



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 10:36 PM UTC

PDB ID : 4JAX / pdb_00004jax
Title : Crystal structure of dimeric KIHxk1 in crystal form X
Authors : Kuettner, E.B.; Strater, N.; Kettner, K.; Otto, A.; Lilie, H.; Golbik, R.P.;
Kriegel, T.M.
Deposited on : 2013-02-19
Resolution : 2.26 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:

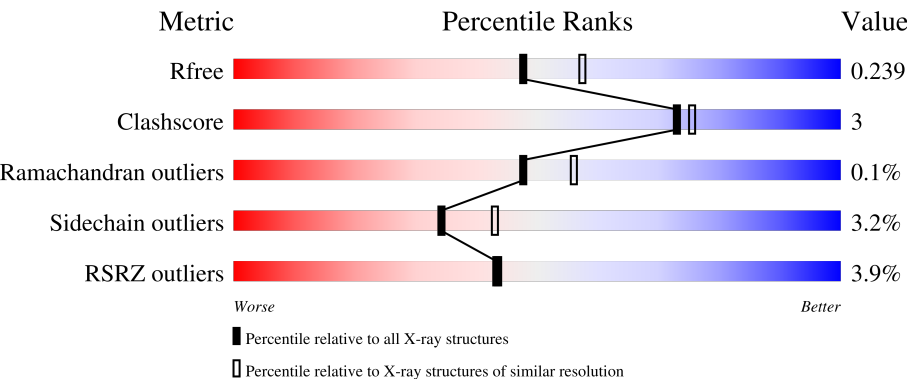
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry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

1 Overall quality at a glance i

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

The reported resolution of this entry is 2.26 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	1898 (2.26-2.26)
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)
RSRZ outliers	180081	1898 (2.26-2.26)

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	485	% 88%8% ..
1	B	485	% 87%9% ..
1	C	485	% 89%8% ..
1	D	485	% 89%8% .
1	E	485	2% 88%9% .

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	F	485	17% 81% 13% • 5%

2 Entry composition [i](#)

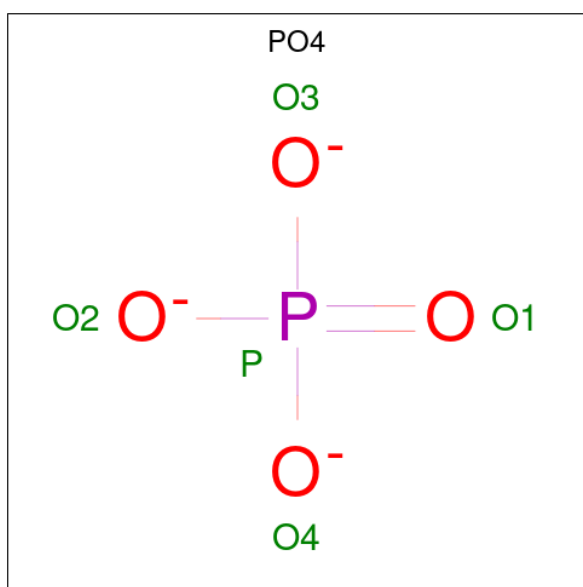
There are 4 unique types of molecules in this entry. The entry contains 22473 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Hexokinase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	470	Total	C	N	O	P	S	0	1	0
			3664	2330	603	714	1	16			
1	B	472	Total	C	N	O	P	S	0	2	0
			3681	2340	606	718	1	16			
1	C	475	Total	C	N	O	P	S	0	0	0
			3698	2349	610	722	1	16			
1	D	473	Total	C	N	O	P	S	0	0	0
			3682	2340	605	720	1	16			
1	E	471	Total	C	N	O	P	S	0	0	0
			3669	2332	602	718	1	16			
1	F	460	Total	C	N	O	P	S	0	0	0
			3592	2287	586	702	1	16			

- 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	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	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	E	1	Total O P 5 4 1	0	0
2	F	1	Total O P 5 4 1	0	0

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		
3	E	1	Total	C	O	0	0
			6	3	3		
3	E	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	110	Total	O	0	0
			110	110		
4	B	55	Total	O	0	0
			55	55		
4	C	72	Total	O	0	0
			72	72		
4	D	47	Total	O	0	0
			47	47		

Continued on next page...

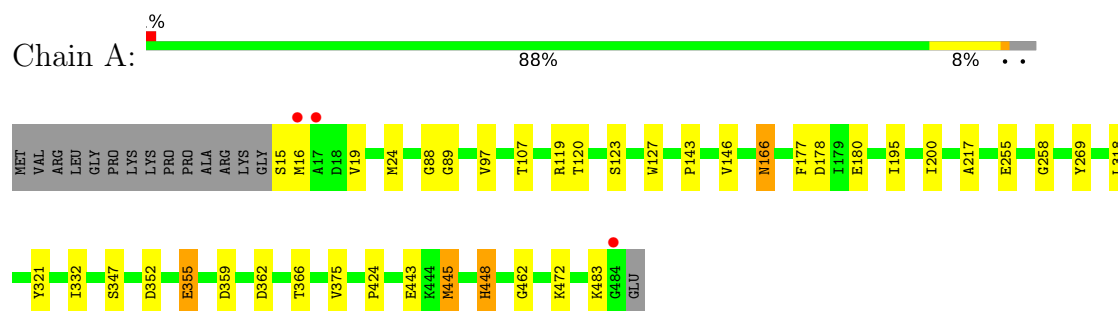
Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	E	36	Total 36	O 36	0	0
4	F	34	Total 34	O 34	0	0

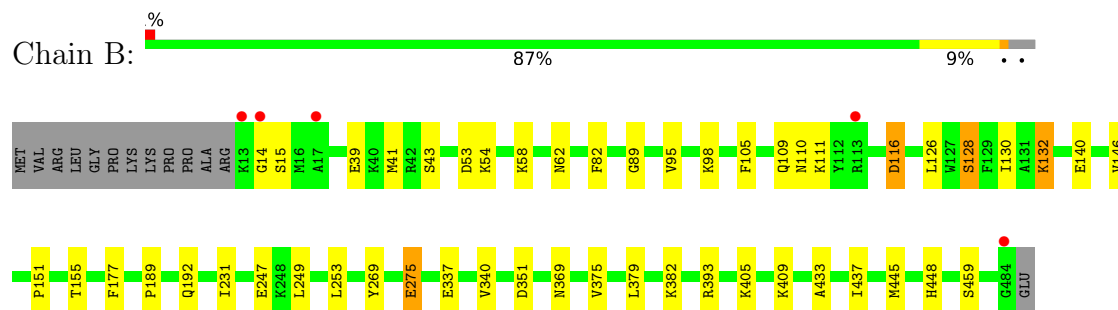
3 Residue-property plots [i](#)

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.

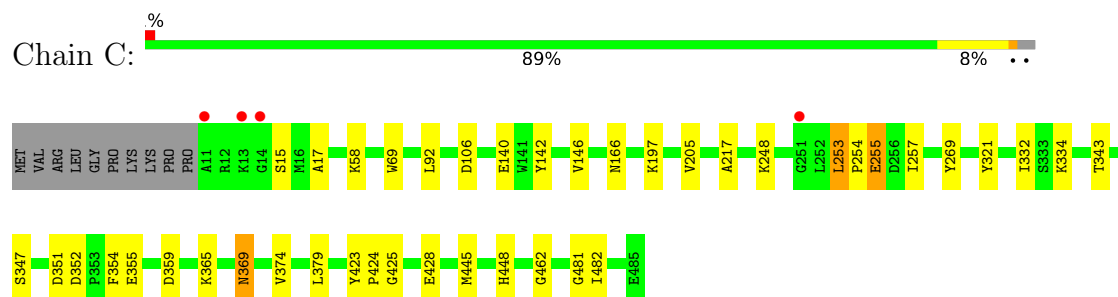
• Molecule 1: Hexokinase



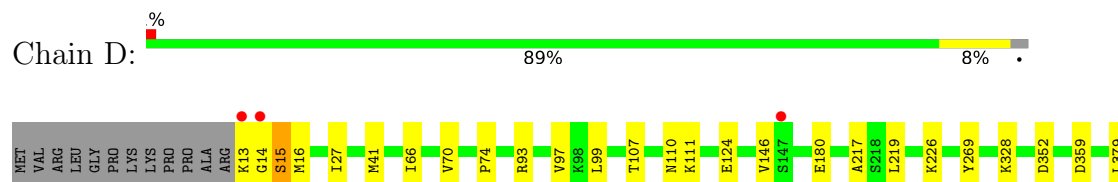
• Molecule 1: Hexokinase



• Molecule 1: Hexokinase

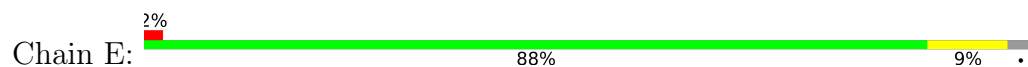


• Molecule 1: Hexokinase

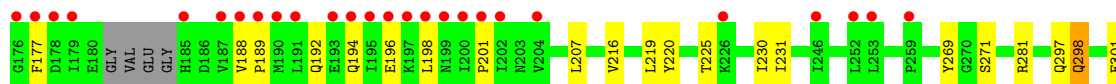
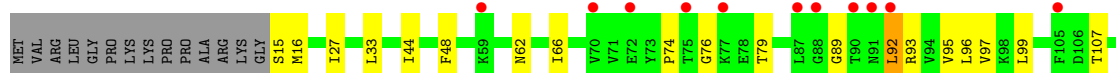
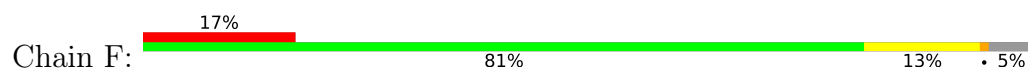




• Molecule 1: Hexokinase



• Molecule 1: Hexokinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	105.81Å 178.30Å 216.21Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.60 – 2.26 29.60 – 2.26	Depositor EDS
% Data completeness (in resolution range)	99.8 (29.60-2.26) 99.8 (29.60-2.26)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.80 (at 2.26Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.200 , 0.240 0.199 , 0.239	Depositor DCC
R_{free} test set	1911 reflections (1.00%)	wwPDB-VP
Wilson B-factor (Å ²)	38.6	Xtriage
Anisotropy	0.365	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 37.0	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	0.95	EDS
Total number of atoms	22473	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.28% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SEP, PO4, GOL

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	1/3733 (0.0%)	0.96	4/5057 (0.1%)
1	B	0.69	0/3754	0.92	0/5082
1	C	0.70	0/3760	0.94	2/5088 (0.0%)
1	D	0.65	0/3744	0.91	2/5067 (0.0%)
1	E	0.62	0/3732	0.89	2/5054 (0.0%)
1	F	0.70	0/3652	0.94	0/4943
All	All	0.69	1/22375 (0.0%)	0.93	10/30291 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	88	GLY	N-CA	5.78	1.49	1.44

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	142	TYR	CA-C-N	-10.88	109.76	120.21
1	C	142	TYR	C-N-CA	-10.88	109.76	120.21
1	A	258	GLY	CA-C-N	7.98	127.70	119.56
1	A	258	GLY	C-N-CA	7.98	127.70	119.56
1	A	448	HIS	CA-C-N	5.95	125.43	119.24
1	A	448	HIS	C-N-CA	5.95	125.43	119.24
1	D	70	VAL	N-CA-C	-5.73	98.98	107.51
1	D	66	ILE	N-CA-C	5.32	113.13	107.76
1	E	222	ASP	CA-C-N	5.00	125.27	119.47
1	E	222	ASP	C-N-CA	5.00	125.27	119.47

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	3664	0	3644	19	0
1	B	3681	0	3659	22	0
1	C	3698	0	3679	26	0
1	D	3682	0	3661	23	0
1	E	3669	0	3647	20	0
1	F	3592	0	3571	35	0
2	A	20	0	0	0	0
2	B	5	0	0	0	0
2	C	20	0	0	0	0
2	D	20	0	0	0	0
2	E	15	0	0	1	0
2	F	5	0	0	0	0
3	A	12	0	16	0	0
3	B	6	0	8	0	0
3	C	12	0	16	0	0
3	D	6	0	8	1	0
3	E	12	0	16	1	0
4	A	110	0	0	1	0
4	B	55	0	0	0	0
4	C	72	0	0	1	0
4	D	47	0	0	0	0
4	E	36	0	0	0	0
4	F	34	0	0	1	0
All	All	22473	0	21925	140	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 3.

All (140) 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:D:14:GLY:O	1:D:15:SEP:HB3	1.59	0.98
1:D:41:MET:HE2	1:D:437:ILE:HD11	1.50	0.93
1:F:97:VAL:HG22	1:F:107:THR:HG22	1.55	0.87
1:C:253:LEU:HD13	1:C:254:PRO:HD2	1.58	0.83

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:166:ASN:HD22	1:A:166:ASN:H	1.29	0.78
1:E:41:MET:HE2	1:E:437:ILE:HD11	1.66	0.76
1:C:334:LYS:HZ2	1:C:369:ASN:HB2	1.51	0.76
1:D:41:MET:HE2	1:D:437:ILE:CD1	2.14	0.76
1:D:16:MET:HE3	1:D:27:ILE:HD13	1.68	0.75
1:F:127:TRP:HB2	1:F:194:GLN:HG3	1.69	0.73
1:B:41:MET:HE3	1:B:393:ARG:O	1.89	0.72
1:B:189:PRO:HA	1:B:192:GLN:HE21	1.52	0.72
1:A:355:GLU:HG2	1:D:110:ASN:OD1	1.91	0.70
1:C:253:LEU:HD11	1:C:257:ILE:HB	1.76	0.67
1:B:116:ASP:OD2	1:B:116:ASP:N	2.26	0.66
1:F:115:PRO:O	1:F:116:ASP:HB3	1.94	0.65
1:D:110:ASN:HD22	1:D:111:LYS:H	1.45	0.65
1:C:352:ASP:CG	1:C:359:ASP:HB2	2.22	0.65
1:D:41:MET:HE1	1:D:433:ALA:HB1	1.80	0.64
1:D:217:ALA:HB2	1:D:462:GLY:HA2	1.81	0.63
1:C:334:LYS:NZ	1:C:369:ASN:HB2	2.14	0.62
1:E:110:ASN:HD22	1:E:111:LYS:H	1.50	0.59
1:A:97:VAL:HG22	1:A:107:THR:HG22	1.85	0.59
1:A:89:GLY:HA2	1:A:177:PHE:HE1	1.68	0.59
1:B:89:GLY:HA2	1:B:177:PHE:HE1	1.68	0.59
1:C:253:LEU:CD1	1:C:257:ILE:HB	2.33	0.59
1:E:41:MET:HE2	1:E:437:ILE:CD1	2.33	0.58
1:B:41:MET:HE1	1:B:433:ALA:HB1	1.86	0.58
1:B:58:LYS:HG3	1:B:247:GLU:HG2	1.86	0.58
1:B:62:ASN:ND2	1:B:275:GLU:OE2	2.38	0.57
1:F:219:LEU:HA	1:F:225:THR:HB	1.87	0.57
1:E:233:THR:HG23	2:E:501:PO4:O3	2.05	0.56
1:F:16:MET:HE1	1:F:27:ILE:HD13	1.87	0.56
1:A:19:VAL:HB	1:A:24:MET:HE2	1.88	0.55
1:D:217:ALA:HB2	1:D:462:GLY:CA	2.36	0.55
1:B:111:LYS:HB2	1:C:354:PHE:CE2	2.42	0.55
1:F:298:GLN:HB2	1:F:301:GLU:HB3	1.89	0.55
1:F:44:ILE:O	1:F:48:PHE:HB2	2.08	0.54
1:F:92:LEU:C	1:F:92:LEU:HD12	2.33	0.54
1:B:95:VAL:HG13	1:B:109:GLN:HB3	1.90	0.53
1:D:74:PRO:HD2	1:D:219:LEU:HD23	1.89	0.53
1:F:297:GLN:O	1:F:298:GLN:HG2	2.08	0.53
1:C:321:TYR:CE2	1:C:332:ILE:HG12	2.45	0.53
1:D:41:MET:HE3	1:D:393:ARG:O	2.09	0.52
1:F:89:GLY:HA2	1:F:177:PHE:CE1	2.45	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:445:MET:HA	1:F:448:HIS:CE1	2.45	0.52
1:C:347:SER:O	1:C:424:PRO:HG2	2.10	0.51
1:D:97:VAL:HG22	1:D:107:THR:HG22	1.92	0.51
1:C:253:LEU:HD13	1:C:254:PRO:CD	2.36	0.51
1:E:148:GLU:HB2	1:E:149:PRO:HD2	1.92	0.51
1:E:58:LYS:HG3	1:E:247:GLU:HG2	1.92	0.51
1:A:143:PRO:HG3	1:D:13:LYS:HD2	1.92	0.51
1:E:234:GLY:HA3	3:E:504:GOL:H31	1.94	0.50
1:F:89:GLY:HA2	1:F:177:PHE:HE1	1.76	0.50
1:F:93:ARG:HG2	1:F:111:LYS:HG3	1.93	0.50
1:B:41:MET:HE2	1:B:437:ILE:HD11	1.92	0.50
1:F:135:LYS:HD2	1:F:198:LEU:HD12	1.93	0.50
1:F:230:ILE:C	1:F:231:ILE:HG13	2.36	0.50
1:E:74:PRO:HD2	1:E:219:LEU:HD23	1.93	0.50
1:F:132:LYS:O	1:F:136:GLU:HG2	2.11	0.50
1:D:379:LEU:C	1:D:379:LEU:HD23	2.37	0.50
1:E:394:LEU:O	1:E:397:CYS:HB2	2.12	0.49
1:B:337[B]:GLU:HB3	1:B:340:VAL:HB	1.95	0.49
1:F:192:GLN:O	1:F:196:GLU:HG2	2.13	0.49
1:C:347:SER:HB2	1:C:424:PRO:HD3	1.95	0.49
1:F:79:THR:HA	1:F:99:LEU:O	2.13	0.48
1:D:110:ASN:ND2	1:D:111:LYS:H	2.10	0.48
1:E:16:MET:HE3	1:E:19:VAL:HG21	1.94	0.48
1:C:448:HIS:HD2	4:C:608:HOH:O	1.96	0.47
1:A:352:ASP:CG	1:A:359:ASP:HB2	2.40	0.47
1:B:155:THR:HG21	1:B:459:SER:HB2	1.96	0.47
1:F:76:GLY:HA2	1:F:220:TYR:CE1	2.49	0.47
1:C:347:SER:O	1:C:351:ASP:HB2	2.15	0.47
1:F:168:GLY:HA3	1:F:207:LEU:HD13	1.96	0.47
1:B:82:PHE:CD1	1:B:151:PRO:HG2	2.50	0.46
1:C:166:ASN:HB3	1:C:205:VAL:O	2.16	0.46
1:A:166:ASN:HD22	1:A:166:ASN:N	2.05	0.46
1:B:53:ASP:OD1	1:B:405:LYS:HE2	2.15	0.46
1:D:328:LYS:HB3	3:D:505:GOL:O3	2.15	0.46
1:B:110:ASN:OD1	1:C:355:GLU:HB2	2.16	0.46
1:F:394:LEU:O	1:F:397:CYS:HB2	2.16	0.46
1:D:226:LYS:HD3	1:D:408:TYR:CG	2.51	0.45
1:E:41:MET:HE2	1:E:437:ILE:CG1	2.47	0.45
1:A:321:TYR:CZ	1:A:332:ILE:HG12	2.52	0.45
1:E:379:LEU:C	1:E:379:LEU:HD23	2.41	0.45
1:F:165:ILE:HG22	1:F:480:VAL:HB	1.98	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:188:VAL:HB	1:F:189:PRO:HD3	1.98	0.45
1:A:119:ARG:HG3	4:A:659:HOH:O	2.17	0.45
1:A:120:THR:HG22	1:A:178:ASP:HB3	1.98	0.45
1:E:363:LEU:O	1:E:367:ASN:HB2	2.17	0.45
1:C:217:ALA:HB2	1:C:462:GLY:HA2	2.00	0.44
1:B:379:LEU:C	1:B:379:LEU:HD23	2.42	0.44
1:E:217:ALA:HB2	1:E:462:GLY:HA2	1.99	0.44
1:D:425:GLY:O	1:D:429:LYS:HG3	2.17	0.44
1:C:255:GLU:H	1:C:255:GLU:HG2	1.54	0.44
1:E:352:ASP:CG	1:E:359:ASP:HB2	2.43	0.44
1:A:217:ALA:HB2	1:A:462:GLY:HA2	2.00	0.43
1:A:445:MET:HA	1:A:448:HIS:CE1	2.53	0.43
1:C:15:SEP:O1P	1:C:17:ALA:HB2	2.17	0.43
1:F:74:PRO:HG3	1:F:216:VAL:HG13	2.00	0.43
1:E:46:LYS:HD3	1:E:46:LYS:HA	1.60	0.43
1:C:92:LEU:C	1:C:92:LEU:HD12	2.44	0.43
1:D:352:ASP:CG	1:D:359:ASP:HB2	2.43	0.43
1:B:445:MET:HA	1:B:448:HIS:CD2	2.54	0.43
1:C:425:GLY:HA2	1:C:428:GLU:OE1	2.19	0.42
1:F:298:GLN:HE21	1:F:298:GLN:HB3	1.53	0.42
1:B:249:LEU:HD13	1:B:253:LEU:HD21	2.01	0.42
1:C:423:TYR:HA	1:C:424:PRO:HD3	1.86	0.42
1:A:166:ASN:H	1:A:166:ASN:ND2	2.08	0.42
1:B:14:GLY:HA2	1:C:140:GLU:HG3	2.02	0.42
1:F:271:SER:HA	1:F:298:GLN:HA	2.02	0.42
1:C:69:TRP:CE3	1:C:481:GLY:HA2	2.54	0.42
1:A:123:SER:HG	1:A:127:TRP:CD1	2.37	0.42
1:E:358:GLU:H	1:E:358:GLU:HG2	1.57	0.42
1:F:142:TYR:CE1	1:F:150:LEU:HD11	2.55	0.42
1:F:281:ARG:HD2	4:F:615:HOH:O	2.19	0.42
1:D:482:ILE:HG13	1:D:485:GLU:HG2	2.02	0.41
1:F:163:LYS:HB3	1:F:485:GLU:HG3	2.00	0.41
1:A:195:ILE:HG23	1:A:200:ILE:HB	2.02	0.41
1:F:174:THR:O	1:F:175:LYS:HB2	2.20	0.41
1:F:116:ASP:OD1	1:F:116:ASP:C	2.63	0.41
1:F:440:TRP:HH2	1:F:450:ILE:HB	1.84	0.41
1:A:19:VAL:HB	1:A:24:MET:CE	2.50	0.41
1:B:98:LYS:O	1:B:105:PHE:HA	2.21	0.41
1:F:446:GLU:H	1:F:446:GLU:HG3	1.61	0.41
1:E:92:LEU:C	1:E:92:LEU:HD12	2.45	0.41
1:A:347:SER:O	1:A:424:PRO:HG2	2.20	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:379:LEU:C	1:C:379:LEU:HD23	2.45	0.41
1:F:297:GLN:C	1:F:298:GLN:HG2	2.46	0.41
1:D:99:LEU:HD21	1:D:465:ILE:HG12	2.03	0.41
1:E:77:LYS:O	1:E:101:GLY:HA2	2.21	0.41
1:B:126:LEU:O	1:B:130:ILE:HG13	2.20	0.40
1:C:445:MET:HA	1:C:448:HIS:CE1	2.56	0.40
1:F:96:LEU:O	1:F:107:THR:HA	2.21	0.40
1:B:128:SER:O	1:B:132:LYS:HB2	2.22	0.40
1:E:65:MET:O	1:E:248:LYS:NZ	2.52	0.40
1:C:58:LYS:HG2	1:C:248:LYS:HA	2.04	0.40
1:D:15:SEP:HB2	1:D:16:MET:H	1.68	0.40
1:D:435:LYS:HA	1:D:440:TRP:CE3	2.56	0.40
1:A:362:ASP:O	1:A:366:THR:HB	2.21	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	469/485 (97%)	450 (96%)	18 (4%)	1 (0%)	43	50
1	B	471/485 (97%)	450 (96%)	21 (4%)	0	100	100
1	C	472/485 (97%)	454 (96%)	18 (4%)	0	100	100
1	D	470/485 (97%)	447 (95%)	23 (5%)	0	100	100
1	E	469/485 (97%)	450 (96%)	19 (4%)	0	100	100
1	F	454/485 (94%)	428 (94%)	25 (6%)	1 (0%)	43	50
All	All	2805/2910 (96%)	2679 (96%)	124 (4%)	2 (0%)	48	56

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	16	MET
1	F	201	PRO

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	400/411 (97%)	388 (97%)	12 (3%)	36	45
1	B	402/411 (98%)	386 (96%)	16 (4%)	28	34
1	C	402/411 (98%)	391 (97%)	11 (3%)	39	49
1	D	401/411 (98%)	395 (98%)	6 (2%)	57	68
1	E	400/411 (97%)	386 (96%)	14 (4%)	32	39
1	F	392/411 (95%)	375 (96%)	17 (4%)	26	31
All	All	2397/2466 (97%)	2321 (97%)	76 (3%)	34	43

All (76) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	146	VAL
1	A	166	ASN
1	A	180	GLU
1	A	255	GLU
1	A	269	TYR
1	A	318	LEU
1	A	355	GLU
1	A	375	VAL
1	A	443	GLU
1	A	445	MET
1	A	472	LYS
1	A	483	LYS
1	B	39	GLU
1	B	43	SER
1	B	54	LYS
1	B	116	ASP
1	B	128	SER

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	132	LYS
1	B	140	GLU
1	B	146	VAL
1	B	231	ILE
1	B	269	TYR
1	B	275	GLU
1	B	351	ASP
1	B	369	ASN
1	B	375	VAL
1	B	382	LYS
1	B	409	LYS
1	C	106	ASP
1	C	146	VAL
1	C	197	LYS
1	C	253	LEU
1	C	255	GLU
1	C	269	TYR
1	C	343	THR
1	C	365	LYS
1	C	369	ASN
1	C	374	VAL
1	C	482	ILE
1	D	93	ARG
1	D	124	GLU
1	D	146	VAL
1	D	180	GLU
1	D	269	TYR
1	D	396	VAL
1	E	18	ASP
1	E	22	ASN
1	E	27	ILE
1	E	54	LYS
1	E	106	ASP
1	E	110	ASN
1	E	231	ILE
1	E	250	GLU
1	E	269	TYR
1	E	275	GLU
1	E	347	SER
1	E	358	GLU
1	E	369	ASN
1	E	472	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	F	33	LEU
1	F	62	ASN
1	F	66	ILE
1	F	92	LEU
1	F	95	VAL
1	F	114	LEU
1	F	116	ASP
1	F	124	GLU
1	F	136	GLU
1	F	161	SER
1	F	269	TYR
1	F	298	GLN
1	F	333	SER
1	F	378	LYS
1	F	396	VAL
1	F	446	GLU
1	F	480	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (37) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	166	ASN
1	A	171	GLN
1	A	194	GLN
1	A	209	ASN
1	A	224	GLN
1	B	22	ASN
1	B	117	HIS
1	B	192	GLN
1	B	203	ASN
1	B	276	HIS
1	B	471	GLN
1	C	47	HIS
1	C	102	ASN
1	C	110	ASN
1	C	367	ASN
1	C	369	ASN
1	C	448	HIS
1	D	102	ASN
1	D	110	ASN
1	D	203	ASN
1	D	356	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	367	ASN
1	D	448	HIS
1	D	471	GLN
1	E	22	ASN
1	E	102	ASN
1	E	110	ASN
1	E	194	GLN
1	E	276	HIS
1	E	471	GLN
1	F	110	ASN
1	F	125	GLN
1	F	194	GLN
1	F	298	GLN
1	F	367	ASN
1	F	369	ASN
1	F	448	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

6 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	SEP	A	15	1	8,9,10	0.56	0	7,12,14	1.84	2 (28%)
1	SEP	F	15	1	8,9,10	0.67	0	7,12,14	2.12	3 (42%)
1	SEP	E	15	1	8,9,10	0.67	0	7,12,14	1.43	1 (14%)
1	SEP	B	15	1	8,9,10	0.75	0	7,12,14	4.00	4 (57%)
1	SEP	D	15	1	8,9,10	0.61	0	7,12,14	2.46	1 (14%)
1	SEP	C	15	1	8,9,10	0.65	0	7,12,14	1.11	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
1	SEP	A	15	1	-	3/6/8/10	-
1	SEP	F	15	1	-	2/6/8/10	-
1	SEP	E	15	1	-	6/6/8/10	-
1	SEP	B	15	1	-	2/6/8/10	-
1	SEP	D	15	1	-	6/6/8/10	-
1	SEP	C	15	1	-	2/6/8/10	-

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	15	SEP	OG-CB-CA	8.96	116.87	108.14
1	D	15	SEP	OG-CB-CA	6.04	114.02	108.14
1	B	15	SEP	O2P-P-OG	-4.25	95.57	106.67
1	F	15	SEP	OG-CB-CA	3.88	111.92	108.14
1	A	15	SEP	OG-CB-CA	2.91	110.98	108.14
1	E	15	SEP	OG-CB-CA	2.69	110.76	108.14
1	B	15	SEP	OG-P-O1P	2.68	113.69	106.44
1	A	15	SEP	O2P-P-OG	-2.49	100.17	106.67
1	F	15	SEP	O3P-P-OG	-2.32	100.62	106.67
1	F	15	SEP	O3P-P-O2P	2.26	116.26	107.80
1	B	15	SEP	O3P-P-O2P	2.12	115.74	107.80

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	15	SEP	CB-OG-P-O1P
1	B	15	SEP	N-CA-CB-OG
1	B	15	SEP	C-CA-CB-OG
1	C	15	SEP	CA-CB-OG-P
1	D	15	SEP	N-CA-CB-OG
1	D	15	SEP	C-CA-CB-OG
1	D	15	SEP	CA-CB-OG-P
1	D	15	SEP	CB-OG-P-O1P
1	D	15	SEP	CB-OG-P-O2P

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
1	E	15	SEP	C-CA-CB-OG
1	E	15	SEP	CA-CB-OG-P
1	E	15	SEP	CB-OG-P-O1P
1	E	15	SEP	CB-OG-P-O3P
1	F	15	SEP	N-CA-CB-OG
1	F	15	SEP	C-CA-CB-OG
1	A	15	SEP	N-CA-CB-OG
1	C	15	SEP	N-CA-CB-OG
1	E	15	SEP	N-CA-CB-OG
1	A	15	SEP	CB-OG-P-O3P
1	D	15	SEP	CB-OG-P-O3P
1	E	15	SEP	CB-OG-P-O2P

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	D	15	SEP	2	0
1	C	15	SEP	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

25 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	PO4	C	504	-	4,4,4	0.74	0	6,6,6	0.69	0
2	PO4	E	502	-	4,4,4	0.84	0	6,6,6	1.09	0
2	PO4	D	501	-	4,4,4	0.92	0	6,6,6	0.52	0
3	GOL	E	505	-	5,5,5	0.40	0	5,5,5	0.49	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	PO4	D	502	-	4,4,4	0.99	0	6,6,6	0.72	0
2	PO4	E	503	-	4,4,4	0.96	0	6,6,6	0.51	0
3	GOL	C	505	-	5,5,5	0.47	0	5,5,5	0.89	0
3	GOL	E	504	-	5,5,5	0.37	0	5,5,5	0.61	0
2	PO4	E	501	-	4,4,4	0.97	0	6,6,6	0.87	0
2	PO4	D	503	-	4,4,4	0.93	0	6,6,6	0.56	0
2	PO4	B	501	-	4,4,4	0.96	0	6,6,6	1.43	1 (16%)
2	PO4	A	502	-	4,4,4	1.03	0	6,6,6	1.31	1 (16%)
3	GOL	A	505	-	5,5,5	0.45	0	5,5,5	1.17	0
2	PO4	A	501	-	4,4,4	0.86	0	6,6,6	0.54	0
3	GOL	A	506	-	5,5,5	0.55	0	5,5,5	0.54	0
2	PO4	A	503	-	4,4,4	1.01	0	6,6,6	0.89	0
2	PO4	A	504	-	4,4,4	0.87	0	6,6,6	0.54	0
2	PO4	D	504	-	4,4,4	0.84	0	6,6,6	0.56	0
3	GOL	D	505	-	5,5,5	0.51	0	5,5,5	0.45	0
3	GOL	C	506	-	5,5,5	0.48	0	5,5,5	0.43	0
2	PO4	F	501	-	4,4,4	0.90	0	6,6,6	0.75	0
2	PO4	C	502	-	4,4,4	0.83	0	6,6,6	0.44	0
3	GOL	B	502	-	5,5,5	0.23	0	5,5,5	1.21	0
2	PO4	C	501	-	4,4,4	0.88	0	6,6,6	0.56	0
2	PO4	C	503	-	4,4,4	0.76	0	6,6,6	0.80	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	GOL	C	506	-	-	2/4/4/4	-
3	GOL	A	506	-	-	2/4/4/4	-
3	GOL	C	505	-	-	2/4/4/4	-
3	GOL	B	502	-	-	4/4/4/4	-
3	GOL	E	505	-	-	4/4/4/4	-
3	GOL	E	504	-	-	3/4/4/4	-
3	GOL	D	505	-	-	0/4/4/4	-
3	GOL	A	505	-	-	4/4/4/4	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	PO4	O4-P-O2	2.51	115.71	107.91
2	A	502	PO4	O3-P-O2	2.42	115.43	107.91

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	505	GOL	O1-C1-C2-C3
3	E	504	GOL	C1-C2-C3-O3
3	E	505	GOL	C1-C2-C3-O3
3	E	505	GOL	O2-C2-C3-O3
3	A	505	GOL	C1-C2-C3-O3
3	A	506	GOL	C1-C2-C3-O3
3	B	502	GOL	O1-C1-C2-C3
3	B	502	GOL	C1-C2-C3-O3
3	C	506	GOL	O1-C1-C2-C3
3	E	505	GOL	O1-C1-C2-C3
3	A	505	GOL	O1-C1-C2-O2
3	A	505	GOL	O2-C2-C3-O3
3	A	506	GOL	O2-C2-C3-O3
3	B	502	GOL	O2-C2-C3-O3
3	C	506	GOL	O1-C1-C2-O2
3	E	504	GOL	O2-C2-C3-O3
3	B	502	GOL	O1-C1-C2-O2
3	C	505	GOL	O2-C2-C3-O3
3	E	504	GOL	O1-C1-C2-O2
3	E	505	GOL	O1-C1-C2-O2
3	C	505	GOL	C1-C2-C3-O3

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	504	GOL	1	0
2	E	501	PO4	1	0
3	D	505	GOL	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	469/485 (96%)	-0.54	3 (0%) 85 86	23, 34, 56, 81	1 (0%)
1	B	471/485 (97%)	-0.29	5 (1%) 78 79	27, 41, 74, 102	2 (0%)
1	C	474/485 (97%)	-0.27	4 (0%) 82 84	19, 42, 80, 98	0
1	D	472/485 (97%)	-0.24	3 (0%) 85 86	25, 46, 74, 105	0
1	E	470/485 (96%)	-0.03	11 (2%) 61 61	23, 52, 97, 156	0
1	F	459/485 (94%)	0.82	84 (18%) 3 3	25, 59, 115, 146	0
All	All	2815/2910 (96%)	-0.10	110 (3%) 43 43	19, 44, 92, 156	3 (0%)

All (110) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	179	ILE	6.1
1	A	484	GLY	5.1
1	F	115	PRO	4.9
1	F	127	TRP	4.8
1	F	165	ILE	4.7
1	C	11	ALA	4.6
1	F	163	LYS	4.5
1	F	176	GLY	4.4
1	F	185	HIS	4.3
1	F	188	VAL	4.3
1	F	126	LEU	4.2
1	F	189	PRO	4.2
1	F	177	PHE	4.2
1	F	170	LEU	4.1
1	F	124	GLU	4.0
1	F	128	SER	4.0
1	F	202	ILE	3.9
1	F	75	THR	3.7
1	F	143	PRO	3.7

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	E	366	THR	3.6
1	E	17	ALA	3.6
1	F	173	TRP	3.6
1	F	190	MET	3.6
1	F	116	ASP	3.4
1	B	13	LYS	3.3
1	F	187	VAL	3.3
1	F	149	PRO	3.2
1	F	474	LEU	3.2
1	F	158	TYR	3.2
1	F	175	LYS	3.2
1	F	114	LEU	3.2
1	F	172	ARG	3.1
1	F	141	TRP	3.1
1	F	166	ASN	3.1
1	F	129	PHE	2.9
1	F	475	ALA	2.9
1	F	87	LEU	2.9
1	F	90	THR	2.9
1	C	13	LYS	2.9
1	F	159	PRO	2.9
1	D	13	LYS	2.8
1	F	201	PRO	2.8
1	F	137	PHE	2.8
1	F	164	LYS	2.8
1	F	195	ILE	2.7
1	F	125	GLN	2.7
1	F	472	LYS	2.7
1	F	253	LEU	2.7
1	E	21	ALA	2.7
1	F	407	GLY	2.7
1	F	198	LEU	2.6
1	F	161	SER	2.6
1	F	191	LEU	2.6
1	F	252	LEU	2.6
1	F	70	VAL	2.6
1	E	59	LYS	2.6
1	F	197	LYS	2.6
1	F	92	LEU	2.5
1	F	476	ALA	2.5
1	A	16	MET	2.5
1	F	194	GLN	2.5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	F	131	ALA	2.5
1	F	144	ASP	2.5
1	F	196	GLU	2.5
1	F	174	THR	2.5
1	F	169	VAL	2.5
1	E	16	MET	2.5
1	A	17	ALA	2.4
1	D	147	SER	2.4
1	F	167	SER	2.4
1	F	259	PRO	2.4
1	E	327	PHE	2.4
1	F	200	ILE	2.4
1	F	72	GLU	2.4
1	F	178	ASP	2.4
1	E	354	PHE	2.4
1	F	146	VAL	2.4
1	F	154	PHE	2.4
1	F	480	VAL	2.4
1	F	482	ILE	2.4
1	E	18	ASP	2.3
1	F	147	SER	2.3
1	F	204	VAL	2.3
1	F	134	LEU	2.3
1	F	199	ASN	2.3
1	D	14	GLY	2.3
1	F	88	GLY	2.3
1	F	77	LYS	2.3
1	F	162	GLN	2.3
1	B	14	GLY	2.2
1	F	171	GLN	2.2
1	C	251	GLY	2.2
1	B	17	ALA	2.2
1	F	59	LYS	2.2
1	B	484	GLY	2.2
1	B	113	ARG	2.1
1	E	24	MET	2.1
1	F	138	VAL	2.1
1	F	463	ALA	2.1
1	F	193	GLU	2.1
1	F	91	ASN	2.1
1	C	14	GLY	2.1
1	F	133	CYS	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	E	409	LYS	2.1
1	F	160	ALA	2.1
1	E	277	LEU	2.1
1	F	246	ILE	2.1
1	F	112	TYR	2.0
1	F	105	PHE	2.0
1	F	226	LYS	2.0

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

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
1	SEP	E	15	10/11	0.50	0.12	103,126,142,144	0
1	SEP	C	15	10/11	0.72	0.15	83,95,111,115	0
1	SEP	F	15	10/11	0.76	0.15	69,75,84,89	0
1	SEP	A	15	10/11	0.79	0.16	69,85,92,97	0
1	SEP	B	15	10/11	0.82	0.13	68,81,87,91	0
1	SEP	D	15	10/11	0.89	0.11	64,73,79,85	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	PO4	D	504	5/5	0.65	0.11	61,62,66,84	5
2	PO4	D	503	5/5	0.72	0.20	51,51,57,63	5
2	PO4	C	501	5/5	0.76	0.10	59,59,73,80	5
2	PO4	C	504	5/5	0.78	0.16	52,54,67,84	5
2	PO4	D	501	5/5	0.80	0.13	55,57,79,85	5

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	PO4	E	502	5/5	0.81	0.11	44,51,64,67	5
3	GOL	E	505	6/6	0.81	0.16	51,68,74,76	0
2	PO4	E	503	5/5	0.82	0.14	44,52,67,73	5
2	PO4	A	504	5/5	0.84	0.12	41,58,62,73	5
3	GOL	D	505	6/6	0.86	0.13	41,45,47,52	0
2	PO4	A	501	5/5	0.86	0.09	46,53,65,71	5
3	GOL	B	502	6/6	0.87	0.19	42,49,59,60	0
3	GOL	C	506	6/6	0.87	0.14	45,65,72,79	0
2	PO4	C	502	5/5	0.89	0.09	47,52,71,78	5
3	GOL	A	506	6/6	0.89	0.14	41,60,71,72	0
3	GOL	C	505	6/6	0.90	0.14	33,43,55,73	0
3	GOL	E	504	6/6	0.93	0.12	38,49,68,74	0
2	PO4	F	501	5/5	0.93	0.08	43,61,69,78	0
3	GOL	A	505	6/6	0.94	0.10	29,35,40,40	0
2	PO4	C	503	5/5	0.94	0.08	40,46,58,65	0
2	PO4	A	503	5/5	0.95	0.08	41,55,60,64	0
2	PO4	A	502	5/5	0.95	0.08	36,39,44,46	0
2	PO4	D	502	5/5	0.95	0.08	47,58,64,65	0
2	PO4	B	501	5/5	0.95	0.08	45,48,55,57	0
2	PO4	E	501	5/5	0.97	0.06	42,43,67,75	0

6.5 Other polymers

There are no such residues in this entry.