A	72	ALA	CYS	ENGINEERED	UNP P0A0N4
A	141	SER	CYS	ENGINEERED	UNP P0A0N4
A	189	HIS	-	EXPRESSION TAG	UNP P0A0N4
A	190	HIS	-	EXPRESSION TAG	UNP P0A0N4
A	191	HIS	-	EXPRESSION TAG	UNP P0A0N4

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Chain	Residue	Modelled	Actual	Comment	Reference
A	192	HIS	-	EXPRESSION TAG	UNP P0A0N4
A	193	HIS	-	EXPRESSION TAG	UNP P0A0N4
A	194	HIS	-	EXPRESSION TAG	UNP P0A0N4
E	72	ALA	CYS	ENGINEERED	UNP P0A0N4
E	141	SER	CYS	ENGINEERED	UNP P0A0N4
E	189	HIS	-	EXPRESSION TAG	UNP P0A0N4
E	190	HIS	-	EXPRESSION TAG	UNP P0A0N4
E	191	HIS	-	EXPRESSION TAG	UNP P0A0N4
E	192	HIS	-	EXPRESSION TAG	UNP P0A0N4
E	193	HIS	-	EXPRESSION TAG	UNP P0A0N4
E	194	HIS	-	EXPRESSION TAG	UNP P0A0N4

- Molecule 2 is 1,5-BIS(4-AMIDINOPHENOXY)PENTANE (three-letter code: PNT) (formula: C₁₉H₂₄N₄O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			25	19	4	2		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

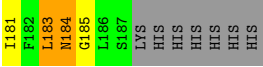
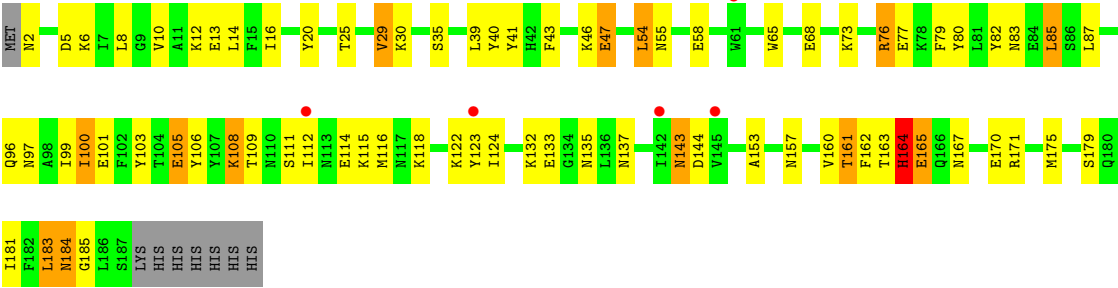
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	35	Total	O	0	0
			35	35		
4	B	10	Total	O	0	0
			10	10		
4	D	12	Total	O	0	0
			12	12		
4	E	23	Total	O	0	0
			23	23		



4 Data and refinement statistics

Property	Value	Source
Space group	P 42 21 2	Depositor
Cell constants a, b, c, α , β , γ	172.05Å 172.05Å 94.43Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	86.03 – 2.62 86.02 – 2.62	Depositor EDS
% Data completeness (in resolution range)	96.0 (86.03-2.62) 96.1 (86.02-2.62)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.77 (at 2.62Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.224 , 0.275 0.224 , 0.274	Depositor DCC
R_{free} test set	2082 reflections (5.04%)	DCC
Wilson B-factor (Å ²)	74.3	Xtriage
Anisotropy	0.005	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 75.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6403	wwPDB-VP
Average B, all atoms (Å ²)	81.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.93% 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, PNT

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	0.76	0/1578	0.78	0/2126
1	B	0.62	0/1578	0.70	0/2126
1	D	0.60	0/1578	0.72	0/2126
1	E	0.67	0/1578	0.77	0/2126
All	All	0.66	0/6312	0.75	0/8504

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1547	0	1536	63	0
1	B	1547	0	1536	113	1
1	D	1547	0	1536	81	0
1	E	1547	0	1536	75	0
2	A	25	0	23	3	0
3	A	25	0	0	2	0
3	B	15	0	0	1	0
3	D	40	0	0	0	0
3	E	30	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	35	0	0	3	0
4	B	10	0	0	0	0
4	D	12	0	0	1	0
4	E	23	0	0	2	0
All	All	6403	0	6167	313	1

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 25.

All (313) 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:E:68:GLU:HG2	1:E:85:LEU:HD21	1.35	1.08
1:E:143:ASN:ND2	1:E:144:ASP:H	1.57	1.03
1:D:157:ASN:O	1:D:161:THR:HG23	1.57	1.01
1:B:36:LYS:H	1:B:36:LYS:HD3	1.27	0.97
1:B:96:GLN:HG3	1:B:157:ASN:HD21	1.29	0.96
1:B:56:ILE:HG22	1:B:60:LYS:HD2	1.47	0.96
1:B:6:LYS:HE3	1:B:6:LYS:HA	1.50	0.93
1:A:85:LEU:O	1:A:89:THR:HG23	1.73	0.88
1:D:165:GLU:HB3	1:E:103:TYR:CE1	2.09	0.87
1:E:157:ASN:O	1:E:161:THR:HG23	1.73	0.87
1:B:106:TYR:HB3	1:B:112:ILE:HG13	1.59	0.84
1:B:167:ASN:ND2	1:B:169:ASN:H	1.77	0.82
1:D:142:ILE:HD11	1:D:186:LEU:HD13	1.59	0.81
1:D:101:GLU:HG3	1:E:97:ASN:HD21	1.46	0.80
1:B:167:ASN:HD22	1:B:168:ILE:N	1.79	0.80
1:A:142:ILE:HD11	1:A:186:LEU:HD13	1.63	0.79
1:A:164:HIS:HA	1:A:171:ARG:CZ	2.12	0.78
1:B:18:ASN:HB2	1:B:23:THR:HG22	1.66	0.78
1:B:4:LYS:O	1:B:8:LEU:HD23	1.84	0.78
1:D:6:LYS:HG3	1:D:32:SER:HB2	1.66	0.76
1:B:58:GLU:HG3	1:B:119:LEU:HD21	1.68	0.75
1:B:47:GLU:OE2	1:B:112:ILE:HD11	1.86	0.75
1:B:39:LEU:H	1:B:39:LEU:HD22	1.51	0.75
1:E:143:ASN:HD22	1:E:144:ASP:H	1.35	0.75
1:B:118:LYS:HD2	1:B:121:ASN:HD22	1.51	0.74
1:B:10:VAL:O	1:B:14:LEU:HD13	1.87	0.73
1:A:100:ILE:O	1:A:104:THR:HG23	1.88	0.73
1:B:72:ALA:HA	1:B:73:LYS:HZ1	1.55	0.72
1:E:106:TYR:HB3	1:E:112:ILE:HG21	1.71	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:163:THR:O	1:A:165:GLU:N	2.22	0.72
1:B:44:LYS:H	1:B:48:ASN:HD22	1.37	0.71
1:D:36:LYS:HG3	1:D:40:TYR:CE1	2.25	0.71
1:A:52:GLU:HG2	3:A:579:SO4:O2	1.90	0.71
1:B:32:SER:O	1:B:33:GLU:HB2	1.89	0.71
1:D:185:GLY:HA2	1:E:184:ASN:HB3	1.73	0.71
1:B:104:THR:O	1:B:105:GLU:HG3	1.92	0.70
1:B:6:LYS:O	1:B:10:VAL:HG23	1.93	0.69
1:D:2:ASN:CG	1:D:3:LEU:H	1.96	0.69
1:D:101:GLU:HG3	1:E:97:ASN:ND2	2.07	0.69
1:D:76:ARG:HG3	1:D:183:LEU:HD13	1.75	0.68
1:B:44:LYS:H	1:B:48:ASN:ND2	1.92	0.68
1:A:62:GLN:HG2	4:A:706:HOH:O	1.93	0.67
1:B:36:LYS:HG2	1:B:37:GLY:H	1.59	0.67
1:B:120:GLU:O	1:B:124:ILE:HG23	1.95	0.67
1:B:14:LEU:HD21	1:B:27:GLU:HG2	1.77	0.67
1:A:96:GLN:O	1:A:100:ILE:HG12	1.95	0.66
1:B:73:LYS:H	1:B:73:LYS:HZ2	1.43	0.66
1:D:2:ASN:ND2	1:D:3:LEU:H	1.93	0.66
1:B:167:ASN:HD21	1:B:169:ASN:H	1.43	0.66
1:B:39:LEU:N	1:B:39:LEU:HD22	2.10	0.66
1:D:184:ASN:HB3	1:E:185:GLY:HA2	1.78	0.66
1:E:97:ASN:O	1:E:100:ILE:HG13	1.96	0.66
1:A:64:GLN:NE2	2:A:225:PNT:H3'	2.11	0.65
1:E:105:GLU:HG2	1:E:106:TYR:CD2	2.31	0.65
1:E:106:TYR:HA	1:E:108:LYS:HE2	1.78	0.65
1:B:103:TYR:HD2	1:B:116:MET:SD	2.19	0.65
1:B:24:THR:HG23	1:B:27:GLU:H	1.61	0.64
1:B:28:ILE:HD13	1:B:49:LEU:HD21	1.79	0.64
1:A:109:THR:O	1:A:113:ASN:HB2	1.98	0.64
1:A:12:LYS:HD2	3:A:199:SO4:O3	1.96	0.64
1:B:167:ASN:C	1:B:167:ASN:HD22	2.00	0.64
1:D:120:GLU:O	1:D:124:ILE:HG12	1.98	0.64
1:D:54:LEU:HD23	1:D:119:LEU:HD13	1.80	0.64
1:E:165:GLU:OE1	1:E:165:GLU:HA	1.95	0.64
1:E:105:GLU:O	1:E:105:GLU:HG3	1.98	0.63
1:B:7:ILE:HD13	1:B:32:SER:OG	1.99	0.63
1:D:165:GLU:HB3	1:E:103:TYR:HE1	1.63	0.63
1:E:132:LYS:HD2	3:E:209:SO4:O3	1.99	0.63
1:B:3:LEU:HD21	1:B:34:SER:OG	1.97	0.63
1:B:72:ALA:HA	1:B:73:LYS:NZ	2.13	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:103:TYR:CE2	1:B:116:MET:HG2	2.34	0.62
1:D:2:ASN:ND2	1:D:3:LEU:N	2.47	0.62
1:B:54:LEU:HB3	1:B:119:LEU:HD11	1.80	0.62
1:B:27:GLU:HG3	1:B:31:LEU:HD23	1.81	0.62
1:B:38:ASN:HB2	1:B:39:LEU:HD22	1.81	0.62
1:B:15:PHE:CE1	1:B:50:PHE:HD2	2.16	0.62
1:D:65:TRP:O	1:D:69:GLN:HB3	2.00	0.62
1:A:64:GLN:HE21	2:A:225:PNT:H3'	1.65	0.62
1:A:97:ASN:ND2	4:A:727:HOH:O	2.33	0.62
1:E:76:ARG:HD3	1:E:183:LEU:HD23	1.82	0.61
1:B:29:VAL:HG13	1:B:34:SER:O	2.01	0.61
1:E:143:ASN:CG	1:E:144:ASP:H	2.03	0.61
1:B:167:ASN:C	1:B:167:ASN:ND2	2.53	0.61
1:D:38:ASN:O	1:D:41:TYR:HB3	2.00	0.61
1:B:177:LYS:HE2	1:A:144:ASP:OD2	2.02	0.60
1:E:6:LYS:O	1:E:10:VAL:HG23	2.00	0.60
1:B:106:TYR:HB3	1:B:112:ILE:CG1	2.31	0.60
1:B:14:LEU:C	1:B:23:THR:HG21	2.21	0.60
1:B:14:LEU:HB3	1:B:23:THR:HB	1.83	0.60
1:B:103:TYR:CD2	1:B:116:MET:HG2	2.35	0.60
1:B:116:MET:O	1:B:116:MET:HG3	2.01	0.60
1:B:103:TYR:CE1	1:B:107:TYR:HB3	2.37	0.59
1:B:144:ASP:OD1	1:A:177:LYS:HE2	2.02	0.59
1:D:73:LYS:HG3	1:D:74:THR:H	1.66	0.59
1:B:39:LEU:CD2	1:B:39:LEU:H	2.15	0.59
1:D:6:LYS:CG	1:D:32:SER:HB2	2.32	0.59
1:A:43:PHE:O	1:A:44:LYS:HB2	2.02	0.59
1:B:112:ILE:C	1:B:114:GLU:H	2.05	0.59
1:B:107:TYR:H	1:B:107:TYR:HD1	1.48	0.59
1:E:114:GLU:O	1:E:118:LYS:HG3	2.02	0.59
1:E:68:GLU:HG2	1:E:85:LEU:CD2	2.23	0.59
1:B:104:THR:C	1:B:105:GLU:HG3	2.24	0.58
1:B:8:LEU:CD2	1:B:43:PHE:HZ	2.16	0.58
1:A:68:GLU:HG2	1:A:85:LEU:HD21	1.85	0.58
1:B:80:TYR:O	1:B:84:GLU:HG3	2.02	0.58
1:D:165:GLU:CB	1:E:103:TYR:CE1	2.84	0.58
1:A:134:GLY:HA3	1:A:140:TRP:CE2	2.39	0.58
1:D:147:ALA:O	1:D:151:ILE:HG13	2.04	0.58
1:D:54:LEU:O	1:D:58:GLU:HB2	2.03	0.58
1:E:164:HIS:CE1	4:E:1021:HOH:O	2.57	0.57
1:B:58:GLU:HG3	1:B:119:LEU:CD2	2.32	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:144:ASP:OD1	1:A:144:ASP:N	2.36	0.57
1:E:157:ASN:O	1:E:161:THR:CG2	2.48	0.57
1:A:90:GLU:HG2	1:A:160:VAL:CG1	2.35	0.57
1:A:164:HIS:HA	1:A:171:ARG:NH2	2.19	0.57
1:B:100:ILE:O	1:B:104:THR:HG23	2.04	0.57
1:B:44:LYS:N	1:B:48:ASN:HD22	2.01	0.57
1:E:108:LYS:HE3	1:E:108:LYS:H	1.69	0.57
1:B:6:LYS:CA	1:B:6:LYS:HE3	2.30	0.56
1:D:125:ASP:O	1:D:129:VAL:HG23	2.04	0.56
1:B:96:GLN:NE2	1:B:157:ASN:OD1	2.37	0.56
1:B:81:LEU:HG	1:B:85:LEU:HD22	1.87	0.56
1:D:177:LYS:HE3	1:D:180:GLN:OE1	2.05	0.56
1:B:36:LYS:CD	1:B:36:LYS:H	2.07	0.56
1:A:103:TYR:C	1:A:105:GLU:H	2.09	0.56
1:D:186:LEU:O	1:D:187:SER:HB2	2.06	0.56
1:B:21:ASN:O	1:B:22:ALA:O	2.24	0.56
1:B:73:LYS:H	1:B:73:LYS:NZ	2.03	0.56
1:B:19:GLY:O	1:B:23:THR:HG23	2.06	0.55
1:E:164:HIS:HA	1:E:171:ARG:NH1	2.20	0.55
1:D:27:GLU:O	1:D:31:LEU:HB2	2.07	0.55
1:B:119:LEU:O	1:B:119:LEU:HD23	2.06	0.55
1:D:165:GLU:CB	1:E:103:TYR:HE1	2.20	0.54
1:D:185:GLY:HA3	1:E:181:ILE:O	2.06	0.54
1:D:45:THR:HG23	1:D:47:GLU:HG3	1.89	0.54
1:A:164:HIS:C	1:A:171:ARG:NH2	2.61	0.54
1:D:167:ASN:ND2	1:D:169:ASN:H	2.06	0.54
1:A:135:ASN:OD1	1:A:141:SER:HA	2.08	0.54
1:E:105:GLU:HG2	1:E:106:TYR:CE2	2.43	0.54
1:E:108:LYS:HE3	1:E:108:LYS:N	2.23	0.54
1:E:83:ASN:HB3	1:E:175:MET:SD	2.48	0.54
1:A:96:GLN:HE21	1:A:100:ILE:HD11	1.74	0.53
1:D:185:GLY:CA	1:E:184:ASN:HB3	2.37	0.53
1:A:113:ASN:HA	1:A:116:MET:HG3	1.91	0.53
1:A:90:GLU:OE2	1:A:96:GLN:HG2	2.08	0.53
1:D:177:LYS:O	1:D:177:LYS:HD3	2.08	0.53
1:A:62:GLN:CG	4:A:706:HOH:O	2.56	0.53
1:D:184:ASN:O	1:D:186:LEU:N	2.42	0.53
1:E:20:TYR:CZ	1:E:46:LYS:HE2	2.43	0.53
1:B:45:THR:OG1	1:B:47:GLU:HG2	2.09	0.53
1:D:186:LEU:O	1:D:187:SER:CB	2.56	0.52
1:E:103:TYR:CD2	1:E:116:MET:HG2	2.45	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:21:ASN:OD1	1:B:105:GLU:OE2	2.27	0.52
1:D:158:GLY:O	1:D:162:PHE:HD1	1.93	0.52
1:E:73:LYS:HB3	1:E:77:GLU:OE1	2.10	0.52
1:A:83:ASN:ND2	1:A:156:VAL:HG21	2.25	0.52
1:B:56:ILE:O	1:B:59:SER:HB2	2.10	0.52
1:D:96:GLN:NE2	1:D:157:ASN:HD21	2.08	0.52
1:B:61:TRP:HZ2	1:B:82:TYR:CE1	2.28	0.52
1:B:112:ILE:O	1:B:116:MET:HB3	2.10	0.52
1:D:144:ASP:HB2	4:D:808:HOH:O	2.10	0.52
1:E:12:LYS:O	1:E:16:ILE:HG13	2.10	0.52
1:A:105:GLU:OE1	1:A:106:TYR:CE2	2.63	0.52
1:E:47:GLU:HG3	1:E:112:ILE:HD11	1.91	0.51
1:D:171:ARG:HH11	1:D:171:ARG:HG2	1.75	0.51
1:E:103:TYR:HD2	1:E:116:MET:HG2	1.75	0.51
1:D:45:THR:CG2	1:D:47:GLU:HG3	2.40	0.51
1:B:61:TRP:HE3	1:B:89:THR:HG21	1.75	0.51
1:B:18:ASN:HB2	1:B:23:THR:CG2	2.40	0.51
1:A:21:ASN:ND2	1:A:101:GLU:OE1	2.45	0.51
1:D:134:GLY:HA3	1:D:140:TRP:CZ2	2.46	0.50
1:E:132:LYS:O	1:E:135:ASN:HB2	2.12	0.50
1:A:38:ASN:ND2	1:E:13:GLU:CD	2.65	0.50
1:B:11:ALA:HB3	1:B:53:ILE:HD11	1.93	0.50
1:A:47:GLU:HG2	1:A:106:TYR:CZ	2.47	0.50
1:B:95:LEU:O	1:B:99:ILE:HG13	2.11	0.50
1:E:96:GLN:O	1:E:100:ILE:HG12	2.11	0.50
1:B:118:LYS:HD2	1:B:121:ASN:ND2	2.24	0.49
1:B:61:TRP:CZ2	1:B:82:TYR:CE1	2.99	0.49
1:B:164:HIS:CD2	1:B:164:HIS:O	2.65	0.49
1:E:35:SER:HA	4:E:1000:HOH:O	2.13	0.49
1:A:3:LEU:O	1:A:7:ILE:HG13	2.13	0.49
1:B:112:ILE:C	1:B:114:GLU:N	2.66	0.49
1:E:76:ARG:CD	1:E:183:LEU:HD23	2.43	0.49
1:A:39:LEU:C	1:A:39:LEU:HD23	2.33	0.49
1:D:127:TYR:O	1:D:131:PHE:CD2	2.65	0.49
1:A:170:GLU:HA	1:A:170:GLU:OE2	2.12	0.49
1:B:110:ASN:ND2	1:B:110:ASN:O	2.46	0.49
1:B:106:TYR:CB	1:B:112:ILE:HG13	2.37	0.49
1:B:132:LYS:O	1:B:136:LEU:HD13	2.13	0.49
1:B:36:LYS:N	1:B:36:LYS:HD3	2.11	0.49
1:D:2:ASN:CG	1:D:3:LEU:N	2.66	0.49
1:D:29:VAL:HG11	1:D:35:SER:HA	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:163:THR:O	1:E:171:ARG:HD3	2.13	0.48
1:A:106:TYR:C	1:A:108:LYS:H	2.16	0.48
1:A:143:ASN:H	1:A:143:ASN:ND2	2.11	0.48
1:A:40:TYR:CE1	1:E:6:LYS:HD3	2.48	0.48
1:E:108:LYS:HE3	1:E:109:THR:H	1.78	0.48
1:E:153:ALA:O	1:E:157:ASN:HB2	2.13	0.48
1:E:41:TYR:CD2	1:E:41:TYR:C	2.85	0.48
1:D:15:PHE:CE2	1:D:23:THR:HB	2.49	0.48
1:E:79:PHE:CD2	1:E:183:LEU:HD13	2.48	0.48
1:A:164:HIS:HA	1:A:171:ARG:NH1	2.28	0.48
1:A:40:TYR:CZ	1:E:6:LYS:HD3	2.49	0.48
1:A:90:GLU:HG2	1:A:160:VAL:HG11	1.95	0.48
1:E:106:TYR:CB	1:E:112:ILE:HG21	2.40	0.48
1:B:56:ILE:O	1:B:60:LYS:HG3	2.14	0.48
1:A:163:THR:O	1:A:166:GLN:HG3	2.13	0.48
1:D:27:GLU:HG2	1:D:31:LEU:HD22	1.95	0.48
1:A:165:GLU:O	1:A:166:GLN:HG2	2.14	0.47
1:E:143:ASN:HD22	1:E:144:ASP:N	2.09	0.47
1:E:54:LEU:HD11	1:E:99:ILE:HG12	1.95	0.47
1:D:150:LYS:HE3	1:D:150:LYS:HB2	1.56	0.47
1:A:49:LEU:HD23	1:A:49:LEU:O	2.14	0.47
1:A:96:GLN:HE21	1:A:100:ILE:CD1	2.27	0.47
1:A:92:TYR:O	1:A:94:PRO:HD3	2.15	0.47
1:D:40:TYR:O	1:D:44:LYS:HA	2.14	0.47
1:D:103:TYR:HE1	1:E:165:GLU:HG2	1.79	0.47
1:B:106:TYR:HB3	1:B:112:ILE:HG21	1.96	0.47
1:B:54:LEU:HB3	1:B:119:LEU:CD1	2.45	0.47
1:B:58:GLU:CG	1:B:119:LEU:HD21	2.42	0.47
1:E:2:ASN:HB2	1:E:5:ASP:HB2	1.97	0.47
1:A:134:GLY:HA3	1:A:140:TRP:CZ2	2.50	0.47
1:D:177:LYS:O	1:D:181:ILE:HG13	2.15	0.47
1:A:57:GLU:OE1	2:A:225:PNT:H6'	2.15	0.46
1:E:39:LEU:C	1:E:39:LEU:HD23	2.35	0.46
1:A:165:GLU:CD	1:A:165:GLU:N	2.69	0.46
1:B:106:TYR:O	1:B:112:ILE:HB	2.15	0.46
1:A:109:THR:HG22	1:A:110:ASN:N	2.30	0.46
1:A:179:SER:O	1:A:183:LEU:HB2	2.15	0.46
1:E:133:GLU:O	1:E:137:ASN:ND2	2.46	0.46
1:D:101:GLU:OE2	1:E:101:GLU:OE2	2.33	0.46
1:B:96:GLN:O	1:B:100:ILE:HG13	2.15	0.46
1:E:76:ARG:NH1	1:E:80:TYR:OH	2.48	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:102:PHE:CZ	1:B:112:ILE:HD12	2.51	0.46
1:D:163:THR:O	1:D:171:ARG:HD2	2.16	0.46
1:A:103:TYR:O	1:A:105:GLU:N	2.48	0.46
1:D:49:LEU:O	1:D:53:ILE:HG13	2.16	0.46
1:E:111:SER:O	1:E:115:LYS:HG3	2.16	0.46
1:E:29:VAL:HG21	1:E:35:SER:C	2.37	0.45
1:A:54:LEU:HB3	1:A:119:LEU:HD21	1.98	0.45
1:A:103:TYR:C	1:A:105:GLU:N	2.68	0.45
1:A:52:GLU:OE1	1:A:52:GLU:HA	2.16	0.45
1:B:167:ASN:ND2	1:B:168:ILE:N	2.57	0.45
1:B:107:TYR:CD1	1:B:107:TYR:N	2.84	0.45
1:B:12:LYS:HD2	3:B:299:SO4:O1	2.17	0.45
1:D:100:ILE:O	1:D:104:THR:HG23	2.16	0.45
1:D:10:VAL:HG21	1:D:32:SER:HB3	1.99	0.45
1:D:43:PHE:O	1:D:44:LYS:HB2	2.16	0.45
1:B:14:LEU:HB3	1:B:23:THR:CB	2.46	0.45
1:B:165:GLU:OE1	1:B:166:GLN:OE1	2.35	0.45
1:B:37:GLY:O	1:B:38:ASN:C	2.54	0.44
1:A:136:LEU:N	1:A:136:LEU:CD1	2.80	0.44
1:A:49:LEU:C	1:A:49:LEU:HD23	2.38	0.44
1:D:177:LYS:C	1:D:177:LYS:HD3	2.37	0.44
1:B:118:LYS:O	1:B:118:LYS:HD2	2.18	0.44
1:B:14:LEU:HB3	1:B:23:THR:CG2	2.47	0.44
1:E:40:TYR:O	1:E:43:PHE:O	2.36	0.44
1:B:36:LYS:HG2	1:B:37:GLY:N	2.27	0.44
1:D:55:ASN:OD1	1:D:119:LEU:HD21	2.18	0.44
1:D:177:LYS:CE	1:D:180:GLN:OE1	2.66	0.44
1:D:39:LEU:HD23	1:D:39:LEU:C	2.38	0.44
1:E:12:LYS:HD2	3:E:899:SO4:O3	2.18	0.44
1:D:117:ASN:O	1:D:121:ASN:ND2	2.51	0.44
1:A:55:ASN:HD22	1:A:122:LYS:NZ	2.16	0.43
1:A:165:GLU:CD	1:A:165:GLU:H	2.22	0.43
1:B:142:ILE:HD11	1:B:186:LEU:HD13	1.99	0.43
1:B:21:ASN:ND2	1:B:101:GLU:CG	2.82	0.43
1:D:84:GLU:CD	1:D:176:ASN:HD21	2.21	0.43
1:D:29:VAL:HG13	1:D:34:SER:C	2.39	0.43
1:A:103:TYR:CZ	1:A:116:MET:HE2	2.54	0.43
1:B:28:ILE:HD13	1:B:49:LEU:CD2	2.47	0.43
1:B:40:TYR:HA	1:B:43:PHE:O	2.18	0.43
1:D:166:GLN:HG2	1:D:166:GLN:H	1.55	0.43
1:D:73:LYS:HG3	1:D:74:THR:N	2.32	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:3:LEU:HD13	1:D:3:LEU:O	2.19	0.43
1:D:96:GLN:HG3	1:E:162:PHE:CE2	2.53	0.43
1:D:96:GLN:HG3	1:E:162:PHE:HE2	1.84	0.43
1:B:147:ALA:O	1:B:151:ILE:HG13	2.19	0.42
1:B:3:LEU:C	1:B:3:LEU:HD13	2.39	0.42
1:D:127:TYR:CE2	1:D:153:ALA:HA	2.53	0.42
1:D:93:TYR:N	1:D:94:PRO:CD	2.82	0.42
1:E:179:SER:O	1:E:183:LEU:HB2	2.19	0.42
1:D:57:GLU:HG3	1:D:95:LEU:HD12	2.01	0.42
1:D:156:VAL:HG13	1:D:175:MET:CE	2.49	0.42
1:D:108:LYS:O	1:D:110:ASN:N	2.52	0.42
1:B:14:LEU:CD2	1:B:27:GLU:HG2	2.46	0.42
1:B:93:TYR:N	1:B:94:PRO:CD	2.82	0.42
1:D:52:GLU:O	1:D:56:ILE:HG12	2.20	0.42
1:D:142:ILE:CD1	1:D:186:LEU:HD13	2.41	0.42
1:D:57:GLU:HG3	1:D:95:LEU:CD1	2.49	0.42
1:E:122:LYS:C	1:E:124:ILE:H	2.22	0.42
1:E:65:TRP:CH2	1:E:82:TYR:HB2	2.55	0.41
1:A:134:GLY:HA3	1:A:140:TRP:NE1	2.35	0.41
1:D:183:LEU:HA	1:D:183:LEU:HD23	1.88	0.41
1:D:36:LYS:HG3	1:D:40:TYR:CD1	2.55	0.41
1:A:142:ILE:CD1	1:A:186:LEU:HD13	2.42	0.41
1:B:114:GLU:C	1:B:114:GLU:OE1	2.59	0.41
1:B:15:PHE:CD1	1:B:50:PHE:HD2	2.38	0.41
1:E:108:LYS:CE	1:E:108:LYS:H	2.34	0.41
1:B:16:ILE:HG13	1:B:95:LEU:CD2	2.51	0.41
1:D:6:LYS:O	1:D:10:VAL:HG23	2.21	0.41
1:B:35:SER:HB3	1:B:38:ASN:OD1	2.20	0.41
1:E:122:LYS:C	1:E:124:ILE:N	2.74	0.41
1:E:8:LEU:HA	1:E:8:LEU:HD23	1.91	0.41
1:B:45:THR:H	1:B:48:ASN:HB2	1.85	0.41
1:E:14:LEU:HD23	1:E:14:LEU:HA	1.88	0.41
1:E:87:LEU:HD23	1:E:160:VAL:HG22	2.03	0.41
1:D:123:TYR:OH	1:D:157:ASN:HB2	2.21	0.40
1:D:127:TYR:O	1:D:131:PHE:HD2	2.03	0.40
1:B:8:LEU:CD2	1:B:43:PHE:CZ	3.01	0.40
1:E:108:LYS:CA	1:E:108:LYS:HE3	2.52	0.40
1:B:144:ASP:OD2	1:B:144:ASP:N	2.53	0.40

All (1) 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	Interatomic distance (Å)	Clash overlap (Å)
1:B:42:HIS:CE1	1:B:42:HIS:CE1[8_445]	2.08	0.12

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	184/194 (95%)	172 (94%)	10 (5%)	2 (1%)	17	32
1	B	184/194 (95%)	150 (82%)	27 (15%)	7 (4%)	4	5
1	D	184/194 (95%)	153 (83%)	25 (14%)	6 (3%)	4	7
1	E	184/194 (95%)	169 (92%)	13 (7%)	2 (1%)	17	32
All	All	736/776 (95%)	644 (88%)	75 (10%)	17 (2%)	7	13

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	22	ALA
1	A	164	HIS
1	E	143	ASN
1	B	44	LYS
1	B	106	TYR
1	D	109	THR
1	D	185	GLY
1	A	104	THR
1	E	164	HIS
1	B	23	THR
1	B	33	GLU
1	B	164	HIS
1	D	44	LYS
1	B	89	THR
1	D	33	GLU
1	D	184	ASN
1	D	186	LEU

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	169/177 (96%)	151 (89%)	18 (11%)	8	13
1	B	169/177 (96%)	143 (85%)	26 (15%)	3	4
1	D	169/177 (96%)	151 (89%)	18 (11%)	8	13
1	E	169/177 (96%)	149 (88%)	20 (12%)	6	10
All	All	676/708 (96%)	594 (88%)	82 (12%)	6	10

All (82) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	6	LYS
1	B	23	THR
1	B	36	LYS
1	B	39	LEU
1	B	51	LEU
1	B	58	GLU
1	B	64	GLN
1	B	73	LYS
1	B	85	LEU
1	B	89	THR
1	B	107	TYR
1	B	109	THR
1	B	110	ASN
1	B	114	GLU
1	B	116	MET
1	B	118	LYS
1	B	120	GLU
1	B	124	ILE
1	B	128	HIS
1	B	140	TRP
1	B	146	ASN
1	B	157	ASN
1	B	165	GLU
1	B	167	ASN

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Mol	Chain	Res	Type
1	B	173	LYS
1	B	180	GLN
1	D	2	ASN
1	D	3	LEU
1	D	31	LEU
1	D	89	THR
1	D	97	ASN
1	D	104	THR
1	D	108	LYS
1	D	109	THR
1	D	116	MET
1	D	120	GLU
1	D	124	ILE
1	D	141	SER
1	D	150	LYS
1	D	161	THR
1	D	167	ASN
1	D	170	GLU
1	D	177	LYS
1	D	184	ASN
1	A	30	LYS
1	A	44	LYS
1	A	48	ASN
1	A	54	LEU
1	A	85	LEU
1	A	89	THR
1	A	90	GLU
1	A	105	GLU
1	A	113	ASN
1	A	115	LYS
1	A	116	MET
1	A	118	LYS
1	A	136	LEU
1	A	143	ASN
1	A	157	ASN
1	A	165	GLU
1	A	167	ASN
1	A	173	LYS
1	E	25	THR
1	E	29	VAL
1	E	30	LYS
1	E	47	GLU

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Mol	Chain	Res	Type
1	E	54	LEU
1	E	55	ASN
1	E	58	GLU
1	E	76	ARG
1	E	85	LEU
1	E	100	ILE
1	E	105	GLU
1	E	108	LYS
1	E	123	TYR
1	E	161	THR
1	E	164	HIS
1	E	165	GLU
1	E	167	ASN
1	E	170	GLU
1	E	183	LEU
1	E	184	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (43) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	21	ASN
1	B	48	ASN
1	B	83	ASN
1	B	96	GLN
1	B	97	ASN
1	B	110	ASN
1	B	121	ASN
1	B	154	ASN
1	B	157	ASN
1	B	164	HIS
1	B	167	ASN
1	D	2	ASN
1	D	18	ASN
1	D	75	ASN
1	D	83	ASN
1	D	96	GLN
1	D	121	ASN
1	D	143	ASN
1	D	154	ASN
1	D	167	ASN
1	D	176	ASN
1	A	21	ASN

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Mol	Chain	Res	Type
1	A	38	ASN
1	A	55	ASN
1	A	64	GLN
1	A	69	GLN
1	A	83	ASN
1	A	96	GLN
1	A	113	ASN
1	A	121	ASN
1	A	143	ASN
1	A	164	HIS
1	A	166	GLN
1	A	180	GLN
1	A	184	ASN
1	E	18	ASN
1	E	62	GLN
1	E	83	ASN
1	E	96	GLN
1	E	97	ASN
1	E	121	ASN
1	E	143	ASN
1	E	184	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

23 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
3	SO4	A	199	-	4,4,4	0.28	0	6,6,6	0.20	0
2	PNT	A	225	-	26,26,26	2.36	8 (30%)	29,33,33	1.69	6 (20%)
3	SO4	A	279	-	4,4,4	0.29	0	6,6,6	0.12	0
3	SO4	A	399	-	4,4,4	0.33	0	6,6,6	0.15	0
3	SO4	A	579	-	4,4,4	0.31	0	6,6,6	0.11	0
3	SO4	A	699	-	4,4,4	0.31	0	6,6,6	0.15	0
3	SO4	B	259	-	4,4,4	0.28	0	6,6,6	0.24	0
3	SO4	B	269	-	4,4,4	0.30	0	6,6,6	0.12	0
3	SO4	B	299	-	4,4,4	0.30	0	6,6,6	0.08	0
3	SO4	D	219	-	4,4,4	0.35	0	6,6,6	0.07	0
3	SO4	D	229	-	4,4,4	0.33	0	6,6,6	0.09	0
3	SO4	D	289	-	4,4,4	0.32	0	6,6,6	0.08	0
3	SO4	D	499	-	4,4,4	0.33	0	6,6,6	0.08	0
3	SO4	D	539	-	4,4,4	0.33	0	6,6,6	0.06	0
3	SO4	D	639	-	4,4,4	0.34	0	6,6,6	0.09	0
3	SO4	D	779	-	4,4,4	0.31	0	6,6,6	0.10	0
3	SO4	D	799	-	4,4,4	0.38	0	6,6,6	0.09	0
3	SO4	E	209	-	4,4,4	0.34	0	6,6,6	0.11	0
3	SO4	E	239	-	4,4,4	0.29	0	6,6,6	0.22	0
3	SO4	E	249	-	4,4,4	0.34	0	6,6,6	0.09	0
3	SO4	E	599	-	4,4,4	0.35	0	6,6,6	0.17	0
3	SO4	E	899	-	4,4,4	0.27	0	6,6,6	0.09	0
3	SO4	E	999	-	4,4,4	0.31	0	6,6,6	0.11	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	SO4	A	199	-	-	0/0/0/0	0/0/0/0
2	PNT	A	225	-	-	0/18/18/18	0/2/2/2
3	SO4	A	279	-	-	0/0/0/0	0/0/0/0
3	SO4	A	399	-	-	0/0/0/0	0/0/0/0
3	SO4	A	579	-	-	0/0/0/0	0/0/0/0
3	SO4	A	699	-	-	0/0/0/0	0/0/0/0
3	SO4	B	259	-	-	0/0/0/0	0/0/0/0
3	SO4	B	269	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	SO4	B	299	-	-	0/0/0/0	0/0/0/0
3	SO4	D	219	-	-	0/0/0/0	0/0/0/0
3	SO4	D	229	-	-	0/0/0/0	0/0/0/0
3	SO4	D	289	-	-	0/0/0/0	0/0/0/0
3	SO4	D	499	-	-	0/0/0/0	0/0/0/0
3	SO4	D	539	-	-	0/0/0/0	0/0/0/0
3	SO4	D	639	-	-	0/0/0/0	0/0/0/0
3	SO4	D	779	-	-	0/0/0/0	0/0/0/0
3	SO4	D	799	-	-	0/0/0/0	0/0/0/0
3	SO4	E	209	-	-	0/0/0/0	0/0/0/0
3	SO4	E	239	-	-	0/0/0/0	0/0/0/0
3	SO4	E	249	-	-	0/0/0/0	0/0/0/0
3	SO4	E	599	-	-	0/0/0/0	0/0/0/0
3	SO4	E	899	-	-	0/0/0/0	0/0/0/0
3	SO4	E	999	-	-	0/0/0/0	0/0/0/0

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	225	PNT	C4-C9	-6.56	1.36	1.47
2	A	225	PNT	C4'-C9'	-5.05	1.39	1.47
2	A	225	PNT	C6'-C1'	2.13	1.43	1.38
2	A	225	PNT	C9'-N1'	2.37	1.36	1.28
2	A	225	PNT	C6-C1	2.79	1.44	1.38
2	A	225	PNT	C3'-C4'	2.97	1.44	1.39
2	A	225	PNT	C9'-N2'	3.27	1.41	1.33
2	A	225	PNT	C3-C4	3.90	1.45	1.39

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	225	PNT	O1-C7-C8	-4.36	91.88	108.34
2	A	225	PNT	C10-C8-C7	-3.98	95.57	113.48
2	A	225	PNT	C7'-O1'-C1'	-2.97	110.03	117.90
2	A	225	PNT	C10-C8'-C7'	-2.29	103.20	113.48
2	A	225	PNT	C4-C9-N2	2.50	121.96	118.12
2	A	225	PNT	C4'-C9'-N2'	2.86	122.51	118.12

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	199	SO4	1	0
2	A	225	PNT	3	0
3	A	579	SO4	1	0
3	B	299	SO4	1	0
3	E	209	SO4	1	0
3	E	899	SO4	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	186/194 (95%)	0.65	6 (3%)	48	41	41, 59, 107, 126	0
1	B	186/194 (95%)	1.00	20 (10%)	6	4	46, 84, 165, 175	0
1	D	186/194 (95%)	0.61	11 (5%)	23	17	53, 84, 115, 127	0
1	E	186/194 (95%)	0.59	5 (2%)	55	48	46, 70, 104, 118	0
All	All	744/776 (95%)	0.71	42 (5%)	25	19	41, 72, 126, 175	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	112	ILE	10.5
1	B	34	SER	9.4
1	B	108	LYS	7.1
1	B	41	TYR	6.0
1	B	42	HIS	5.7
1	B	107	TYR	4.8
1	B	117	ASN	4.7
1	B	40	TYR	4.6
1	A	107	TYR	4.1
1	B	116	MET	4.1
1	B	7	ILE	3.9
1	B	103	TYR	3.9
1	B	120	GLU	3.8
1	B	187	SER	3.8
1	D	123	TYR	3.6
1	B	39	LEU	3.6
1	A	113	ASN	3.6
1	E	123	TYR	3.4
1	D	186	LEU	3.4
1	A	112	ILE	3.3
1	E	112	ILE	3.2

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Mol	Chain	Res	Type	RSRZ
1	D	39	LEU	3.1
1	D	112	ILE	3.1
1	D	119	LEU	2.9
1	B	28	ILE	2.8
1	B	142	ILE	2.8
1	D	3	LEU	2.7
1	A	164	HIS	2.6
1	B	32	SER	2.6
1	A	187	SER	2.5
1	D	116	MET	2.5
1	E	142	ILE	2.5
1	D	164	HIS	2.3
1	B	35	SER	2.3
1	E	145	VAL	2.3
1	B	124	ILE	2.2
1	B	102	PHE	2.2
1	D	36	LYS	2.1
1	D	73	LYS	2.1
1	E	61	TRP	2.1
1	D	115	LYS	2.0
1	A	168	ILE	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 carbohydrates 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. 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	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	SO4	E	899	5/5	0.87	0.54	19.99	182,185,186,186	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	SO4	A	199	5/5	0.91	0.58	12.85	142,143,146,146	0
3	SO4	B	299	5/5	0.80	0.64	9.90	198,199,199,199	0
3	SO4	E	999	5/5	0.95	0.38	9.78	134,135,137,137	0
3	SO4	D	639	5/5	0.64	0.52	6.24	198,199,199,200	0
3	SO4	B	259	5/5	0.80	0.43	3.88	142,144,145,146	0
3	SO4	E	239	5/5	0.84	0.36	3.65	123,125,129,132	0
3	SO4	A	279	5/5	0.88	0.28	2.71	150,152,155,155	0
3	SO4	D	289	5/5	0.82	0.31	0.78	172,173,173,174	0
3	SO4	D	229	5/5	0.64	0.28	0.40	185,186,186,186	0
3	SO4	D	499	5/5	0.85	0.23	0.21	166,166,167,167	0
3	SO4	D	219	5/5	0.84	0.27	0.21	155,155,156,156	0
3	SO4	A	699	5/5	0.93	0.19	-0.61	117,120,126,127	0
2	PNT	A	225	25/25	0.93	0.23	-0.71	58,68,85,98	0
3	SO4	D	799	5/5	0.80	0.18	-1.23	148,149,149,150	0
3	SO4	D	779	5/5	0.84	0.20	-1.25	173,173,174,174	0
3	SO4	A	399	5/5	0.87	0.22	-	140,142,142,144	0
3	SO4	E	209	5/5	0.80	0.15	-	171,171,172,172	0
3	SO4	E	599	5/5	0.86	0.18	-	136,138,139,139	0
3	SO4	D	539	5/5	0.67	0.17	-	186,186,187,187	0
3	SO4	A	579	5/5	0.40	0.59	-	188,188,189,189	0
3	SO4	E	249	5/5	0.58	0.23	-	181,182,182,183	0
3	SO4	B	269	5/5	0.59	0.21	-	179,179,180,181	0

6.5 Other polymers [i](#)

There are no such residues in this entry.