



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 10:08 AM UTC

PDB ID : 6G2U / pdb_00006g2u
Title : Crystal structure of the human glutamate dehydrogenase 2 (hGDH2)
Authors : Fadoulglou, V.F.; Dimovasili, C.; Providaki, M.; Kotsifaki, D.; Sarrou, I.;
Plaitakis, A.; Zaganas, I.; Kokkinidis, M.
Deposited on : 2018-03-23
Resolution : 2.93 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

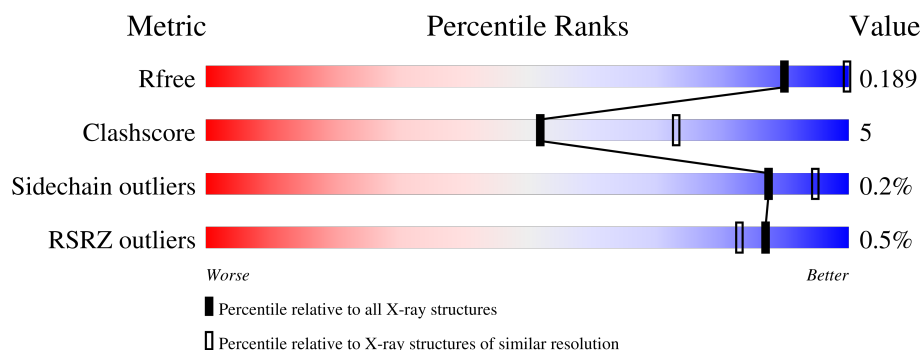
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.93 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	1159 (2.96-2.92)
Clashscore	190562	1184 (2.96-2.92)
Sidechain outliers	187428	1131 (2.96-2.92)
RSRZ outliers	180081	1159 (2.96-2.92)

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	505	 84% 14% .
1	B	505	 84% 13% .
1	C	505	 83% 14% .
1	D	505	 88% 10% .
1	E	505	 86% 11% .
1	F	505	 86% 12% .

2 Entry composition [i](#)

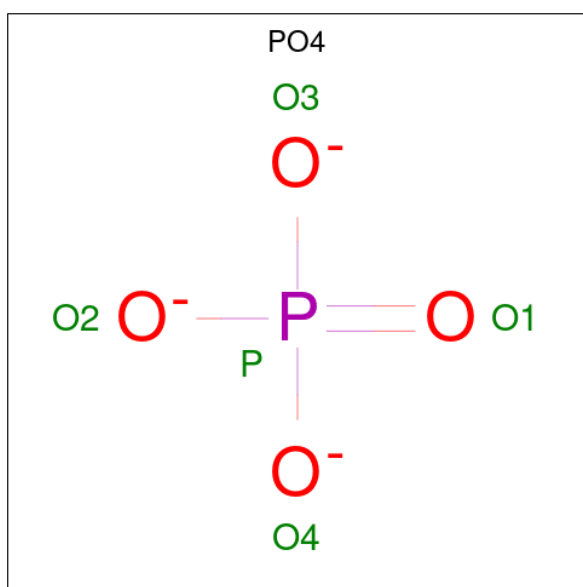
There are 5 unique types of molecules in this entry. The entry contains 24160 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Glutamate dehydrogenase 2, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	494	Total	C	N	O	S	0	0	0
			3809	2422	664	706	17			
1	B	491	Total	C	N	O	S	0	0	0
			3605	2267	643	679	16			
1	C	491	Total	C	N	O	S	0	0	0
			3532	2206	641	669	16			
1	D	496	Total	C	N	O	S	0	0	0
			3821	2431	659	715	16			
1	E	492	Total	C	N	O	S	0	0	0
			3805	2420	664	704	17			
1	F	496	Total	C	N	O	S	0	0	0
			3809	2431	659	702	17			

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total O P 5 4 1	0	0
2	A	1	Total O P 5 4 1	0	0
2	A	1	Total O P 5 4 1	0	0
2	A	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0
2	C	1	Total O P 5 4 1	0	0
2	C	1	Total O P 5 4 1	0	0
2	C	1	Total O P 5 4 1	0	0
2	C	1	Total O P 5 4 1	0	0
2	C	1	Total O P 5 4 1	0	0
2	D	1	Total O P 5 4 1	0	0
2	D	1	Total O P 5 4 1	0	0
2	E	1	Total O P 5 4 1	0	0
2	E	1	Total O P 5 4 1	0	0
2	F	1	Total O P 5 4 1	0	0
2	F	1	Total O P 5 4 1	0	0
2	F	1	Total O P 5 4 1	0	0
2	F	1	Total O P 5 4 1	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	F	1	Total	O	P	0	0
			5	4	1		

- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Na	0	0
			2	2		
3	C	2	Total	Na	0	0
			2	2		
3	D	11	Total	Na	0	0
			11	11		
3	E	3	Total	Na	0	0
			3	3		
3	F	3	Total	Na	0	0
			3	3		

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	3	Total	Cl	0	0
			3	3		
4	B	2	Total	Cl	0	0
			2	2		
4	C	2	Total	Cl	0	0
			2	2		
4	D	5	Total	Cl	0	0
			5	5		
4	E	6	Total	Cl	0	0
			6	6		
4	F	1	Total	Cl	0	0
			1	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	296	Total	O	0	0
			296	296		
5	B	197	Total	O	0	0
			197	197		
5	C	235	Total	O	0	0
			235	235		

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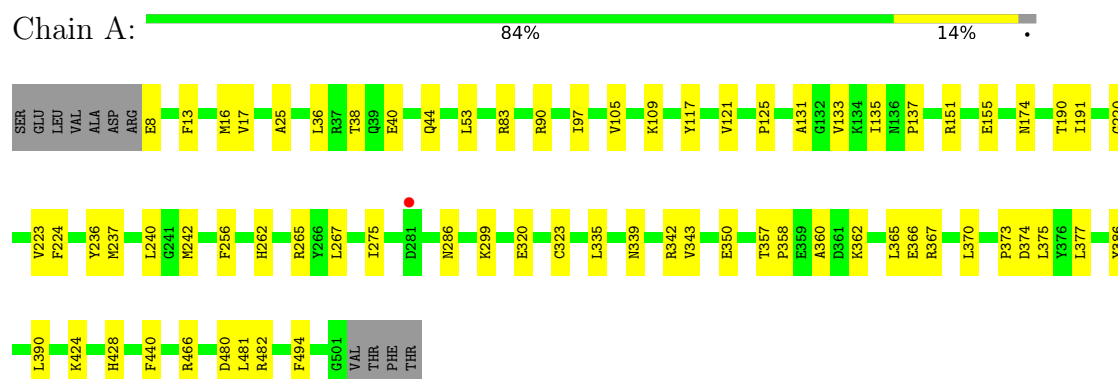
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	D	365	Total 365	O 365	0	0
5	E	296	Total 296	O 296	0	0
5	F	235	Total 235	O 235	0	0

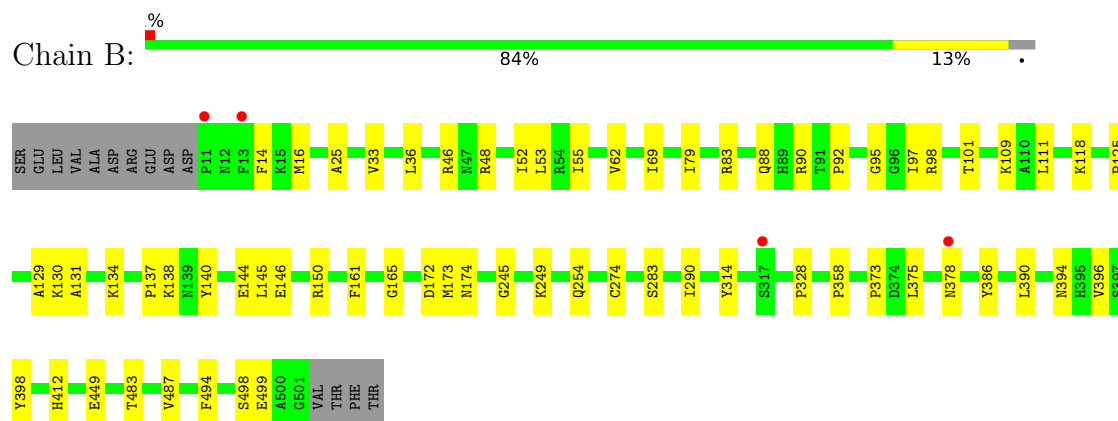
3 Residue-property plots [i](#)

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.

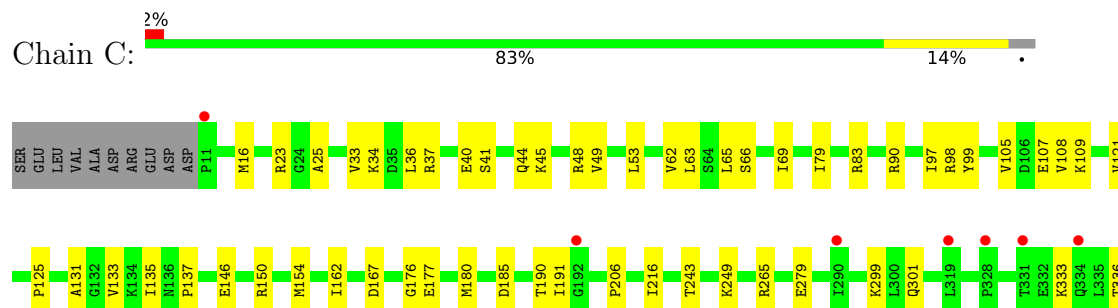
- Molecule 1: Glutamate dehydrogenase 2, mitochondrial



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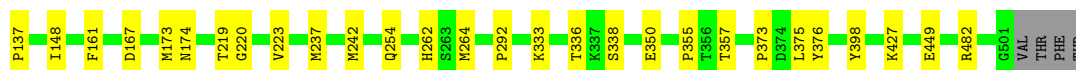
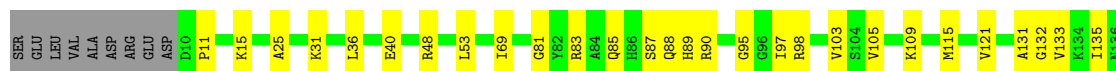
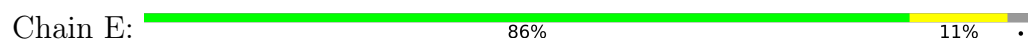




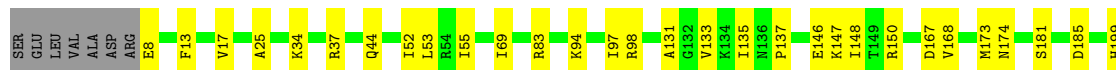
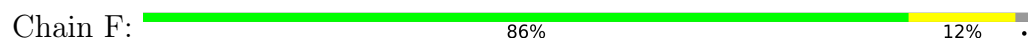
- Molecule 1: Glutamate dehydrogenase 2, mitochondrial



- Molecule 1: Glutamate dehydrogenase 2, mitochondrial



- Molecule 1: Glutamate dehydrogenase 2, mitochondrial



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	121.80Å 149.30Å 433.31Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.65 – 2.93 47.65 – 2.93	Depositor EDS
% Data completeness (in resolution range)	98.6 (47.65-2.93) 98.7 (47.65-2.93)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.51 (at 2.96Å)	Xtriage
Refinement program	PHENIX 1.12_2829, PHENIX 1.12_2829	Depositor
R, R_{free}	0.193 , 0.249 (Not available) , 0.189	Depositor DCC
R_{free} test set	4191 reflections (2.58%)	wwPDB-VP
Wilson B-factor (Å ²)	55.0	Xtriage
Anisotropy	0.318	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 47.9	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	0.93	EDS
Total number of atoms	24160	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.95% 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: CL, NA, PO4

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	0.08	0/3893	0.24	0/5253
1	B	0.10	0/3688	0.26	0/4960
1	C	0.08	0/3609	0.24	0/4848
1	D	0.09	0/3905	0.23	0/5270
1	E	0.10	0/3889	0.24	0/5249
1	F	0.09	0/3893	0.25	0/5252
All	All	0.09	0/22877	0.24	0/30832

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

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	3809	0	3707	45	0
1	B	3605	0	3227	41	0
1	C	3532	0	3109	42	0
1	D	3821	0	3717	30	0
1	E	3805	0	3737	32	0
1	F	3809	0	3714	38	0
2	A	20	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	25	0	0	0	0
2	C	25	0	0	0	0
2	D	10	0	0	0	0
2	E	10	0	0	0	0
2	F	25	0	0	0	0
3	A	2	0	0	0	0
3	C	2	0	0	0	0
3	D	11	0	0	0	0
3	E	3	0	0	0	0
3	F	3	0	0	0	0
4	A	3	0	0	0	0
4	B	2	0	0	0	0
4	C	2	0	0	0	0
4	D	5	0	0	0	0
4	E	6	0	0	0	0
4	F	1	0	0	0	0
5	A	296	0	0	2	0
5	B	197	0	0	2	0
5	C	235	0	0	1	0
5	D	365	0	0	0	0
5	E	296	0	0	1	0
5	F	235	0	0	1	0
All	All	24160	0	21211	221	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 (221) 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:83:ARG:HD2	1:E:131:ALA:HB2	1.65	0.79
1:B:83:ARG:HD2	1:B:131:ALA:HB2	1.67	0.76
1:F:69:ILE:HD12	1:F:148:ILE:HG13	1.68	0.74
1:C:69:ILE:HG12	1:C:79:ILE:HD11	1.71	0.72
1:C:216:ILE:HD12	1:C:216:ILE:H	1.55	0.71
1:F:69:ILE:HD11	1:F:147:LYS:HB2	1.73	0.70
1:E:98:ARG:NH1	1:E:173:MET:SD	2.67	0.67
1:B:254:GLN:HB3	1:B:328:PRO:HA	1.74	0.67
1:E:95:GLY:HA2	1:E:115:MET:HE3	1.76	0.66
1:F:227:ILE:HD11	1:F:267:LEU:HD11	1.77	0.66
1:F:52:ILE:HA	1:F:55:ILE:HD12	1.78	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:98:ARG:HG3	1:F:173:MET:HB3	1.77	0.65
1:C:98:ARG:NH1	1:C:107:GLU:OE1	2.29	0.65
1:F:327:ILE:HG12	1:F:349:ALA:HB3	1.80	0.64
1:F:398:TYR:HB2	1:F:449:GLU:HG3	1.79	0.64
1:C:333:LYS:H	1:C:355:PRO:HA	1.63	0.62
1:C:83:ARG:HD3	1:C:131:ALA:HB2	1.81	0.62
1:A:220:GLY:HA3	1:A:262:HIS:HB2	1.81	0.61
1:C:243:THR:O	1:C:249:LYS:NZ	2.33	0.61
1:F:137:PRO:HG2	1:F:174:ASN:HB3	1.81	0.61
1:A:25:ALA:HB1	1:A:53:LEU:HD13	1.82	0.60
1:A:424:LYS:HD3	1:D:432:ILE:HG23	1.82	0.60
1:C:190:THR:OG1	1:C:191:ILE:N	2.31	0.60
1:E:398:TYR:HB2	1:E:449:GLU:HG3	1.84	0.60
1:D:71:ARG:NH2	1:D:139:ASN:O	2.33	0.60
1:A:343:VAL:O	1:A:367:ARG:NH2	2.32	0.59
1:F:34:LYS:O	1:F:37:ARG:NH1	2.36	0.59
1:B:130:LYS:NZ	1:B:172:ASP:OD2	2.36	0.58
1:F:94:LYS:HD2	1:F:168:VAL:HB	1.84	0.58
1:E:137:PRO:HG2	1:E:174:ASN:HB3	1.84	0.58
1:A:320:GLU:OE2	1:A:342:ARG:NH1	2.33	0.57
1:E:36:LEU:HD21	1:E:48:ARG:HD2	1.86	0.57
1:C:437:THR:OG1	1:C:440:PHE:HB2	2.05	0.57
1:B:118:LYS:HD2	1:B:378:ASN:O	2.05	0.57
1:C:150:ARG:NH2	1:C:185:ASP:OD2	2.37	0.56
1:B:16:MET:SD	1:B:358:PRO:HD3	2.45	0.56
1:E:83:ARG:HG2	1:E:161:PHE:HD2	1.69	0.56
1:B:97:ILE:HG12	1:B:131:ALA:HB3	1.88	0.56
1:D:286:ASN:ND2	1:D:310:LYS:O	2.39	0.56
1:F:261:LEU:HD11	1:F:296:GLU:HG3	1.87	0.55
1:D:25:ALA:HB1	1:D:53:LEU:HD13	1.88	0.55
1:F:83:ARG:HD3	1:F:131:ALA:HB2	1.89	0.55
1:F:336:THR:HG22	1:F:338:SER:H	1.72	0.55
1:C:177:GLU:HB3	1:C:206:PRO:HG3	1.89	0.54
1:A:97:ILE:HG12	1:A:131:ALA:HB3	1.89	0.54
1:F:150:ARG:NH2	1:F:185:ASP:OD2	2.40	0.54
1:B:14:PHE:HD1	1:B:16:MET:HE2	1.73	0.54
1:C:279:GLU:OE2	1:C:299:LYS:NZ	2.40	0.54
1:F:25:ALA:HB1	1:F:53:LEU:HD13	1.89	0.54
1:B:52:ILE:HA	1:B:55:ILE:HD12	1.90	0.53
1:B:98:ARG:HG3	1:B:173:MET:HB3	1.89	0.53
1:C:97:ILE:HG12	1:C:131:ALA:HB3	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:133:VAL:HG23	1:F:135:ILE:HG12	1.91	0.53
1:F:335:LEU:HB2	1:F:356:THR:HG22	1.90	0.53
1:B:245:GLY:O	1:B:249:LYS:NZ	2.41	0.53
1:E:220:GLY:HA3	1:E:262:HIS:HB2	1.90	0.53
1:A:16:MET:HE3	1:A:358:PRO:HD3	1.91	0.53
1:F:322:ASP:HA	1:F:344:LYS:HB2	1.91	0.53
1:A:83:ARG:HH11	1:A:131:ALA:HB2	1.74	0.53
1:A:90:ARG:HD3	1:A:125:PRO:HA	1.91	0.53
1:E:133:VAL:HG23	1:E:135:ILE:HG12	1.91	0.53
1:C:34:LYS:O	1:C:37:ARG:NH1	2.42	0.52
1:A:240:LEU:HD21	1:A:370:LEU:HB3	1.90	0.52
1:D:373:PRO:HB2	1:D:375:LEU:HD23	1.91	0.52
1:E:25:ALA:HB1	1:E:53:LEU:HD13	1.90	0.52
1:C:133:VAL:HG12	1:C:135:ILE:HG12	1.90	0.52
1:C:90:ARG:HD3	1:C:125:PRO:HA	1.91	0.52
1:C:40:GLU:OE2	1:C:48:ARG:NH2	2.42	0.52
1:B:494:PHE:O	1:B:498:SER:OG	2.26	0.52
1:A:83:ARG:HD2	1:A:131:ALA:HB2	1.92	0.52
1:D:111:LEU:HD13	1:D:130:LYS:HE3	1.92	0.51
1:A:275:ILE:HD11	1:A:323:CYS:HB3	1.92	0.51
1:D:16:MET:HE3	1:D:358:PRO:HD3	1.92	0.51
1:D:223:VAL:HA	1:D:377:LEU:HG	1.93	0.51
1:F:146:GLU:OE2	1:F:150:ARG:NH1	2.44	0.51
1:A:137:PRO:HG2	1:A:174:ASN:HB3	1.93	0.51
1:E:373:PRO:HB2	1:E:375:LEU:HD23	1.92	0.51
1:F:373:PRO:HB2	1:F:375:LEU:HD23	1.93	0.51
1:A:286:ASN:ND2	5:A:702:HOH:O	2.33	0.51
1:F:348:ILE:HB	1:F:371:VAL:HG22	1.93	0.51
1:A:36:LEU:HD11	1:A:494:PHE:HZ	1.76	0.50
1:C:471:THR:HA	1:C:474:LYS:HB3	1.92	0.50
1:B:398:TYR:HB2	1:B:449:GLU:HG3	1.93	0.50
1:B:83:ARG:HG2	1:B:161:PHE:HD2	1.76	0.50
1:F:44:GLN:NE2	5:F:705:HOH:O	2.44	0.50
1:B:48:ARG:NH2	5:B:704:HOH:O	2.45	0.50
1:E:85:GLN:OE1	1:E:88:GLN:NE2	2.45	0.49
1:A:373:PRO:HB2	1:A:375:LEU:HD23	1.94	0.49
1:B:95:GLY:HA3	1:B:129:ALA:O	2.12	0.49
1:B:483:THR:O	1:B:487:VAL:HG12	2.11	0.49
1:A:36:LEU:HB3	1:A:38:THR:HG23	1.94	0.49
1:C:499:GLU:O	1:F:181:SER:OG	2.26	0.49
1:B:118:LYS:HD3	1:B:375:LEU:HA	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:227:ILE:HG22	1:F:372:ILE:HD12	1.92	0.49
1:B:25:ALA:HB1	1:B:53:LEU:HD13	1.94	0.49
1:A:8:GLU:N	1:A:8:GLU:OE1	2.46	0.49
1:A:105:VAL:O	1:A:109:LYS:HG3	2.13	0.49
1:C:460:THR:OG1	1:F:400:ARG:NH1	2.45	0.49
1:B:274:CYS:O	1:B:290:ILE:N	2.45	0.48
1:D:361:ASP:OD2	1:D:482:ARG:NH2	2.46	0.48
1:F:97:ILE:HA	1:F:131:ALA:HB3	1.94	0.48
1:D:83:ARG:HH11	1:D:131:ALA:HB2	1.78	0.48
1:E:11:PRO:HB3	1:E:15:LYS:NZ	2.28	0.48
1:E:237:MET:HE1	1:E:242:MET:HE3	1.95	0.48
1:A:133:VAL:HG12	1:A:135:ILE:HG12	1.95	0.48
1:E:31:LYS:NZ	5:E:706:HOH:O	2.47	0.48
1:E:83:ARG:NH2	1:E:167:ASP:OD1	2.40	0.48
1:C:62:VAL:HG21	1:C:109:LYS:HD3	1.95	0.48
1:C:162:ILE:HA	1:C:167:ASP:HB3	1.95	0.48
1:C:33:VAL:HG22	1:C:49:VAL:HG21	1.95	0.48
1:F:372:ILE:HG21	1:F:377:LEU:HD13	1.96	0.48
1:B:111:LEU:HB2	1:B:130:LYS:HD2	1.95	0.48
1:E:336:THR:HG23	1:E:338:SER:H	1.79	0.48
1:E:350:GLU:OE2	1:E:482:ARG:NH2	2.47	0.47
1:C:265:ARG:NH1	5:C:710:HOH:O	2.47	0.47
1:B:101:THR:O	1:B:134:LYS:NZ	2.47	0.47
1:E:121:VAL:HG11	1:E:376:TYR:HB2	1.96	0.47
1:A:362:LYS:O	1:A:366:GLU:HG3	2.14	0.47
1:B:394:ASN:HB3	1:B:396:VAL:HG12	1.96	0.47
1:E:264:MET:HE2	1:E:292:PRO:HG3	1.97	0.47
1:A:256:PHE:HD2	1:A:299:LYS:HD2	1.79	0.47
1:C:121:VAL:HG21	1:C:375:LEU:HD12	1.96	0.47
1:A:236:TYR:O	1:A:240:LEU:HD13	2.15	0.47
1:D:221:ARG:HG2	1:D:266:TYR:CE2	2.50	0.47
1:E:333:LYS:HG2	1:E:357:THR:HG22	1.97	0.47
1:A:242:MET:HA	1:E:427:LYS:HE2	1.96	0.46
1:D:151:ARG:NH1	1:D:155:GLU:OE1	2.48	0.46
1:A:223:VAL:HA	1:A:377:LEU:HG	1.97	0.46
1:A:262:HIS:CD2	1:A:265:ARG:HH21	2.32	0.46
1:C:154:MET:HE2	1:C:154:MET:HB3	1.78	0.46
1:A:357:THR:HG23	1:A:360:ALA:H	1.80	0.46
1:F:83:ARG:HH21	1:F:167:ASP:CG	2.23	0.46
1:E:97:ILE:HG12	1:E:131:ALA:HB3	1.97	0.46
1:A:190:THR:OG1	1:A:191:ILE:N	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:386:TYR:CZ	1:A:390:LEU:HD11	2.51	0.46
1:A:224:PHE:HD1	1:A:267:LEU:HD22	1.80	0.46
1:B:373:PRO:HB2	1:B:375:LEU:HD23	1.97	0.46
1:D:211:GLY:O	1:D:391:LYS:NZ	2.48	0.46
1:D:97:ILE:HG12	1:D:131:ALA:HB3	1.98	0.46
1:B:111:LEU:HD13	1:B:130:LYS:HE2	1.97	0.45
1:D:372:ILE:HG21	1:D:377:LEU:HD13	1.98	0.45
1:E:69:ILE:HG21	1:E:148:ILE:HG12	1.96	0.45
1:E:87:SER:OG	1:E:89:HIS:ND1	2.43	0.45
1:A:40:GLU:HB3	1:A:44:GLN:HB2	1.99	0.45
1:D:254:GLN:NE2	1:D:319:LEU:HD21	2.31	0.45
1:A:151:ARG:NH1	1:A:155:GLU:OE1	2.49	0.45
1:C:105:VAL:HA	1:C:108:VAL:HG22	1.97	0.45
1:C:386:TYR:CZ	1:C:390:LEU:HD11	2.51	0.45
1:E:105:VAL:O	1:E:109:LYS:HG3	2.17	0.45
1:A:374:ASP:N	1:A:374:ASP:OD1	2.49	0.45
1:D:217:SER:O	1:D:221:ARG:HG3	2.16	0.45
1:B:90:ARG:HD3	1:B:125:PRO:HA	1.99	0.45
1:C:41:SER:HB2	1:C:44:GLN:HG3	1.98	0.45
1:C:336:THR:OG1	1:C:337:LYS:N	2.50	0.45
1:D:83:ARG:HH21	1:D:167:ASP:CG	2.25	0.45
1:A:350:GLU:OE2	1:A:482:ARG:NH2	2.50	0.45
1:B:138:LYS:HD3	1:B:138:LYS:HA	1.57	0.44
1:E:333:LYS:H	1:E:355:PRO:HA	1.83	0.44
1:A:440:PHE:CG	1:B:412:HIS:HB3	2.53	0.44
1:F:83:ARG:HH11	1:F:131:ALA:HB2	1.83	0.44
1:C:23:ARG:HE	1:C:23:ARG:HB2	1.66	0.44
1:C:66:SER:O	1:C:66:SER:OG	2.29	0.44
1:D:101:THR:O	1:D:134:LYS:NZ	2.50	0.44
1:D:150:ARG:NH2	1:D:185:ASP:OD2	2.50	0.44
1:A:466:ARG:NH1	5:A:712:HOH:O	2.47	0.44
1:B:69:ILE:HG12	1:B:79:ILE:HD11	2.00	0.44
1:C:146:GLU:OE1	1:C:150:ARG:NH1	2.50	0.44
1:D:154:MET:HE2	1:D:154:MET:HB3	1.91	0.44
1:B:88:GLN:NE2	5:B:707:HOH:O	2.50	0.44
1:B:140:TYR:HB3	1:B:144:GLU:HB3	2.00	0.44
1:B:172:ASP:OD1	1:B:173:MET:N	2.51	0.43
1:C:440:PHE:O	1:C:444:ILE:HG13	2.18	0.43
1:A:117:TYR:O	1:A:121:VAL:HG23	2.17	0.43
1:F:240:LEU:HD12	1:F:346:LYS:HB3	2.00	0.43
1:A:365:LEU:HD21	1:A:480:ASP:HB2	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:386:TYR:CZ	1:B:390:LEU:HD11	2.52	0.43
1:B:499:GLU:O	1:D:181:SER:OG	2.30	0.43
1:E:89:HIS:CD2	1:E:90:ARG:HG3	2.53	0.43
1:C:63:LEU:HD21	1:C:65:LEU:HD21	2.01	0.43
1:F:199:HIS:O	1:F:205:LYS:NZ	2.46	0.43
1:A:428:HIS:HB3	1:F:302:HIS:CD2	2.53	0.43
1:D:92:PRO:HB3	1:D:165:GLY:C	2.43	0.43
1:C:399:GLY:HA2	1:C:402:THR:OG1	2.18	0.43
1:D:136:ASN:C	1:D:136:ASN:OD1	2.61	0.43
1:E:219:THR:O	1:E:223:VAL:HG13	2.19	0.43
1:F:262:HIS:HD2	1:F:265:ARG:HH21	1.67	0.43
1:A:109:LYS:HB2	1:A:109:LYS:HE2	1.82	0.43
1:B:283:SER:OG	1:B:314:TYR:O	2.36	0.43
1:C:16:MET:SD	1:C:358:PRO:HD3	2.59	0.43
1:C:176:GLY:O	1:C:180:MET:HG2	2.19	0.43
1:C:146:GLU:O	1:C:150:ARG:HG3	2.18	0.42
1:B:62:VAL:HG21	1:B:109:LYS:HD2	2.01	0.42
1:D:49:VAL:HG12	1:D:494:PHE:CZ	2.54	0.42
1:B:137:PRO:HG2	1:B:174:ASN:HB3	1.99	0.42
1:F:220:GLY:HA3	1:F:262:HIS:HB2	2.01	0.42
1:D:48:ARG:HE	1:D:48:ARG:HB2	1.73	0.42
1:F:374:ASP:N	1:F:374:ASP:OD1	2.47	0.42
1:A:36:LEU:HD21	1:A:494:PHE:HE2	1.84	0.42
1:B:36:LEU:HA	1:B:36:LEU:HD23	1.79	0.42
1:C:350:GLU:OE1	1:C:355:PRO:HD2	2.19	0.42
1:E:103:VAL:HG21	1:E:132:GLY:HA3	2.01	0.42
1:F:13:PHE:O	1:F:17:VAL:HG23	2.19	0.42
1:B:137:PRO:HA	1:B:145:LEU:HD21	2.02	0.41
1:A:370:LEU:HD11	1:A:481:LEU:CD1	2.50	0.41
1:A:237:MET:HE1	1:A:242:MET:HE3	2.02	0.41
1:D:13:PHE:HZ	1:D:114:LEU:HD22	1.86	0.41
1:D:79:ILE:HD13	1:D:135:ILE:HG21	2.02	0.41
1:E:81:GLY:HA2	1:E:133:VAL:HA	2.02	0.41
1:F:466:ARG:NH1	1:F:469:MET:SD	2.93	0.41
1:B:92:PRO:HB3	1:B:165:GLY:C	2.45	0.41
1:D:94:LYS:HD2	1:D:168:VAL:HB	2.03	0.41
1:D:421:LEU:HB3	1:D:432:ILE:HD12	2.03	0.41
1:E:40:GLU:OE1	1:E:48:ARG:NH2	2.39	0.41
1:B:146:GLU:O	1:B:150:ARG:HG3	2.20	0.40
1:C:36:LEU:O	1:C:45:LYS:NZ	2.43	0.40
1:A:335:LEU:HA	1:A:339:ASN:HD21	1.85	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:99:TYR:HB3	1:C:137:PRO:HG3	2.03	0.40
1:A:13:PHE:O	1:A:17:VAL:HG23	2.21	0.40
1:B:33:VAL:HG21	1:B:46:ARG:HG2	2.03	0.40
1:C:25:ALA:HB1	1:C:53:LEU:HD13	2.03	0.40
1:F:97:ILE:HG12	1:F:131:ALA:HB3	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

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	B	67/423 (16%)	67 (100%)	0	100	100
1	C	119/423 (28%)	118 (99%)	1 (1%)	73	84
1	D	372/423 (88%)	372 (100%)	0	100	100
1	E	392/423 (93%)	391 (100%)	1 (0%)	86	93
1	F	384/423 (91%)	383 (100%)	1 (0%)	86	93
All	All	1334/2115 (63%)	1331 (100%)	3 (0%)	87	95

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	301	GLN
1	E	254	GLN
1	F	8	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	88	GLN
1	C	262	HIS
1	D	174	ASN
1	D	199	HIS
1	D	254	GLN
1	D	268	HIS
1	D	441	GLN
1	E	193	HIS
1	E	209	GLN
1	E	441	GLN
1	F	262	HIS
1	F	441	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 63 ligands modelled in this entry, 40 are monoatomic - leaving 23 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	$\# Z > 2$
2	PO4	F	604	-	4,4,4	0.93	0	6,6,6	0.48	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	PO4	F	605	-	4,4,4	0.95	0	6,6,6	0.46	0
2	PO4	D	602	-	4,4,4	0.96	0	6,6,6	0.46	0
2	PO4	C	602	-	4,4,4	0.96	0	6,6,6	0.47	0
2	PO4	C	605	-	4,4,4	0.96	0	6,6,6	0.47	0
2	PO4	E	601	-	4,4,4	0.95	0	6,6,6	0.44	0
2	PO4	E	602	-	4,4,4	0.96	0	6,6,6	0.44	0
2	PO4	F	601	-	4,4,4	0.95	0	6,6,6	0.46	0
2	PO4	F	602	-	4,4,4	0.94	0	6,6,6	0.50	0
2	PO4	B	602	-	4,4,4	0.97	0	6,6,6	0.45	0
2	PO4	B	601	-	4,4,4	0.94	0	6,6,6	0.46	0
2	PO4	B	604	-	4,4,4	0.94	0	6,6,6	0.47	0
2	PO4	B	603	-	4,4,4	0.94	0	6,6,6	0.49	0
2	PO4	A	604	-	4,4,4	0.93	0	6,6,6	0.46	0
2	PO4	A	603	-	4,4,4	0.94	0	6,6,6	0.48	0
2	PO4	B	605	-	4,4,4	0.95	0	6,6,6	0.45	0
2	PO4	D	601	-	4,4,4	0.96	0	6,6,6	0.46	0
2	PO4	A	601	-	4,4,4	0.94	0	6,6,6	0.47	0
2	PO4	A	602	-	4,4,4	0.95	0	6,6,6	0.44	0
2	PO4	F	603	-	4,4,4	0.96	0	6,6,6	0.46	0
2	PO4	C	604	-	4,4,4	0.96	0	6,6,6	0.45	0
2	PO4	C	603	-	4,4,4	0.95	0	6,6,6	0.46	0
2	PO4	C	601	-	4,4,4	0.94	0	6,6,6	0.49	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.

No monomer is involved in short contacts.

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	494/505 (97%)	-0.19	1 (0%) 91 89	28, 46, 71, 83	0
1	B	491/505 (97%)	0.02	4 (0%) 82 77	30, 53, 81, 109	1 (0%)
1	C	491/505 (97%)	0.17	9 (1%) 67 59	34, 56, 94, 106	1 (0%)
1	D	496/505 (98%)	-0.30	0 100 100	29, 43, 62, 81	0
1	E	492/505 (97%)	-0.10	0 100 100	30, 48, 73, 91	1 (0%)
1	F	496/505 (98%)	-0.02	2 (0%) 88 86	31, 52, 80, 106	0
All	All	2960/3030 (97%)	-0.07	16 (0%) 87 83	28, 49, 81, 109	3 (0%)

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	503	THR	3.3
1	C	11	PRO	3.2
1	B	13	PHE	3.0
1	B	378	ASN	3.0
1	A	281	ASP	2.6
1	B	11	PRO	2.5
1	C	290	ILE	2.4
1	C	334	GLN	2.4
1	C	356	THR	2.4
1	C	368	ASN	2.4
1	C	319	LEU	2.4
1	C	192	GLY	2.2
1	B	317	SER	2.2
1	C	331	THR	2.1
1	F	502	VAL	2.0
1	C	328	PRO	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($\text{\AA}^2$)	Q<0.9
2	PO4	C	605	5/5	0.44	0.10	134,141,148,149	0
3	NA	C	607	1/1	0.70	0.19	92,92,92,92	0
3	NA	D	606	1/1	0.72	0.14	80,80,80,80	0
2	PO4	B	603	5/5	0.75	0.12	75,84,90,104	0
3	NA	A	605	1/1	0.79	0.17	58,58,58,58	0
3	NA	D	608	1/1	0.80	0.13	83,83,83,83	0
4	CL	D	615	1/1	0.82	0.08	74,74,74,74	0
2	PO4	C	604	5/5	0.83	0.15	85,86,104,110	0
3	NA	E	605	1/1	0.84	0.13	71,71,71,71	0
4	CL	A	609	1/1	0.84	0.09	87,87,87,87	0
2	PO4	B	605	5/5	0.84	0.15	72,73,96,97	0
3	NA	D	604	1/1	0.85	0.17	63,63,63,63	0
4	CL	A	608	1/1	0.85	0.09	83,83,83,83	0
4	CL	E	609	1/1	0.85	0.12	82,82,82,82	0
4	CL	B	607	1/1	0.87	0.08	89,89,89,89	0
2	PO4	F	602	5/5	0.87	0.12	65,71,76,80	0
4	CL	E	606	1/1	0.87	0.07	80,80,80,80	0
4	CL	E	608	1/1	0.87	0.14	84,84,84,84	0
2	PO4	F	605	5/5	0.87	0.20	89,95,97,116	0
3	NA	F	607	1/1	0.88	0.20	65,65,65,65	0
4	CL	D	618	1/1	0.88	0.18	93,93,93,93	0
3	NA	D	609	1/1	0.88	0.14	57,57,57,57	0
3	NA	D	610	1/1	0.88	0.08	61,61,61,61	0
2	PO4	E	602	5/5	0.88	0.09	59,71,79,84	0
4	CL	E	610	1/1	0.88	0.15	78,78,78,78	0
3	NA	D	611	1/1	0.89	0.16	51,51,51,51	0
4	CL	D	614	1/1	0.89	0.08	70,70,70,70	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NA	E	603	1/1	0.89	0.20	65,65,65,65	0
3	NA	E	604	1/1	0.89	0.15	66,66,66,66	0
3	NA	D	607	1/1	0.89	0.15	72,72,72,72	0
4	CL	E	607	1/1	0.89	0.13	81,81,81,81	0
2	PO4	F	603	5/5	0.89	0.13	53,59,65,66	0
2	PO4	B	602	5/5	0.89	0.10	62,65,72,81	0
2	PO4	C	603	5/5	0.89	0.11	56,62,74,79	0
4	CL	C	609	1/1	0.90	0.08	74,74,74,74	0
2	PO4	B	604	5/5	0.90	0.21	67,76,86,94	0
2	PO4	E	601	5/5	0.90	0.14	58,62,67,78	0
4	CL	F	609	1/1	0.90	0.10	80,80,80,80	0
2	PO4	F	604	5/5	0.91	0.31	60,61,67,73	0
3	NA	D	605	1/1	0.91	0.07	66,66,66,66	0
2	PO4	A	604	5/5	0.91	0.15	56,59,69,81	0
2	PO4	C	602	5/5	0.91	0.09	69,78,91,91	0
2	PO4	A	602	5/5	0.91	0.11	65,66,76,83	0
2	PO4	B	601	5/5	0.92	0.10	56,60,67,79	0
2	PO4	C	601	5/5	0.92	0.12	49,56,66,72	0
2	PO4	D	601	5/5	0.92	0.09	51,55,62,67	0
4	CL	B	606	1/1	0.93	0.11	73,73,73,73	0
3	NA	C	606	1/1	0.93	0.17	61,61,61,61	0
3	NA	A	606	1/1	0.93	0.10	58,58,58,58	0
3	NA	D	612	1/1	0.93	0.07	47,47,47,47	0
2	PO4	D	602	5/5	0.94	0.07	65,66,71,79	0
2	PO4	A	601	5/5	0.94	0.15	49,55,67,74	0
4	CL	D	616	1/1	0.95	0.09	74,74,74,74	0
2	PO4	F	601	5/5	0.95	0.09	67,70,80,89	0
3	NA	F	608	1/1	0.95	0.14	53,53,53,53	0
3	NA	D	603	1/1	0.95	0.04	36,36,36,36	0
4	CL	D	617	1/1	0.96	0.05	65,65,65,65	0
3	NA	F	606	1/1	0.97	0.08	53,53,53,53	0
4	CL	A	607	1/1	0.97	0.10	59,59,59,59	0
2	PO4	A	603	5/5	0.97	0.16	42,44,56,62	0
4	CL	C	608	1/1	0.97	0.15	60,60,60,60	0
4	CL	E	611	1/1	0.98	0.12	56,56,56,56	0
3	NA	D	613	1/1	0.98	0.23	34,34,34,34	0

6.5 Other polymers [i](#)

There are no such residues in this entry.