Full wwPDB X-ray Structure Validation Report (i)

Dec 6, 2023 – 01:37 am GMT

PDB ID	:	2WFM
Title	:	Crystal structure of polyneuridine aldehyde esterase mutant (H244A)
Authors	:	Yang, L.; Hill, M.; Panjikar, S.; Wang, M.; Stoeckigt, J.
Deposited on	:	2009-04-08
Resolution	:	2.20 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 (i) 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 (1)) were used in the production of this report:

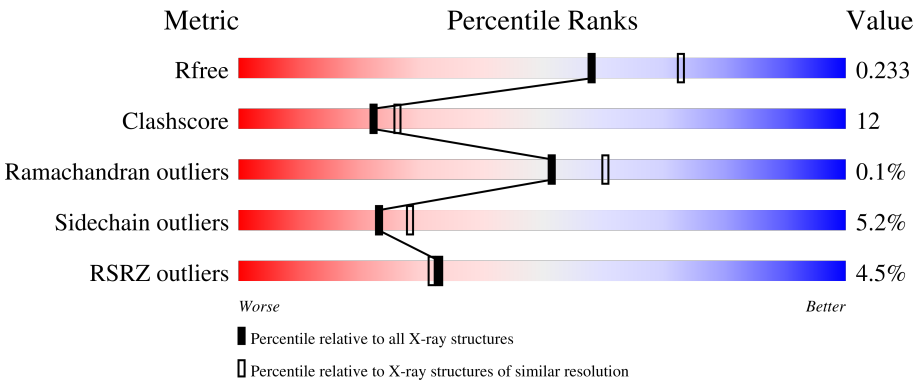
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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 <=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	264	3% 76% 17%
1	B	264	2% 82% 12%
1	C	264	2% 80% 14%
1	D	264	3% 78% 14%
1	E	264	12% 63% 30%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 10221 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 POLYNEURIDINE ALDEHYDE ESTERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	255	2006	1287	326	371	22	0	0	0
1	B	255	2009	1289	326	372	22	0	1	0
1	C	255	2006	1287	326	371	22	0	0	0
1	D	255	2006	1287	326	371	22	0	0	0
1	E	255	2006	1287	326	371	22	0	0	0

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	244	ALA	HIS	engineered mutation	UNP Q9SE93
B	244	ALA	HIS	engineered mutation	UNP Q9SE93
C	244	ALA	HIS	engineered mutation	UNP Q9SE93
D	244	ALA	HIS	engineered mutation	UNP Q9SE93
E	244	ALA	HIS	engineered mutation	UNP Q9SE93

- Molecule 2 is water.

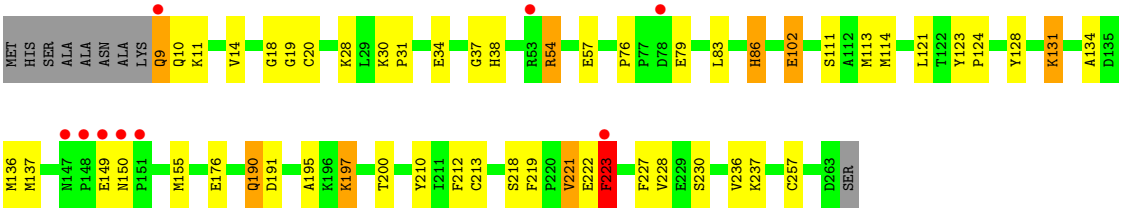
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	47	47	47	0	0
2	B	41	41	41	0	0
2	C	41	41	41	0	0
2	D	44	44	44	0	0
2	E	15	15	15	0	0

3 Residue-property plots (i)

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.

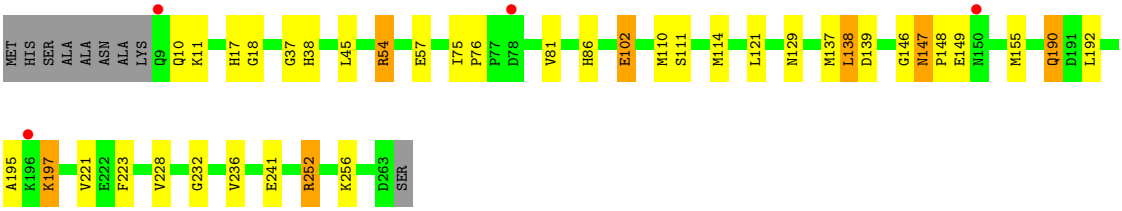
• Molecule 1: POLYNEURIDINE ALDEHYDE ESTERASE

Chain A:



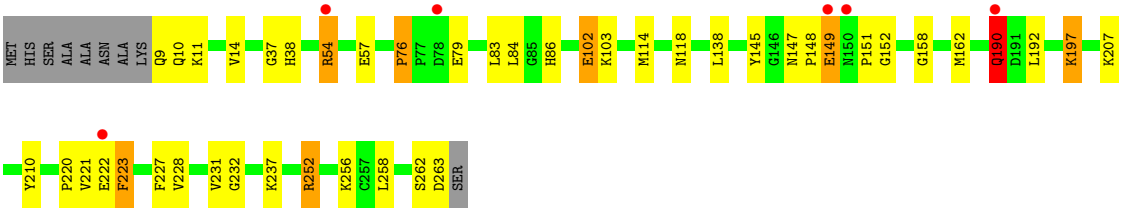
• Molecule 1: POLYNEURIDINE ALDEHYDE ESTERASE

Chain B:



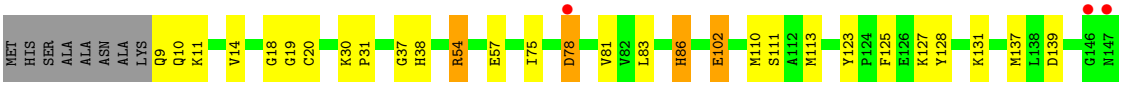
• Molecule 1: POLYNEURIDINE ALDEHYDE ESTERASE

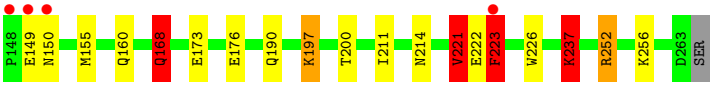
Chain C:



• Molecule 1: POLYNEURIDINE ALDEHYDE ESTERASE

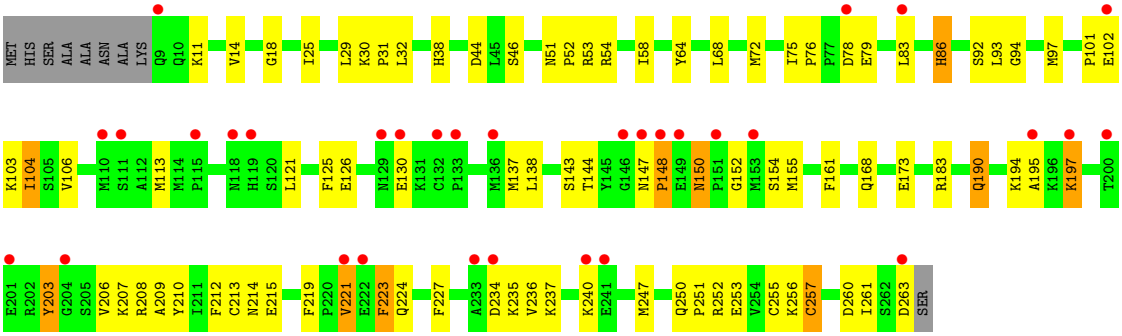
Chain D:





- Molecule 1: POLYNEURIDINE ALDEHYDE ESTERASE

Chain E:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	170.25Å 46.81Å 184.17Å 90.00° 104.89° 90.00°	Depositor
Resolution (Å)	19.94 – 2.20 19.93 – 2.20	Depositor EDS
% Data completeness (in resolution range)	96.8 (19.94-2.20) 96.8 (19.93-2.20)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.13 (at 2.19Å)	Xtriage
Refinement program	REFMAC 5.4.0069	Depositor
R, R_{free}	0.198 , 0.236 0.192 , 0.233	Depositor DCC
R_{free} test set	613 reflections (0.88%)	wwPDB-VP
Wilson B-factor (Å^2)	40.9	Xtriage
Anisotropy	0.061	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.39 , 45.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.45$, $< L^2 > = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	10221	wwPDB-VP
Average B, all atoms (Å^2)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.11% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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)

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	0.86	4/2056 (0.2%)	0.86	5/2772 (0.2%)
1	B	0.92	4/2062 (0.2%)	0.85	5/2780 (0.2%)
1	C	0.83	4/2056 (0.2%)	0.86	8/2772 (0.3%)
1	D	0.86	5/2056 (0.2%)	1.08	10/2772 (0.4%)
1	E	0.62	1/2056 (0.0%)	0.71	1/2772 (0.0%)
All	All	0.82	18/10286 (0.2%)	0.88	29/13868 (0.2%)

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	223	PHE	CB-CG	-12.36	1.30	1.51
1	D	223	PHE	CB-CG	-11.28	1.32	1.51
1	B	223	PHE	CB-CG	-10.99	1.32	1.51
1	A	223	PHE	CB-CG	-10.55	1.33	1.51
1	A	222	GLU	CB-CG	-9.37	1.34	1.52
1	B	223	PHE	CD2-CE2	-7.93	1.23	1.39
1	C	223	PHE	CD2-CE2	-7.34	1.24	1.39
1	B	223	PHE	CE2-CZ	-6.79	1.24	1.37
1	A	223	PHE	CD2-CE2	-6.77	1.25	1.39
1	A	257	CYS	CB-SG	-6.32	1.71	1.82
1	D	168	GLN	CD-OE1	-6.29	1.10	1.24
1	C	223	PHE	CE2-CZ	-6.07	1.25	1.37
1	D	223	PHE	CD2-CE2	-5.91	1.27	1.39
1	D	168	GLN	CD-NE2	-5.75	1.18	1.32
1	C	223	PHE	CD1-CE1	-5.53	1.28	1.39
1	B	241	GLU	CG-CD	-5.13	1.44	1.51
1	D	222	GLU	CG-CD	-5.12	1.44	1.51
1	E	257	CYS	CB-SG	-5.11	1.73	1.81

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	54	ARG	NE-CZ-NH1	-18.15	111.22	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	54	ARG	NE-CZ-NH2	17.04	128.82	120.30
1	D	252	ARG	NE-CZ-NH1	-15.35	112.62	120.30
1	D	252	ARG	NE-CZ-NH2	15.23	127.92	120.30
1	A	223	PHE	CB-CG-CD2	-10.76	113.27	120.80
1	D	223	PHE	CB-CG-CD2	-9.06	114.46	120.80
1	D	54	ARG	CD-NE-CZ	8.15	135.01	123.60
1	D	252	ARG	CD-NE-CZ	7.94	134.72	123.60
1	D	237	LYS	CD-CE-NZ	6.87	127.49	111.70
1	D	221	VAL	CB-CA-C	-6.49	99.06	111.40
1	A	54	ARG	NE-CZ-NH2	-6.45	117.08	120.30
1	A	237	LYS	CD-CE-NZ	6.39	126.39	111.70
1	C	54	ARG	NE-CZ-NH2	-6.37	117.11	120.30
1	B	54	ARG	NE-CZ-NH1	6.23	123.41	120.30
1	A	221	VAL	CB-CA-C	-6.20	99.61	111.40
1	C	252	ARG	NE-CZ-NH2	-5.98	117.31	120.30
1	E	44	ASP	CB-CG-OD1	5.84	123.56	118.30
1	C	237	LYS	CD-CE-NZ	5.82	125.09	111.70
1	C	252	ARG	NE-CZ-NH1	5.80	123.20	120.30
1	C	223	PHE	CB-CG-CD1	-5.66	116.83	120.80
1	C	190	GLN	N-CA-CB	5.52	120.53	110.60
1	B	114	MET	CG-SD-CE	5.48	108.96	100.20
1	D	222	GLU	CG-CD-OE1	-5.40	107.50	118.30
1	B	54	ARG	NE-CZ-NH2	-5.33	117.64	120.30
1	C	54	ARG	NE-CZ-NH1	5.27	122.94	120.30
1	A	54	ARG	NE-CZ-NH1	5.20	122.90	120.30
1	C	223	PHE	CB-CA-C	-5.19	100.02	110.40
1	B	252	ARG	NE-CZ-NH2	-5.07	117.77	120.30
1	B	223	PHE	CB-CG-CD1	-5.00	117.30	120.80

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	2006	0	1985	64	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2009	0	1990	29	0
1	C	2006	0	1985	46	0
1	D	2006	0	1985	38	0
1	E	2006	0	1985	70	0
2	A	47	0	0	4	0
2	B	41	0	0	1	0
2	C	41	0	0	4	0
2	D	44	0	0	3	0
2	E	15	0	0	3	0
All	All	10221	0	9930	233	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 12.

All (233) 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:A:136:MET:HE2	2:A:2020:HOH:O	1.26	1.34
1:A:230:SER:CB	1:C:118:ASN:HD21	1.52	1.22
1:A:137:MET:HE1	1:A:155:MET:CE	1.71	1.20
1:A:190:GLN:H	1:A:190:GLN:NE2	1.42	1.15
1:E:147:ASN:HB2	1:E:148:PRO:HD2	1.18	1.14
1:E:137:MET:HE1	1:E:155:MET:HE1	1.13	1.11
1:A:230:SER:HB3	1:C:118:ASN:HD21	1.09	1.10
1:A:230:SER:HB3	1:C:118:ASN:ND2	1.70	1.05
1:A:230:SER:CB	1:C:118:ASN:ND2	2.19	1.05
1:A:137:MET:CE	1:A:155:MET:HE1	1.86	1.04
1:C:114:MET:CE	1:C:210:TYR:HD1	1.70	1.03
1:E:137:MET:CE	1:E:155:MET:HE1	1.87	1.02
1:A:230:SER:CA	1:C:118:ASN:HD21	1.75	1.00
1:A:114:MET:CE	1:A:210:TYR:HD1	1.76	0.97
1:A:190:GLN:NE2	1:A:190:GLN:N	2.12	0.97
1:C:11:LYS:H	1:C:38:HIS:HD2	1.06	0.95
1:D:11:LYS:H	1:D:38:HIS:HD2	1.12	0.94
1:B:11:LYS:H	1:B:38:HIS:HD2	1.09	0.94
1:D:190:GLN:H	1:D:190:GLN:NE2	1.66	0.94
1:E:147:ASN:CB	1:E:148:PRO:HD2	1.95	0.94
1:A:11:LYS:H	1:A:38:HIS:HD2	1.13	0.92
1:A:137:MET:HE1	1:A:155:MET:HE1	0.92	0.92
1:B:190:GLN:H	1:B:190:GLN:NE2	1.67	0.91
1:A:137:MET:CE	1:A:155:MET:CE	2.48	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:137:MET:HE3	1:D:155:MET:HE1	1.55	0.88
1:A:230:SER:HA	1:C:118:ASN:ND2	1.90	0.87
1:D:197:LYS:NZ	2:D:2039:HOH:O	2.02	0.85
1:D:137:MET:CE	1:D:155:MET:HE1	2.07	0.83
1:E:147:ASN:HB2	1:E:148:PRO:CD	2.05	0.82
1:C:114:MET:CE	1:C:210:TYR:CD1	2.61	0.81
1:D:160:GLN:OE1	2:D:2025:HOH:O	1.99	0.81
1:E:11:LYS:H	1:E:38:HIS:HD2	1.25	0.81
1:C:149:GLU:HA	1:C:149:GLU:OE2	1.79	0.80
1:A:230:SER:CA	1:C:118:ASN:ND2	2.44	0.80
1:C:223:PHE:HE1	2:C:2019:HOH:O	1.66	0.79
1:C:149:GLU:OE2	1:C:149:GLU:CA	2.28	0.79
1:C:114:MET:HE3	1:C:210:TYR:HD1	1.45	0.78
1:E:150:ASN:OD1	1:E:194:LYS:NZ	2.18	0.76
1:A:230:SER:HA	1:C:118:ASN:HD21	1.48	0.76
1:C:11:LYS:H	1:C:38:HIS:CD2	1.98	0.76
1:A:11:LYS:H	1:A:38:HIS:CD2	2.01	0.76
1:B:17:HIS:HE1	1:B:45:LEU:H	1.31	0.75
1:B:11:LYS:H	1:B:38:HIS:CD2	1.99	0.75
1:A:136:MET:CE	2:A:2020:HOH:O	2.01	0.75
1:C:147:ASN:HD21	1:C:149:GLU:HG2	1.52	0.74
1:A:114:MET:HE2	1:A:210:TYR:HD1	1.51	0.74
1:C:114:MET:HE3	1:C:210:TYR:CD1	2.22	0.74
1:E:75:ILE:HG23	1:E:76:PRO:HD2	1.69	0.73
1:D:11:LYS:H	1:D:38:HIS:CD2	2.03	0.73
1:E:197:LYS:H	1:E:197:LYS:HD2	1.54	0.72
1:A:114:MET:CE	1:A:210:TYR:CD1	2.68	0.72
1:E:75:ILE:CG2	1:E:76:PRO:HD2	2.20	0.71
1:A:190:GLN:H	1:A:190:GLN:CD	1.92	0.71
1:B:17:HIS:HD2	1:B:18:GLY:O	1.74	0.71
1:A:128:TYR:OH	1:A:218:SER:O	2.07	0.69
1:C:262:SER:O	1:C:263:ASP:C	2.32	0.68
1:E:252:ARG:O	1:E:256:LYS:HG2	1.94	0.68
1:A:114:MET:HE3	1:A:210:TYR:HD1	1.58	0.68
1:C:114:MET:HE2	1:C:210:TYR:HD1	1.55	0.68
1:E:137:MET:HE1	1:E:155:MET:CE	2.08	0.67
1:C:11:LYS:N	1:C:38:HIS:HD2	1.88	0.66
1:C:145:TYR:O	1:C:152:GLY:N	2.24	0.66
1:D:137:MET:CE	1:D:155:MET:CE	2.73	0.66
1:D:190:GLN:H	1:D:190:GLN:CD	1.98	0.65
1:E:257:CYS:O	1:E:260:ASP:HB2	1.96	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:9:GLN:OE1	1:A:9:GLN:HA	1.97	0.65
1:B:54:ARG:O	1:B:57:GLU:HG2	1.97	0.65
1:A:18:GLY:H	1:A:86:HIS:CE1	2.15	0.64
1:C:147:ASN:HD22	1:C:148:PRO:HD2	1.63	0.64
1:E:213:CYS:H	1:E:224:GLN:HE22	1.43	0.64
1:C:102:GLU:CD	1:C:102:GLU:H	2.01	0.64
1:E:54:ARG:HH11	1:E:54:ARG:HG2	1.61	0.64
1:A:128:TYR:HA	1:A:223:PHE:HZ	1.61	0.64
1:A:128:TYR:HA	1:A:223:PHE:CZ	2.33	0.64
1:A:114:MET:HE2	1:A:210:TYR:CD1	2.32	0.64
1:E:137:MET:CE	1:E:155:MET:CE	2.71	0.64
1:D:128:TYR:HA	1:D:223:PHE:CZ	2.34	0.63
1:E:121:LEU:HD12	1:E:195:ALA:HB3	1.80	0.63
1:E:46:SER:HB3	1:E:58:ILE:CD1	2.29	0.63
1:B:197:LYS:HB2	2:B:2033:HOH:O	1.99	0.62
1:E:72:MET:HA	1:E:75:ILE:HD12	1.81	0.62
1:E:76:PRO:HG2	1:E:79:GLU:HB2	1.81	0.61
1:A:200:THR:OG1	1:C:207:LYS:HG3	1.99	0.61
1:E:102:GLU:H	1:E:102:GLU:CD	2.03	0.60
1:C:54:ARG:O	1:C:57:GLU:HG2	2.01	0.60
1:E:30:LYS:HB3	1:E:31:PRO:HD3	1.83	0.60
1:E:214:ASN:HA	1:E:221:VAL:CG2	2.31	0.60
1:A:131:LYS:HB2	1:A:223:PHE:HE2	1.66	0.59
1:C:114:MET:HE2	1:C:210:TYR:CD1	2.33	0.59
1:E:14:VAL:HB	1:E:83:LEU:HD23	1.84	0.59
1:A:19:GLY:O	1:A:20:CYS:HB2	2.01	0.59
1:E:138:LEU:HD12	1:E:161:PHE:HA	1.85	0.59
1:D:137:MET:HE1	1:D:155:MET:CE	2.32	0.59
1:E:223:PHE:O	1:E:227:PHE:CD1	2.56	0.58
1:C:252:ARG:O	1:C:256:LYS:HG2	2.04	0.58
1:B:137:MET:HE3	1:B:155:MET:HE1	1.84	0.58
1:C:190:GLN:HG2	2:C:2032:HOH:O	2.04	0.58
1:E:214:ASN:HA	1:E:221:VAL:HG21	1.84	0.58
1:C:9:GLN:NE2	2:C:2001:HOH:O	2.36	0.58
1:A:128:TYR:HE2	1:A:136:MET:SD	2.26	0.58
1:E:255:CYS:HB3	1:E:256:LYS:HE2	1.85	0.58
1:D:128:TYR:HA	1:D:223:PHE:HZ	1.68	0.57
1:A:114:MET:HE3	1:A:210:TYR:CD1	2.37	0.56
1:E:76:PRO:CG	1:E:79:GLU:HB2	2.35	0.56
1:E:208:ARG:O	1:E:234:ASP:HB2	2.05	0.56
1:B:190:GLN:H	1:B:190:GLN:CD	2.09	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:235:LYS:O	1:E:236:VAL:HG13	2.05	0.55
1:A:18:GLY:H	1:A:86:HIS:HE1	1.55	0.55
1:E:203:TYR:C	1:E:203:TYR:CD2	2.80	0.55
1:D:113:MET:HE3	1:D:125:PHE:CE1	2.42	0.55
1:A:134:ALA:H	1:B:129:ASN:ND2	2.05	0.54
1:A:14:VAL:HB	1:A:83:LEU:HD23	1.89	0.54
1:A:149:GLU:CD	1:A:149:GLU:H	2.10	0.54
1:E:46:SER:HB3	1:E:58:ILE:HD13	1.88	0.54
1:E:68:LEU:HD12	1:E:68:LEU:O	2.08	0.54
1:C:10:GLN:HG3	1:C:37:GLY:O	2.07	0.54
1:E:30:LYS:HB3	1:E:31:PRO:CD	2.37	0.54
1:A:128:TYR:CE2	1:A:136:MET:SD	3.01	0.53
1:B:147:ASN:C	1:B:147:ASN:HD22	2.13	0.53
1:D:75:ILE:HD13	1:D:81:VAL:HG13	1.91	0.53
1:C:149:GLU:OE2	1:C:149:GLU:N	2.42	0.52
1:B:17:HIS:CE1	1:B:45:LEU:H	2.19	0.52
1:E:18:GLY:H	1:E:86:HIS:HD2	1.58	0.52
1:A:137:MET:CE	1:A:155:MET:HE3	2.39	0.52
1:E:46:SER:HB3	1:E:58:ILE:HD11	1.93	0.51
1:A:176:GLU:HG2	1:D:31:PRO:HD3	1.92	0.51
1:B:102:GLU:H	1:B:102:GLU:CD	2.14	0.51
1:B:228:VAL:O	1:B:232:GLY:HA2	2.11	0.51
1:D:127:LYS:HG3	1:D:226:TRP:CH2	2.45	0.51
1:A:102:GLU:CD	1:A:102:GLU:H	2.14	0.51
1:B:146:GLY:O	1:B:148:PRO:HD3	2.11	0.51
1:E:97:MET:HB3	1:E:206:VAL:HG21	1.93	0.51
1:E:126:GLU:O	1:E:130:GLU:HG3	2.11	0.50
1:A:123:TYR:N	1:A:124:PRO:CD	2.75	0.50
1:E:209:ALA:HB1	1:E:261:ILE:HD13	1.93	0.50
1:D:211:ILE:HA	1:D:237:LYS:O	2.12	0.50
1:A:131:LYS:HB2	1:A:223:PHE:CE2	2.47	0.49
1:A:136:MET:HG3	1:A:137:MET:HG3	1.94	0.49
1:E:214:ASN:HB2	1:E:240:LYS:O	2.12	0.49
1:E:72:MET:HA	1:E:75:ILE:CD1	2.43	0.49
1:B:252:ARG:O	1:B:256:LYS:HG2	2.13	0.49
1:E:210:TYR:HB3	1:E:236:VAL:HG12	1.94	0.49
1:D:102:GLU:CD	1:D:102:GLU:H	2.15	0.48
1:E:252:ARG:HG2	2:E:2015:HOH:O	2.12	0.48
1:D:54:ARG:O	1:D:57:GLU:HG2	2.13	0.48
1:E:18:GLY:N	1:E:86:HIS:HD2	2.09	0.48
1:C:14:VAL:HB	1:C:83:LEU:HD23	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:10:GLN:HG3	1:D:37:GLY:O	2.13	0.48
1:A:11:LYS:N	1:A:38:HIS:HD2	1.96	0.47
1:C:76:PRO:HG2	1:C:79:GLU:CD	2.35	0.47
1:E:104:ILE:HD11	1:E:106:VAL:O	2.14	0.47
1:E:68:LEU:O	1:E:72:MET:HG2	2.14	0.47
1:D:19:GLY:O	1:D:20:CYS:HB2	2.15	0.47
1:D:14:VAL:HB	1:D:83:LEU:HD23	1.97	0.47
1:A:54:ARG:O	1:A:57:GLU:HG2	2.15	0.46
1:A:197:LYS:HB2	2:A:2016:HOH:O	2.14	0.46
1:D:110:MET:O	1:D:111:SER:C	2.52	0.46
1:D:139:ASP:OD2	1:D:160:GLN:HB2	2.15	0.46
1:D:252:ARG:O	1:D:256:LYS:HG2	2.15	0.46
1:B:137:MET:CE	1:B:155:MET:HE1	2.46	0.46
1:D:214:ASN:HA	1:D:221:VAL:HG13	1.98	0.46
1:E:46:SER:HB2	1:E:64:TYR:HA	1.97	0.46
1:B:121:LEU:HD12	1:B:195:ALA:HB3	1.97	0.45
1:D:78:ASP:OD2	1:D:78:ASP:C	2.55	0.45
1:E:29:LEU:O	1:E:32:LEU:HB2	2.16	0.45
1:E:152:GLY:HA2	1:E:190:GLN:NE2	2.32	0.45
1:B:192:LEU:HA	1:B:192:LEU:HD23	1.73	0.45
1:A:190:GLN:HE21	1:A:191:ASP:H	1.64	0.45
1:A:31:PRO:HD3	1:D:176:GLU:HG2	1.99	0.45
1:E:68:LEU:O	1:E:68:LEU:HG	2.15	0.45
1:E:183:ARG:HB2	2:E:2001:HOH:O	2.17	0.45
1:E:207:LYS:HA	1:E:207:LYS:HD3	1.74	0.45
1:A:149:GLU:CD	1:A:149:GLU:N	2.71	0.45
1:D:200:THR:HG21	2:D:2018:HOH:O	2.17	0.45
1:C:84:LEU:HD22	1:C:258:LEU:HD22	1.99	0.45
1:E:68:LEU:O	1:E:68:LEU:CG	2.64	0.45
1:A:10:GLN:HG3	1:A:37:GLY:O	2.16	0.45
1:A:121:LEU:HD12	1:A:195:ALA:HB3	1.98	0.44
1:C:158:GLY:O	1:C:162:MET:HG3	2.17	0.44
1:E:197:LYS:HD2	1:E:197:LYS:N	2.27	0.44
1:C:79:GLU:O	1:C:103:LYS:HE3	2.18	0.44
1:E:93:LEU:O	1:E:94:GLY:C	2.55	0.44
1:A:136:MET:HG3	1:A:137:MET:N	2.32	0.44
1:C:227:PHE:O	1:C:231:VAL:HB	2.17	0.43
1:D:139:ASP:CG	1:D:160:GLN:HB2	2.38	0.43
1:D:18:GLY:H	1:D:86:HIS:CD2	2.36	0.43
1:E:75:ILE:CG2	1:E:103:LYS:HE3	2.48	0.43
1:D:190:GLN:NE2	1:D:190:GLN:N	2.49	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:25:ILE:HD12	1:E:247:MET:CE	2.48	0.43
1:E:78:ASP:HB2	2:E:2007:HOH:O	2.18	0.43
1:B:129:ASN:HD22	1:B:129:ASN:HA	1.55	0.43
1:E:101:PRO:HD2	1:E:102:GLU:OE2	2.19	0.43
1:E:168:GLN:NE2	1:E:215:GLU:O	2.51	0.43
1:B:75:ILE:HA	1:B:76:PRO:HD3	1.85	0.42
1:C:190:GLN:H	1:C:190:GLN:NE2	2.17	0.42
1:E:68:LEU:O	1:E:68:LEU:CD1	2.67	0.42
1:D:168:GLN:HE21	1:D:168:GLN:H	1.67	0.42
1:A:176:GLU:OE2	1:D:30:LYS:HE3	2.19	0.42
1:A:190:GLN:NE2	1:A:190:GLN:CA	2.81	0.42
1:B:75:ILE:HD13	1:B:81:VAL:HG13	2.01	0.42
1:A:111:SER:OG	1:A:213:CYS:HB2	2.20	0.42
1:B:147:ASN:ND2	1:B:149:GLU:OE2	2.44	0.42
1:A:212:PHE:CZ	1:A:228:VAL:HG21	2.54	0.42
1:B:10:GLN:HG3	1:B:37:GLY:O	2.19	0.42
1:B:137:MET:HE2	1:B:137:MET:HB3	1.86	0.42
1:D:131:LYS:O	1:D:131:LYS:HG3	2.20	0.42
1:A:190:GLN:N	1:A:190:GLN:CD	2.60	0.42
1:C:192:LEU:HA	1:C:192:LEU:HD23	1.83	0.41
1:C:147:ASN:O	1:C:151:PRO:N	2.53	0.41
1:D:131:LYS:O	1:D:131:LYS:CG	2.67	0.41
1:E:250:GLN:N	1:E:251:PRO:CD	2.83	0.41
1:C:197:LYS:H	1:C:197:LYS:HG2	1.75	0.41
1:E:212:PHE:HA	1:E:224:GLN:NE2	2.35	0.41
1:C:228:VAL:O	1:C:232:GLY:HA2	2.20	0.41
1:A:30:LYS:NZ	1:A:34:GLU:OE2	2.52	0.41
1:B:147:ASN:HD22	1:B:149:GLU:H	1.67	0.41
1:E:223:PHE:O	1:E:227:PHE:HD1	2.00	0.41
1:A:76:PRO:HG2	1:A:79:GLU:CD	2.41	0.41
1:E:54:ARG:HH11	1:E:54:ARG:CG	2.33	0.41
1:E:113:MET:HG3	1:E:125:PHE:HE1	1.85	0.41
1:D:123:TYR:O	1:D:127:LYS:HG2	2.21	0.41
1:A:200:THR:HG21	2:A:2015:HOH:O	2.20	0.41
1:A:219:PHE:HZ	1:A:227:PHE:CE2	2.39	0.41
1:B:190:GLN:NE2	1:B:190:GLN:N	2.52	0.41
1:C:147:ASN:O	1:C:151:PRO:CA	2.69	0.41
1:B:138:LEU:HB3	1:B:139:ASP:H	1.72	0.40
1:E:75:ILE:HG22	1:E:76:PRO:HD2	2.02	0.40
1:C:190:GLN:NE2	2:C:2032:HOH:O	2.14	0.40
1:C:220:PRO:O	1:C:223:PHE:HB2	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:51:ASN:HA	1:E:52:PRO:HD3	1.76	0.40
1:E:251:PRO:C	1:E:253:GLU:H	2.24	0.40
1:A:28:LYS:HG2	1:D:173:GLU:HG2	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	253/264 (96%)	245 (97%)	8 (3%)	0	100	100
1	B	254/264 (96%)	243 (96%)	11 (4%)	0	100	100
1	C	253/264 (96%)	241 (95%)	12 (5%)	0	100	100
1	D	253/264 (96%)	244 (96%)	9 (4%)	0	100	100
1	E	253/264 (96%)	228 (90%)	24 (10%)	1 (0%)	34	37
All	All	1266/1320 (96%)	1201 (95%)	64 (5%)	1 (0%)	51	60

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	148	PRO

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	221/227 (97%)	210 (95%)	11 (5%)	24	30
1	B	222/227 (98%)	212 (96%)	10 (4%)	27	34
1	C	221/227 (97%)	212 (96%)	9 (4%)	30	39
1	D	221/227 (97%)	210 (95%)	11 (5%)	24	30
1	E	221/227 (97%)	204 (92%)	17 (8%)	13	13
All	All	1106/1135 (97%)	1048 (95%)	58 (5%)	23	28

All (58) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	GLN
1	A	86	HIS
1	A	102	GLU
1	A	113	MET
1	A	131	LYS
1	A	150	ASN
1	A	190	GLN
1	A	197	LYS
1	A	221	VAL
1	A	223	PHE
1	A	236	VAL
1	B	86	HIS
1	B	102	GLU
1	B	111[A]	SER
1	B	111[B]	SER
1	B	138	LEU
1	B	147	ASN
1	B	190	GLN
1	B	197	LYS
1	B	221	VAL
1	B	236	VAL
1	C	76	PRO
1	C	86	HIS
1	C	102	GLU
1	C	138	LEU
1	C	149	GLU
1	C	190	GLN
1	C	197	LYS
1	C	221	VAL
1	C	222	GLU
1	D	9	GLN

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Mol	Chain	Res	Type
1	D	78	ASP
1	D	86	HIS
1	D	102	GLU
1	D	149	GLU
1	D	150	ASN
1	D	168	GLN
1	D	197	LYS
1	D	221	VAL
1	D	223	PHE
1	D	237	LYS
1	E	53	ARG
1	E	86	HIS
1	E	92	SER
1	E	104	ILE
1	E	143	SER
1	E	144	THR
1	E	150	ASN
1	E	154	SER
1	E	173	GLU
1	E	190	GLN
1	E	197	LYS
1	E	203	TYR
1	E	219	PHE
1	E	221	VAL
1	E	223	PHE
1	E	237	LYS
1	E	263	ASP

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	38	HIS
1	A	86	HIS
1	A	129	ASN
1	A	190	GLN
1	B	17	HIS
1	B	38	HIS
1	B	129	ASN
1	B	147	ASN
1	B	190	GLN
1	B	250	GLN
1	C	9	GLN

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Mol	Chain	Res	Type
1	C	38	HIS
1	C	86	HIS
1	C	118	ASN
1	C	129	ASN
1	C	141	GLN
1	C	147	ASN
1	C	190	GLN
1	C	250	GLN
1	D	38	HIS
1	D	150	ASN
1	D	168	GLN
1	D	190	GLN
1	D	250	GLN
1	E	38	HIS
1	E	86	HIS
1	E	224	GLN

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	255/264 (96%)	-0.06	9 (3%)	44	42	19, 23, 30, 38	0
1	B	255/264 (96%)	-0.16	4 (1%)	72	70	18, 23, 29, 37	0
1	C	255/264 (96%)	-0.11	6 (2%)	59	56	19, 23, 30, 34	0
1	D	255/264 (96%)	-0.13	7 (2%)	54	52	17, 23, 31, 36	0
1	E	255/264 (96%)	0.57	32 (12%)	3	3	20, 28, 35, 38	0
All	All	1275/1320 (96%)	0.02	58 (4%)	33	32	17, 24, 32, 38	0

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	149	GLU	6.0
1	E	78	ASP	5.6
1	D	150	ASN	5.1
1	D	78	ASP	5.0
1	E	148	PRO	4.7
1	B	9	GLN	4.6
1	E	119	HIS	4.4
1	B	150	ASN	4.3
1	E	132	CYS	4.0
1	E	233	ALA	3.9
1	B	78	ASP	3.9
1	E	83	LEU	3.8
1	E	151	PRO	3.7
1	C	149	GLU	3.6
1	E	9	GLN	3.6
1	A	150	ASN	3.5
1	E	130	GLU	3.4
1	A	78	ASP	3.4
1	E	115	PRO	3.3
1	D	223	PHE	3.3

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Mol	Chain	Res	Type	RSRZ
1	D	149	GLU	3.3
1	A	149	GLU	3.1
1	A	223	PHE	3.1
1	E	204	GLY	3.1
1	E	263	ASP	3.1
1	C	78	ASP	3.0
1	E	133	PRO	2.9
1	E	147	ASN	2.8
1	C	190	GLN	2.8
1	E	102	GLU	2.8
1	A	147	ASN	2.8
1	C	222	GLU	2.8
1	E	241	GLU	2.8
1	E	118	ASN	2.8
1	E	221	VAL	2.6
1	B	196	LYS	2.6
1	E	234	ASP	2.6
1	A	148	PRO	2.6
1	E	240	LYS	2.6
1	A	9	GLN	2.6
1	E	136	MET	2.6
1	E	201	GLU	2.5
1	D	148	PRO	2.5
1	D	146	GLY	2.5
1	E	146	GLY	2.5
1	A	53	ARG	2.4
1	C	150	ASN	2.4
1	E	200	THR	2.4
1	D	147	ASN	2.4
1	E	195	ALA	2.3
1	C	54	ARG	2.3
1	E	111	SER	2.3
1	E	222	GLU	2.3
1	E	110	MET	2.3
1	E	197	LYS	2.2
1	A	151	PRO	2.2
1	E	129	ASN	2.1
1	E	153	MET	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 monosaccharides in this entry.

6.4 Ligands (i)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.