



Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 10:52 pm BST

PDB ID : 3TL6
Title : Crystal structure of purine nucleoside phosphorylase from *Entamoeba histolytica*
Authors : Edwards, T.E.; Clifton, M.C.; Seattle Structural Genomics Center for Infectious Disease (SSGCID)
Deposited on : 2011-08-29
Resolution : 2.65 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

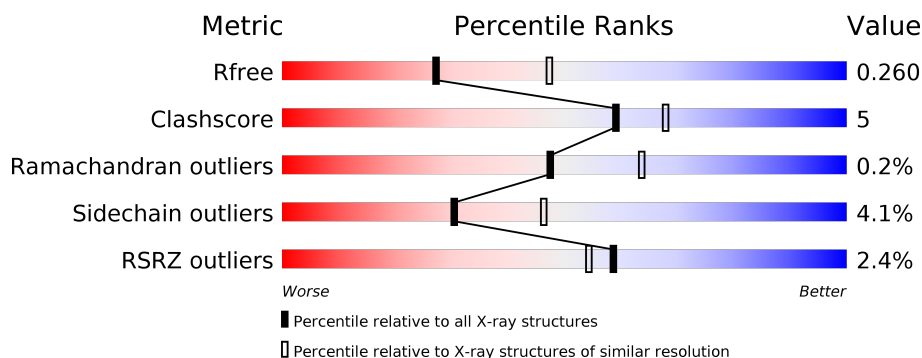
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.65 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}	130704	1332 (2.68-2.64)
Clashscore	141614	1374 (2.68-2.64)
Ramachandran outliers	138981	1349 (2.68-2.64)
Sidechain outliers	138945	1349 (2.68-2.64)
RSRZ outliers	127900	1318 (2.68-2.64)

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 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	242	<div> <div>86%</div> <div>8% • 5%</div> </div>
1	B	242	<div> <div>2%</div> <div>76%</div> <div>14% • 9%</div> </div>
1	C	242	<div> <div>2%</div> <div>81%</div> <div>10% • 8%</div> </div>
1	D	242	<div> <div>87%</div> <div>9% • •</div> </div>
1	E	242	<div> <div>5%</div> <div>82%</div> <div>10% 8%</div> </div>
1	F	242	<div> <div>4%</div> <div>78%</div> <div>13% • 8%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10010 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 Purine nucleoside phosphorylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	231	Total	C	N	O	S	0	0	0
			1715	1095	280	324	16			
1	B	221	Total	C	N	O	S	0	0	0
			1625	1040	264	306	15			
1	C	222	Total	C	N	O	S	0	0	0
			1605	1024	261	305	15			
1	D	237	Total	C	N	O	S	0	0	0
			1735	1104	283	331	17			
1	E	223	Total	C	N	O	S	0	0	0
			1615	1033	265	301	16			
1	F	222	Total	C	N	O	S	0	0	0
			1594	1014	265	300	15			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	EXPRESSION TAG	UNP C4LXG4
A	-2	PRO	-	EXPRESSION TAG	UNP C4LXG4
A	-1	GLY	-	EXPRESSION TAG	UNP C4LXG4
A	0	SER	-	EXPRESSION TAG	UNP C4LXG4
B	-3	GLY	-	EXPRESSION TAG	UNP C4LXG4
B	-2	PRO	-	EXPRESSION TAG	UNP C4LXG4
B	-1	GLY	-	EXPRESSION TAG	UNP C4LXG4
B	0	SER	-	EXPRESSION TAG	UNP C4LXG4
C	-3	GLY	-	EXPRESSION TAG	UNP C4LXG4
C	-2	PRO	-	EXPRESSION TAG	UNP C4LXG4
C	-1	GLY	-	EXPRESSION TAG	UNP C4LXG4
C	0	SER	-	EXPRESSION TAG	UNP C4LXG4
D	-3	GLY	-	EXPRESSION TAG	UNP C4LXG4
D	-2	PRO	-	EXPRESSION TAG	UNP C4LXG4
D	-1	GLY	-	EXPRESSION TAG	UNP C4LXG4
D	0	SER	-	EXPRESSION TAG	UNP C4LXG4
E	-3	GLY	-	EXPRESSION TAG	UNP C4LXG4

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-2	PRO	-	EXPRESSION TAG	UNP C4LXG4
E	-1	GLY	-	EXPRESSION TAG	UNP C4LXG4
E	0	SER	-	EXPRESSION TAG	UNP C4LXG4
F	-3	GLY	-	EXPRESSION TAG	UNP C4LXG4
F	-2	PRO	-	EXPRESSION TAG	UNP C4LXG4
F	-1	GLY	-	EXPRESSION TAG	UNP C4LXG4
F	0	SER	-	EXPRESSION TAG	UNP C4LXG4

- Molecule 2 is SULFATE ION (three-letter code: 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		
2	D	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		

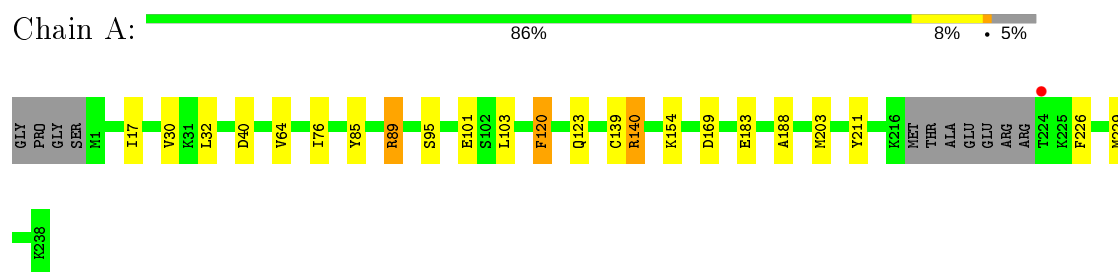
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	19	Total 19	O 19	0	0
3	B	16	Total 16	O 16	0	0
3	C	19	Total 19	O 19	0	0
3	D	18	Total 18	O 18	0	0
3	E	7	Total 7	O 7	0	0
3	F	7	Total 7	O 7	0	0

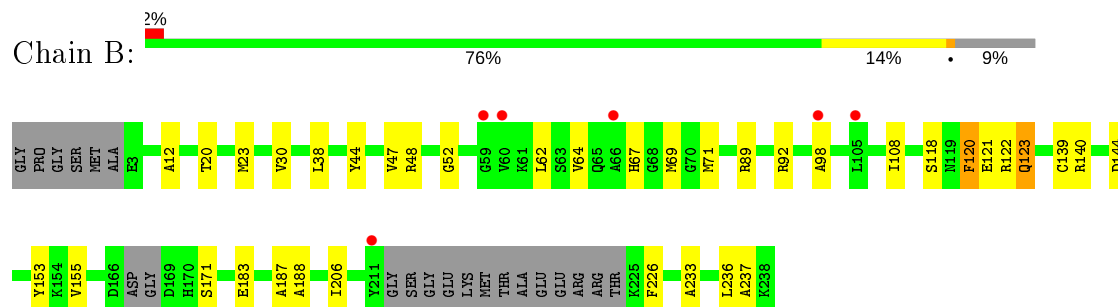
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

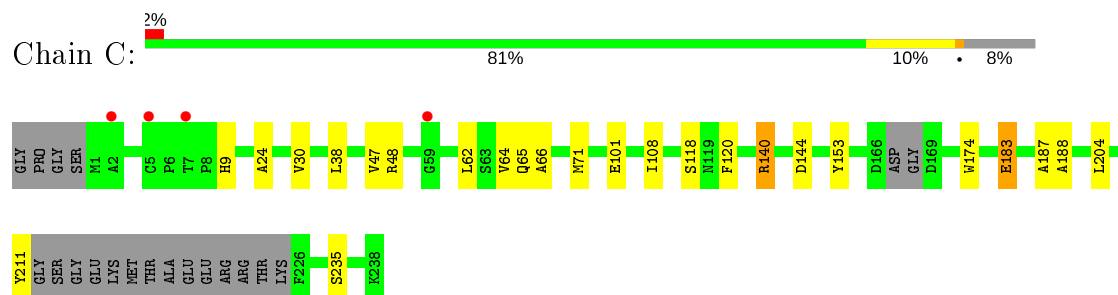
• Molecule 1: Purine nucleoside phosphorylase



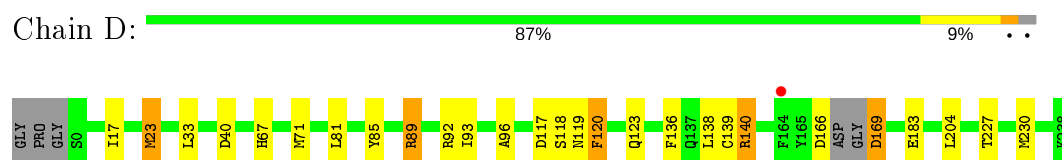
• Molecule 1: Purine nucleoside phosphorylase




• Molecule 1: Purine nucleoside phosphorylase

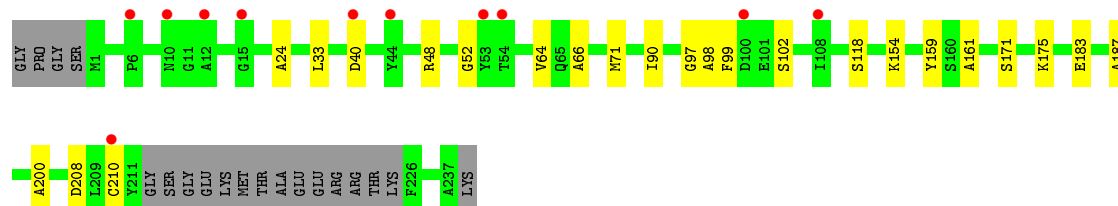


• Molecule 1: Purine nucleoside phosphorylase




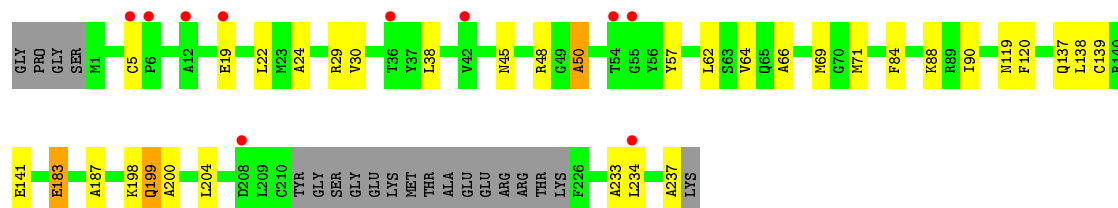
- Molecule 1: Purine nucleoside phosphorylase

Chain E: 



- Molecule 1: Purine nucleoside phosphorylase

Chain F: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	97.97Å 86.56Å 102.56Å 90.00° 117.88° 90.00°	Depositor
Resolution (Å)	50.00 – 2.65 45.33 – 2.65	Depositor EDS
% Data completeness (in resolution range)	99.0 (50.00-2.65) 99.0 (45.33-2.65)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.00 (at 2.65Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.213 , 0.257 0.216 , 0.260	Depositor DCC
R_{free} test set	2200 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	55.5	Xtriage
Anisotropy	0.034	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 47.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.018 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	10010	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.85% 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: SO4

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	0.48	0/1750	0.64	0/2370
1	B	0.47	0/1659	0.60	0/2253
1	C	0.46	1/1637 (0.1%)	0.58	0/2226
1	D	0.48	0/1768	0.61	0/2397
1	E	0.46	0/1648	0.57	0/2237
1	F	0.46	0/1627	0.59	0/2213
All	All	0.47	1/10089 (0.0%)	0.60	0/13696

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	174	TRP	CD2-CE2	5.02	1.47	1.41

There are no bond angle outliers.

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	1715	0	1620	13	0
1	B	1625	0	1492	21	0
1	C	1605	0	1461	13	0
1	D	1735	0	1621	16	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	1615	0	1498	11	0
1	F	1594	0	1444	23	0
2	A	5	0	0	1	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
2	D	10	0	0	0	0
2	E	5	0	0	0	0
2	F	5	0	0	0	0
3	A	19	0	0	0	0
3	B	16	0	0	0	0
3	C	19	0	0	0	0
3	D	18	0	0	1	0
3	E	7	0	0	0	0
3	F	7	0	0	0	0
All	All	10010	0	9136	89	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 5.

All (89) 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:71:MET:HE3	1:B:187:ALA:HB3	1.54	0.89
1:C:30:VAL:HG13	1:C:64:VAL:HG22	1.67	0.76
1:F:119:ASN:ND2	1:F:120:PHE:H	1.84	0.75
1:D:120:PHE:HB2	1:D:123:GLN:HG2	1.73	0.70
1:C:71:MET:HE3	1:C:187:ALA:HB3	1.75	0.67
1:A:140:ARG:NH2	1:B:144:ASP:OD2	2.30	0.65
1:B:23:MET:HB3	1:B:67:HIS:HD2	1.63	0.64
1:A:103:LEU:HD21	1:A:154:LYS:HG3	1.81	0.62
1:B:120:PHE:CZ	1:B:188:ALA:HB1	2.35	0.62
1:F:71:MET:HE3	1:F:187:ALA:HB3	1.81	0.61
1:D:93:ILE:O	1:D:93:ILE:HG13	2.02	0.60
1:B:120:PHE:HZ	1:B:188:ALA:HB1	1.67	0.59
1:E:99:PHE:HB3	1:E:175:LYS:HG2	1.86	0.58
1:F:24:ALA:HB1	1:F:29:ARG:NH2	2.20	0.56
1:F:137:GLN:O	1:F:141:GLU:HG3	2.06	0.55
1:E:171:SER:O	1:E:175:LYS:HG3	2.07	0.55
1:F:30:VAL:HG13	1:F:64:VAL:HG22	1.88	0.54
1:B:47:VAL:HG12	1:B:48:ARG:HG2	1.89	0.54
1:A:17:ILE:HD12	1:A:85:TYR:CD1	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:95:SER:HB3	1:A:226:PHE:HZ	1.71	0.54
1:E:102:SER:OG	1:E:154:LYS:NZ	2.34	0.54
1:B:98:ALA:HB2	1:B:206:ILE:HD12	1.89	0.54
2:A:239:SO4:O1	1:F:48:ARG:NH1	2.40	0.53
1:A:120:PHE:CZ	1:A:188:ALA:HB1	2.43	0.53
1:E:97:GLY:HA2	1:E:208:ASP:O	2.09	0.53
1:E:71:MET:HE3	1:E:187:ALA:HB3	1.90	0.52
1:D:17:ILE:HD12	1:D:85:TYR:CD1	2.44	0.52
1:F:45:ASN:ND2	1:F:50:ALA:O	2.43	0.52
1:C:47:VAL:HG12	1:C:48:ARG:HG2	1.91	0.51
1:F:19:GLU:HB3	1:F:88:LYS:HG2	1.93	0.51
1:B:38:LEU:HD21	1:B:62:LEU:HD22	1.93	0.51
1:F:90:ILE:O	1:F:200:ALA:HA	2.11	0.51
1:B:30:VAL:HG13	1:B:64:VAL:HG22	1.93	0.50
1:A:120:PHE:HZ	1:A:188:ALA:HB1	1.77	0.50
1:F:24:ALA:HB1	1:F:29:ARG:HH21	1.77	0.50
1:B:44:TYR:CZ	1:B:52:GLY:HA3	2.47	0.49
1:B:118:SER:HB3	1:C:118:SER:HB3	1.93	0.49
1:F:88:LYS:HA	1:F:198:LYS:HD2	1.95	0.49
1:F:24:ALA:O	1:F:66:ALA:HA	2.12	0.49
1:C:144:ASP:OD2	1:D:140:ARG:NH2	2.46	0.49
1:C:24:ALA:O	1:C:66:ALA:HA	2.13	0.49
1:B:120:PHE:O	1:B:123:GLN:HG2	2.12	0.48
1:A:203:MET:HE3	1:A:229:MET:HG3	1.95	0.48
1:B:120:PHE:HB2	1:B:123:GLN:HG2	1.96	0.47
1:C:120:PHE:CZ	1:C:188:ALA:HB1	2.49	0.47
1:C:101:GLU:HA	1:C:211:TYR:CZ	2.50	0.47
1:B:108:ILE:O	1:B:153:TYR:HA	2.15	0.47
1:B:62:LEU:HD12	1:B:237:ALA:HB2	1.96	0.46
1:B:69:MET:CE	1:C:9:HIS:CD2	2.98	0.46
1:F:119:ASN:HD22	1:F:120:PHE:H	1.59	0.46
1:F:138:LEU:HD11	1:F:199:GLN:HB3	1.98	0.46
1:F:38:LEU:HD11	1:F:62:LEU:HD22	1.97	0.46
1:D:92:ARG:HG3	1:D:92:ARG:HH11	1.82	0.45
1:B:12:ALA:HB2	1:B:44:TYR:HB2	1.99	0.45
1:F:5:CYS:SG	1:F:84:PHE:HB3	2.57	0.45
1:C:183:GLU:HG2	1:C:204:LEU:HD21	1.99	0.45
1:D:23:MET:HG2	1:D:67:HIS:HB3	2.00	0.43
1:E:24:ALA:O	1:E:66:ALA:HA	2.18	0.43
1:A:95:SER:HB3	1:A:226:PHE:CZ	2.52	0.43
1:D:89:ARG:HG2	1:D:138:LEU:HD21	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:140:ARG:HA	1:D:136:PHE:HZ	1.84	0.42
1:B:121:GLU:HG2	1:B:122:ARG:N	2.35	0.42
1:D:118:SER:HB3	1:E:118:SER:HB3	2.02	0.42
1:D:23:MET:CE	1:D:81:LEU:HD11	2.50	0.42
1:E:98:ALA:O	1:E:210:CYS:HB2	2.19	0.42
1:D:71:MET:SD	1:D:119:ASN:ND2	2.92	0.42
1:E:90:ILE:O	1:E:200:ALA:HA	2.19	0.42
1:A:76:ILE:HD13	1:F:69:MET:HE1	2.01	0.42
1:B:233:ALA:O	1:B:236:LEU:HB2	2.20	0.42
1:C:38:LEU:HD21	1:C:62:LEU:HD22	2.02	0.42
1:D:117:ASP:HA	3:D:241:HOH:O	2.20	0.42
1:A:89:ARG:HB2	1:A:89:ARG:HH11	1.84	0.42
1:A:101:GLU:HA	1:A:211:TYR:CZ	2.54	0.42
1:D:96:ALA:HB2	1:D:204:LEU:HD13	2.02	0.41
1:F:57:TYR:CE1	1:F:234:LEU:HB3	2.55	0.41
1:C:108:ILE:O	1:C:153:TYR:HA	2.20	0.41
1:F:22:LEU:HD11	1:F:233:ALA:HB1	2.01	0.41
1:D:23:MET:HE3	1:D:81:LEU:HD11	2.03	0.41
1:E:52:GLY:HA2	1:E:64:VAL:O	2.21	0.41
1:F:119:ASN:CG	1:F:120:PHE:H	2.18	0.41
1:A:30:VAL:HG13	1:A:64:VAL:HG22	2.03	0.41
1:B:20:THR:HG23	1:B:89:ARG:HB3	2.02	0.41
1:D:166:ASP:OD2	1:D:169:ASP:N	2.54	0.40
1:F:183:GLU:HG2	1:F:204:LEU:HD11	2.02	0.40
1:F:62:LEU:HD12	1:F:237:ALA:CB	2.51	0.40
1:D:33:LEU:HD12	1:D:230:MET:HG2	2.03	0.40
1:A:76:ILE:HD13	1:F:69:MET:CE	2.51	0.40
1:B:20:THR:HG21	1:B:237:ALA:HA	2.04	0.40
1:E:159:TYR:CE2	1:E:161:ALA:HB2	2.57	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	227/242 (94%)	220 (97%)	7 (3%)	0	100	100
1	B	215/242 (89%)	205 (95%)	9 (4%)	1 (0%)	29	43
1	C	216/242 (89%)	205 (95%)	11 (5%)	0	100	100
1	D	233/242 (96%)	223 (96%)	10 (4%)	0	100	100
1	E	219/242 (90%)	214 (98%)	5 (2%)	0	100	100
1	F	218/242 (90%)	210 (96%)	7 (3%)	1 (0%)	29	43
All	All	1328/1452 (92%)	1277 (96%)	49 (4%)	2 (0%)	47	64

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	171	SER
1	F	50	ALA

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	164/188 (87%)	155 (94%)	9 (6%)	21	33
1	B	150/188 (80%)	142 (95%)	8 (5%)	22	35
1	C	145/188 (77%)	141 (97%)	4 (3%)	43	61
1	D	163/188 (87%)	154 (94%)	9 (6%)	21	33
1	E	147/188 (78%)	143 (97%)	4 (3%)	44	63
1	F	142/188 (76%)	139 (98%)	3 (2%)	53	72
All	All	911/1128 (81%)	874 (96%)	37 (4%)	30	46

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	LEU
1	A	40	ASP
1	A	89	ARG

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Mol	Chain	Res	Type
1	A	120	PHE
1	A	123	GLN
1	A	139	CYS
1	A	140	ARG
1	A	169	ASP
1	A	183	GLU
1	B	92	ARG
1	B	120	PHE
1	B	123	GLN
1	B	139	CYS
1	B	140	ARG
1	B	155	VAL
1	B	183	GLU
1	B	226	PHE
1	C	65	GLN
1	C	140	ARG
1	C	183	GLU
1	C	235	SER
1	D	23	MET
1	D	40	ASP
1	D	89	ARG
1	D	120	PHE
1	D	139	CYS
1	D	140	ARG
1	D	169	ASP
1	D	183	GLU
1	D	227	THR
1	E	33	LEU
1	E	40	ASP
1	E	48	ARG
1	E	183	GLU
1	F	139	CYS
1	F	183	GLU
1	F	199	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	9	HIS

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](#)

7 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	D	239	-	4,4,4	0.55	0	6,6,6	0.32	0
2	SO4	D	240	-	4,4,4	0.44	0	6,6,6	0.35	0
2	SO4	B	239	-	4,4,4	0.42	0	6,6,6	0.14	0
2	SO4	A	239	-	4,4,4	0.44	0	6,6,6	0.30	0
2	SO4	C	239	-	4,4,4	0.38	0	6,6,6	0.08	0
2	SO4	E	239	-	4,4,4	0.44	0	6,6,6	0.19	0
2	SO4	F	239	-	4,4,4	0.38	0	6,6,6	0.15	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	239	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	231/242 (95%)	-0.10	1 (0%) 92 93	32, 48, 67, 92	0
1	B	221/242 (91%)	0.04	6 (2%) 54 50	32, 60, 87, 100	0
1	C	222/242 (91%)	0.07	4 (1%) 68 65	33, 59, 93, 103	0
1	D	237/242 (97%)	-0.10	1 (0%) 92 93	32, 47, 79, 110	0
1	E	223/242 (92%)	0.32	11 (4%) 29 26	34, 67, 96, 131	0
1	F	222/242 (91%)	0.34	10 (4%) 33 30	37, 67, 97, 107	0
All	All	1356/1452 (93%)	0.09	33 (2%) 59 54	32, 56, 91, 131	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	60	VAL	3.8
1	F	12	ALA	3.8
1	F	208	ASP	3.5
1	E	40	ASP	3.5
1	C	59	GLY	3.5
1	E	100	ASP	3.2
1	C	2	ALA	3.1
1	B	59	GLY	3.1
1	F	54	THR	3.0
1	E	53	TYR	2.9
1	F	36	THR	2.8
1	E	44	TYR	2.6
1	F	6	PRO	2.6
1	F	42	VAL	2.6
1	E	12	ALA	2.5
1	B	105	LEU	2.5
1	B	211	TYR	2.4
1	B	98	ALA	2.4
1	E	210	CYS	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	7	THR	2.4
1	D	164	PHE	2.3
1	E	6	PRO	2.3
1	E	54	THR	2.3
1	B	66	ALA	2.3
1	E	108	ILE	2.3
1	E	15	GLY	2.2
1	A	224	THR	2.2
1	F	19	GLU	2.2
1	F	55	GLY	2.1
1	F	234	LEU	2.1
1	F	5	CYS	2.1
1	C	5	CYS	2.1
1	E	10	ASN	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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	E	239	5/5	0.83	0.15	84,86,95,99	0
2	SO4	B	239	5/5	0.88	0.14	85,89,97,98	0
2	SO4	F	239	5/5	0.89	0.18	88,92,96,97	0
2	SO4	D	240	5/5	0.90	0.37	71,73,77,78	0
2	SO4	A	239	5/5	0.93	0.11	56,57,59,62	0
2	SO4	C	239	5/5	0.93	0.12	90,91,99,99	0
2	SO4	D	239	5/5	0.95	0.09	55,61,63,68	0

6.5 Other polymers [i](#)

There are no such residues in this entry.