



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:20 PM UTC

PDB ID : 4KBR / pdb_00004kbr
Title : Crystal structure of mouse Ceramide-1-phosphate transfer protein (apo-form)
Authors : Simanshu, D.K.; Brown, R.E.; Patel, D.J.
Deposited on : 2013-04-23
Resolution : 2.55 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

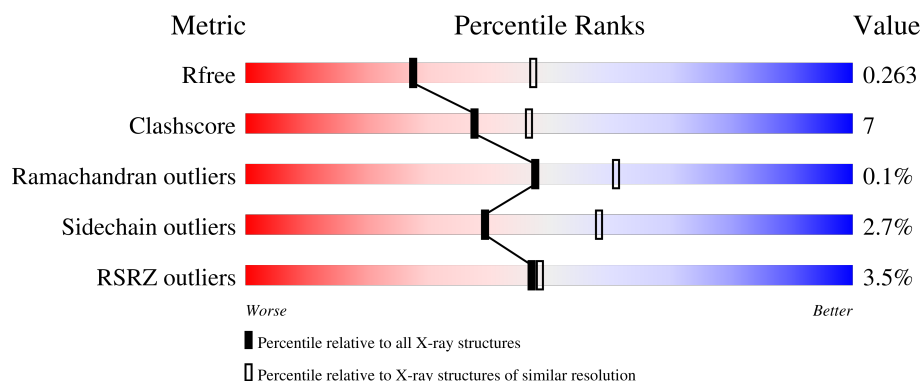
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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	217	
1	B	217	
1	C	217	
1	D	217	
1	E	217	

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Mol	Chain	Length	Quality of chain
1	F	217	7% 76% 17% 7%
1	G	217	% 79% 14% 6%
1	H	217	6% 76% 17% 6%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13204 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 Glycolipid transfer protein domain-containing protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	206	Total	C	N	O	S	0	0	0
			1629	1040	280	299	10			
1	B	207	Total	C	N	O	S	0	0	0
			1619	1034	283	292	10			
1	C	205	Total	C	N	O	S	0	0	0
			1633	1042	282	299	10			
1	D	202	Total	C	N	O	S	0	0	0
			1607	1028	276	293	10			
1	E	207	Total	C	N	O	S	0	0	0
			1658	1058	288	302	10			
1	F	202	Total	C	N	O	S	0	0	0
			1584	1011	273	290	10			
1	G	203	Total	C	N	O	S	0	0	0
			1620	1035	280	296	9			
1	H	203	Total	C	N	O	S	0	0	0
			1560	997	268	285	10			

There are 16 discrepancies between the modelled and reference sequences:

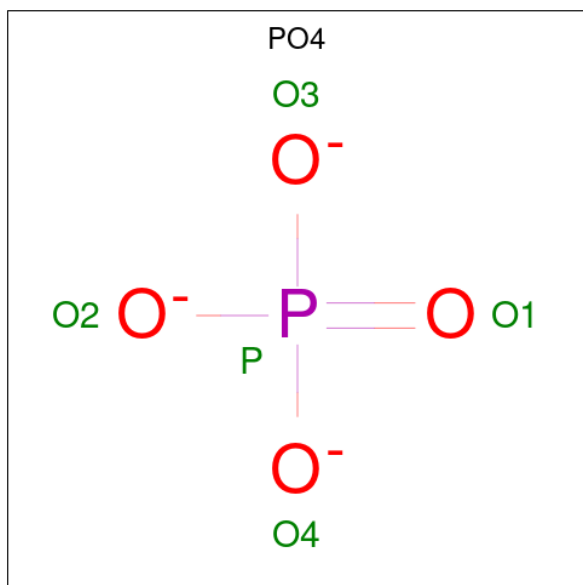
Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP Q8BS40
A	77	THR	ALA	conflict	UNP Q8BS40
B	0	SER	-	expression tag	UNP Q8BS40
B	77	THR	ALA	conflict	UNP Q8BS40
C	0	SER	-	expression tag	UNP Q8BS40
C	77	THR	ALA	conflict	UNP Q8BS40
D	0	SER	-	expression tag	UNP Q8BS40
D	77	THR	ALA	conflict	UNP Q8BS40
E	0	SER	-	expression tag	UNP Q8BS40
E	77	THR	ALA	conflict	UNP Q8BS40
F	0	SER	-	expression tag	UNP Q8BS40
F	77	THR	ALA	conflict	UNP Q8BS40
G	0	SER	-	expression tag	UNP Q8BS40

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Chain	Residue	Modelled	Actual	Comment	Reference
G	77	THR	ALA	conflict	UNP Q8BS40
H	0	SER	-	expression tag	UNP Q8BS40
H	77	THR	ALA	conflict	UNP Q8BS40

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



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

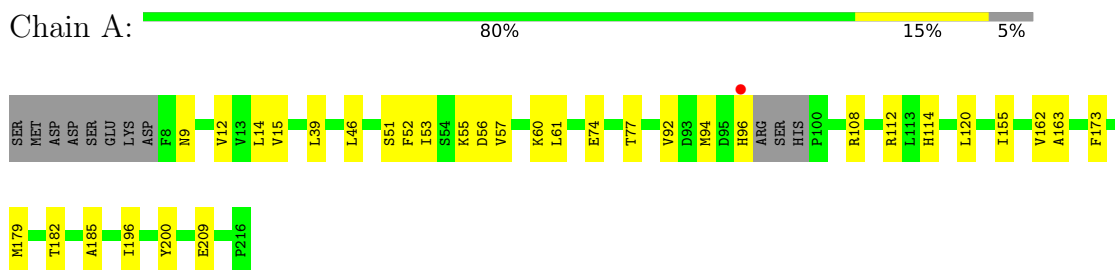
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	54	Total 54	O 54	0	0
3	B	38	Total 38	O 38	0	0
3	C	46	Total 46	O 46	0	0
3	D	15	Total 15	O 15	0	0
3	E	35	Total 35	O 35	0	0
3	F	15	Total 15	O 15	0	0
3	G	29	Total 29	O 29	0	0
3	H	22	Total 22	O 22	0	0

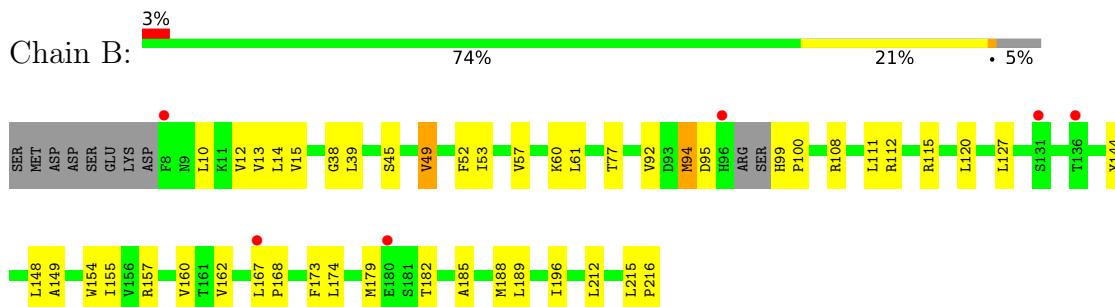
3 Residue-property plots [i](#)

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.

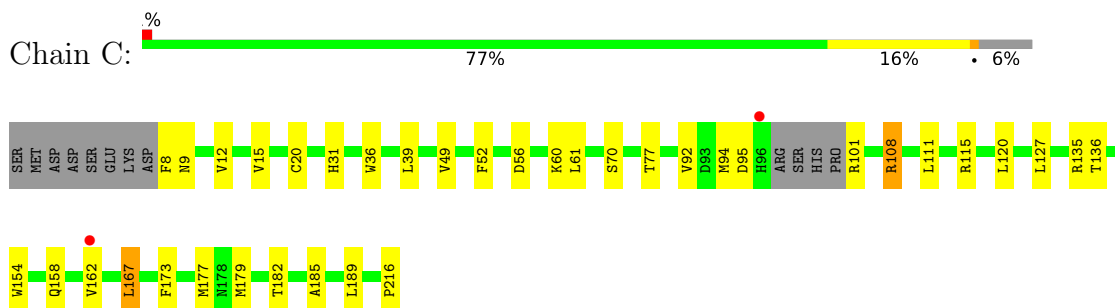
- Molecule 1: Glycolipid transfer protein domain-containing protein 1



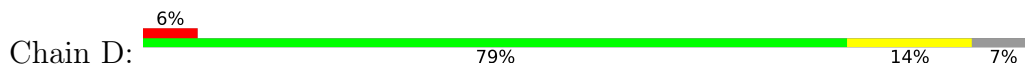
- Molecule 1: Glycolipid transfer protein domain-containing protein 1

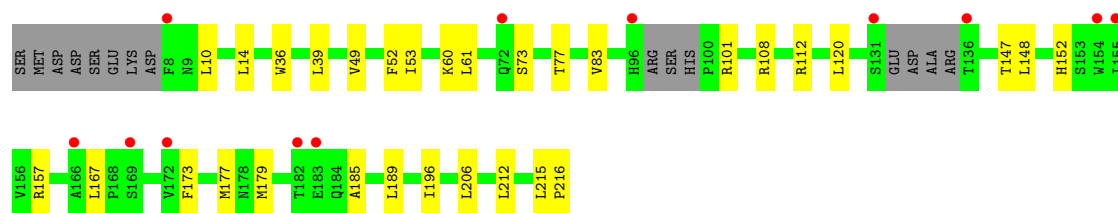


- Molecule 1: Glycolipid transfer protein domain-containing protein 1

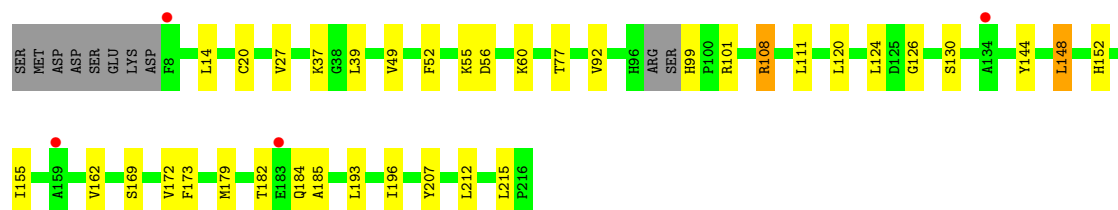
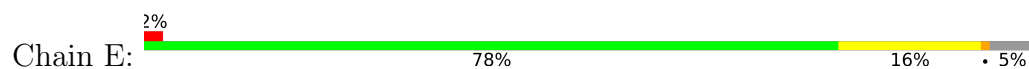


- Molecule 1: Glycolipid transfer protein domain-containing protein 1

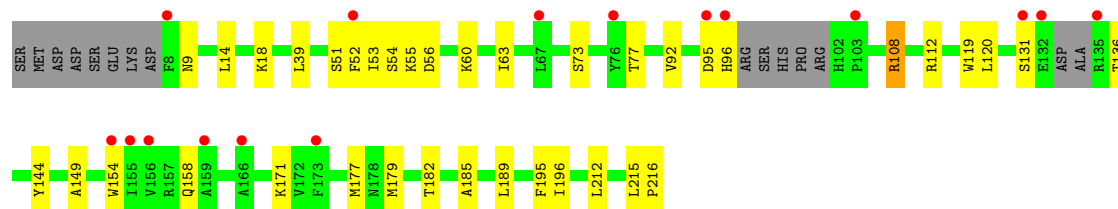
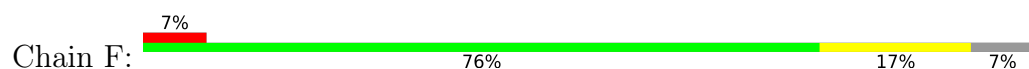




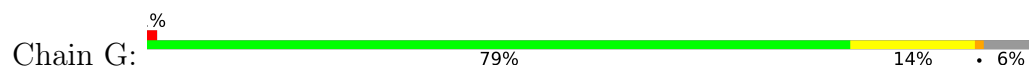
- Molecule 1: Glycolipid transfer protein domain-containing protein 1



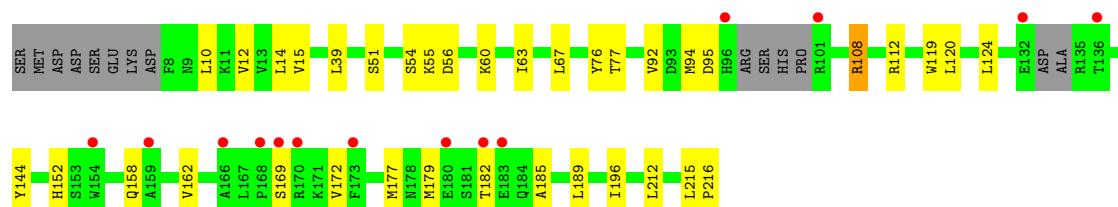
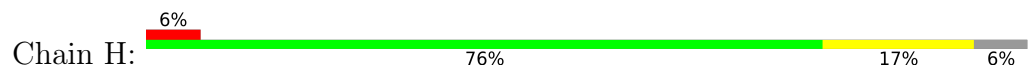
- Molecule 1: Glycolipid transfer protein domain-containing protein 1



- Molecule 1: Glycolipid transfer protein domain-containing protein 1



- Molecule 1: Glycolipid transfer protein domain-containing protein 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	82.00Å 83.88Å 268.72Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.51 – 2.55 39.51 – 2.55	Depositor EDS
% Data completeness (in resolution range)	93.3 (39.51-2.55) 93.4 (39.51-2.55)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.51 (at 2.54Å)	Xtriage
Refinement program	PHENIX 1.8.2_1309	Depositor
R, R_{free}	0.201 , 0.271 (Not available) , 0.263	Depositor DCC
R_{free} test set	2916 reflections (4.73%)	wwPDB-VP
Wilson B-factor (Å ²)	45.8	Xtriage
Anisotropy	0.574	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 47.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.028 for k,h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	13204	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.12% 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: 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	0.61	0/1665	0.85	0/2256
1	B	0.59	0/1655	0.87	1/2246 (0.0%)
1	C	0.57	0/1668	0.93	4/2258 (0.2%)
1	D	0.54	0/1642	0.87	1/2223 (0.0%)
1	E	0.57	0/1695	0.88	1/2294 (0.0%)
1	F	0.51	0/1615	0.87	2/2185 (0.1%)
1	G	0.50	0/1654	0.87	2/2238 (0.1%)
1	H	0.52	0/1590	0.86	0/2156
All	All	0.55	0/13184	0.88	11/17856 (0.1%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	167	LEU	CA-C-N	7.53	128.55	120.11
1	C	167	LEU	C-N-CA	7.53	128.55	120.11
1	C	70	SER	CA-C-N	-7.25	112.33	119.87
1	C	70	SER	C-N-CA	-7.25	112.33	119.87
1	G	167	LEU	CA-C-N	6.78	127.71	120.11
1	G	167	LEU	C-N-CA	6.78	127.71	120.11
1	E	148	LEU	N-CA-C	5.78	117.58	111.28
1	F	136	THR	N-CA-C	5.77	117.25	111.07
1	B	148	LEU	N-CA-C	5.46	117.23	111.28
1	F	73	SER	N-CA-C	5.34	117.79	111.33
1	D	73	SER	N-CA-C	5.15	117.56	111.33

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	1629	0	1602	23	0
1	B	1619	0	1592	31	0
1	C	1633	0	1619	25	0
1	D	1607	0	1593	15	0
1	E	1658	0	1650	29	0
1	F	1584	0	1573	23	0
1	G	1620	0	1618	19	0
1	H	1560	0	1526	25	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
2	D	5	0	0	0	0
2	E	5	0	0	0	0
2	F	5	0	0	0	0
2	G	5	0	0	0	0
2	H	5	0	0	1	0
3	A	54	0	0	2	0
3	B	38	0	0	0	0
3	C	46	0	0	2	0
3	D	15	0	0	0	0
3	E	35	0	0	2	0
3	F	15	0	0	1	0
3	G	29	0	0	0	0
3	H	22	0	0	1	0
All	All	13204	0	12773	179	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 7.

All (179) 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:39:LEU:HD13	1:D:120:LEU:HD23	1.66	0.76
1:E:99:HIS:N	3:E:415:HOH:O	2.21	0.74
1:F:14:LEU:HD22	1:F:196:ILE:HD11	1.70	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:14:LEU:HD22	1:A:196:ILE:HD11	1.69	0.73
1:F:51:SER:HA	1:F:54:SER:HB3	1.71	0.73
1:F:39:LEU:HD13	1:F:120:LEU:HD23	1.71	0.70
1:E:37:LYS:NZ	3:E:410:HOH:O	2.24	0.70
1:E:39:LEU:HD13	1:E:120:LEU:HD23	1.74	0.69
1:F:60:LYS:HE2	1:F:112:ARG:HG2	1.76	0.67
1:B:179:MET:HG3	1:B:185:ALA:HA	1.77	0.66
1:D:14:LEU:HD22	1:D:196:ILE:HD11	1.79	0.65
1:D:60:LYS:HE2	1:D:112:ARG:HG2	1.78	0.64
1:E:14:LEU:HD22	1:E:196:ILE:HD11	1.81	0.62
1:C:101:ARG:N	3:C:439:HOH:O	2.32	0.62
1:A:39:LEU:HD13	1:A:120:LEU:HD23	1.81	0.61
1:G:14:LEU:HD22	1:G:196:ILE:HD11	1.81	0.61
1:B:149:ALA:HB1	1:B:157:ARG:HD3	1.82	0.61
1:E:212:LEU:HD22	1:E:215:LEU:HD11	1.83	0.60
1:A:163:ALA:HB1	1:C:162:VAL:HG12	1.84	0.59
1:F:179:MET:HG3	1:F:185:ALA:HA	1.85	0.58
1:A:12:VAL:O	1:A:15:VAL:HG12	2.03	0.57
1:B:14:LEU:HD22	1:B:196:ILE:HD11	1.84	0.57
1:B:162:VAL:HG11	1:E:49:VAL:HA	1.86	0.57
1:F:60:LYS:HA	1:F:63:ILE:HD12	1.88	0.56
1:H:14:LEU:HD22	1:H:196:ILE:HD11	1.86	0.56
1:H:95:ASP:O	1:H:216:PRO:HG3	2.06	0.56
1:C:9:ASN:OD1	1:C:179:MET:HE1	2.06	0.55
1:C:36:TRP:HB3	1:C:61:LEU:HD11	1.89	0.55
1:E:39:LEU:HD22	1:E:120:LEU:HD21	1.88	0.55
1:B:52:PHE:CE2	1:B:160:VAL:HG22	2.42	0.55
1:C:177:MET:HE1	1:C:189:LEU:HD11	1.87	0.55
1:G:9:ASN:HB2	1:G:179:MET:CE	2.37	0.55
1:G:179:MET:HG3	1:G:185:ALA:HA	1.89	0.55
1:F:212:LEU:O	1:F:215:LEU:HG	2.06	0.54
1:H:179:MET:HG3	1:H:185:ALA:HA	1.89	0.54
1:A:179:MET:HG3	1:A:185:ALA:HA	1.89	0.53
1:C:179:MET:HG3	1:C:185:ALA:HA	1.90	0.53
1:E:179:MET:HG3	1:E:185:ALA:HA	1.89	0.53
1:A:9:ASN:HB2	1:A:179:MET:CE	2.38	0.53
1:H:39:LEU:HD13	1:H:120:LEU:HD23	1.89	0.53
1:D:179:MET:HG3	1:D:185:ALA:HA	1.91	0.53
1:E:169:SER:OG	1:E:172:VAL:HG23	2.08	0.53
1:F:56:ASP:OD1	1:F:60:LYS:NZ	2.42	0.53
1:H:56:ASP:OD1	1:H:60:LYS:NZ	2.42	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:12:VAL:O	1:H:15:VAL:HG12	2.09	0.52
1:E:144:TYR:OH	1:E:152:HIS:ND1	2.35	0.52
1:C:39:LEU:HD13	1:C:120:LEU:HD23	1.91	0.52
1:C:49:VAL:HG21	1:C:167:LEU:HD23	1.92	0.52
1:A:162:VAL:HG11	1:C:49:VAL:HA	1.92	0.52
1:B:95:ASP:O	1:B:216:PRO:HG3	2.11	0.51
1:H:60:LYS:HE2	1:H:112:ARG:HG2	1.91	0.51
1:G:177:MET:HE1	1:G:189:LEU:HD11	1.92	0.50
1:A:46:LEU:HD13	1:A:173:PHE:CE1	2.46	0.50
1:G:39:LEU:HD13	1:G:120:LEU:HD23	1.91	0.50
1:C:56:ASP:OD1	1:C:60:LYS:NZ	2.42	0.50
1:C:95:ASP:O	1:C:216:PRO:HG3	2.11	0.50
1:H:67:LEU:HD22	1:H:76:TYR:CZ	2.47	0.50
1:B:99:HIS:CG	1:B:100:PRO:HD2	2.47	0.49
1:A:96:HIS:O	3:A:438:HOH:O	2.20	0.49
1:E:179:MET:HB3	1:E:184:GLN:OE1	2.12	0.49
1:G:185:ALA:O	1:G:189:LEU:HG	2.11	0.49
1:E:126:GLY:O	1:E:130:SER:HB3	2.12	0.49
1:B:49:VAL:HA	1:E:162:VAL:HG11	1.94	0.49
1:G:158:GLN:O	1:G:162:VAL:HG23	2.13	0.49
1:B:14:LEU:HD21	1:B:39:LEU:HD21	1.95	0.48
1:H:54:SER:OG	3:H:410:HOH:O	2.20	0.48
1:B:155:ILE:HD12	1:E:52:PHE:HB2	1.96	0.47
1:F:52:PHE:CE2	1:F:53:ILE:HG13	2.49	0.47
1:H:144:TYR:OH	1:H:152:HIS:ND1	2.40	0.47
1:G:51:SER:O	1:G:55:LYS:HG3	2.14	0.47
1:B:154:TRP:CZ2	1:E:55:LYS:HB2	2.50	0.47
1:B:154:TRP:HA	1:B:157:ARG:NH2	2.30	0.47
1:F:95:ASP:O	1:F:216:PRO:HG3	2.15	0.47
1:F:96:HIS:HA	3:F:409:HOH:O	2.14	0.47
1:G:171:LYS:NZ	1:G:182:THR:HG21	2.30	0.47
1:B:39:LEU:HD13	1:B:120:LEU:HD23	1.95	0.47
1:G:9:ASN:OD1	1:G:179:MET:HE1	2.15	0.47
1:G:46:LEU:HD13	1:G:173:PHE:CZ	2.50	0.47
1:B:12:VAL:O	1:B:15:VAL:HG12	2.15	0.47
1:F:9:ASN:OD1	1:F:179:MET:HE1	2.15	0.47
1:B:154:TRP:HZ2	1:E:55:LYS:HB2	1.79	0.46
1:B:212:LEU:HD22	1:B:215:LEU:HD11	1.96	0.46
1:B:144:TYR:CZ	1:B:160:VAL:HG11	2.51	0.46
1:B:185:ALA:O	1:B:189:LEU:HG	2.15	0.46
1:E:108:ARG:HD3	1:E:108:ARG:C	2.41	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:52:PHE:CE2	1:D:53:ILE:HG13	2.51	0.46
1:H:94:MET:HE1	1:H:212:LEU:HD23	1.98	0.46
1:H:212:LEU:O	1:H:215:LEU:HG	2.15	0.46
1:D:177:MET:HE1	1:D:189:LEU:HD11	1.97	0.46
1:A:57:VAL:O	1:A:61:LEU:HG	2.16	0.46
1:H:51:SER:HB2	1:H:55:LYS:HE2	1.97	0.46
1:D:147:THR:OG1	1:D:148:LEU:N	2.48	0.46
1:H:10:LEU:O	1:H:14:LEU:HG	2.15	0.45
1:C:136:THR:HG21	1:C:173:PHE:CE2	2.51	0.45
1:A:120:LEU:HD12	1:A:120:LEU:HA	1.70	0.45
1:C:111:LEU:HD21	1:C:115:ARG:NH2	2.31	0.45
1:B:57:VAL:O	1:B:61:LEU:HG	2.16	0.45
1:C:185:ALA:O	1:C:189:LEU:HG	2.16	0.45
1:H:158:GLN:O	1:H:162:VAL:HG23	2.17	0.45
1:A:209:GLU:HG2	1:B:45:SER:O	2.17	0.45
1:F:108:ARG:NH2	1:F:216:PRO:OXT	2.41	0.45
1:D:83:VAL:HG21	1:D:206:LEU:HD21	1.98	0.44
1:A:56:ASP:OD1	1:A:60:LYS:NZ	2.45	0.44
1:E:101:ARG:HE	1:E:101:ARG:HB3	1.42	0.44
1:C:12:VAL:O	1:C:15:VAL:HG12	2.17	0.44
1:F:179:MET:HA	1:F:179:MET:HE2	1.98	0.44
1:A:155:ILE:HD12	1:C:52:PHE:HB2	1.98	0.44
1:F:177:MET:HE1	1:F:189:LEU:HD11	2.00	0.44
1:E:179:MET:HE2	1:E:179:MET:HA	1.99	0.44
1:H:51:SER:HA	1:H:54:SER:HB3	1.99	0.44
1:B:52:PHE:CE2	1:B:53:ILE:HG13	2.53	0.44
1:H:177:MET:HE1	1:H:189:LEU:HD11	2.00	0.44
1:B:60:LYS:HE2	1:B:112:ARG:HG2	2.00	0.43
1:A:9:ASN:OD1	1:A:179:MET:HE1	2.17	0.43
1:B:111:LEU:HD21	1:B:115:ARG:NH2	2.34	0.43
1:C:94:MET:HE3	1:C:94:MET:HB3	1.84	0.43
1:F:120:LEU:HD12	1:F:120:LEU:HA	1.82	0.43
1:F:119:TRP:CZ3	1:F:120:LEU:HD13	2.53	0.43
1:G:169:SER:OG	1:G:172:VAL:HG23	2.18	0.43
1:A:52:PHE:CE2	1:A:53:ILE:HG13	2.54	0.43
1:B:94:MET:HE3	1:B:94:MET:HB3	1.91	0.43
1:C:158:GLN:O	1:C:162:VAL:HG23	2.19	0.43
1:D:212:LEU:HD22	1:D:215:LEU:HD11	2.01	0.43
1:F:144:TYR:CE2	1:F:149:ALA:HB2	2.54	0.43
1:A:114:HIS:CE1	1:A:200:TYR:CE1	3.07	0.42
1:C:136:THR:HG21	1:C:173:PHE:HE2	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:120:LEU:HD12	1:D:120:LEU:HA	1.74	0.42
1:F:154:TRP:O	1:F:158:GLN:HG2	2.18	0.42
1:A:74:GLU:HG3	3:A:440:HOH:O	2.18	0.42
1:D:152:HIS:O	1:D:157:ARG:NH2	2.52	0.42
1:H:60:LYS:HA	1:H:63:ILE:HD12	2.00	0.42
1:C:8:PHE:N	3:C:432:HOH:O	2.52	0.42
1:F:51:SER:HB2	1:F:55:LYS:HE2	2.02	0.42
1:H:124:LEU:HD23	1:H:124:LEU:HA	1.86	0.42
1:A:94:MET:HE3	1:A:94:MET:HB3	1.76	0.42
1:B:52:PHE:HD2	1:E:155:ILE:HD11	1.84	0.42
1:D:10:LEU:HA	1:D:10:LEU:HD12	1.82	0.42
1:E:148:LEU:HA	1:E:148:LEU:HD23	1.84	0.42
1:B:13:VAL:HG13	1:B:38:GLY:HA3	2.02	0.42
1:D:36:TRP:HB3	1:D:61:LEU:HD11	2.02	0.42
1:G:120:LEU:HD12	1:G:120:LEU:HA	1.80	0.42
1:D:49:VAL:HG11	1:D:167:LEU:HD23	2.00	0.42
1:A:9:ASN:HB2	1:A:179:MET:HE1	2.01	0.42
1:A:51:SER:HB2	1:A:55:LYS:HE2	2.01	0.42
1:C:108:ARG:C	1:C:108:ARG:HD3	2.45	0.42
1:E:111:LEU:HB2	1:E:207:TYR:CE2	2.54	0.42
1:H:108:ARG:NH1	2:H:301:PO4:O2	2.45	0.42
1:G:179:MET:HB3	1:G:184:GLN:OE1	2.19	0.41
1:C:20:CYS:HB3	1:C:31:HIS:HB2	2.01	0.41
1:E:56:ASP:OD1	1:E:60:LYS:NZ	2.52	0.41
1:F:18:LYS:HD2	1:F:195:PHE:CZ	2.55	0.41
1:C:179:MET:HA	1:C:179:MET:HE2	2.03	0.41
1:G:167:LEU:HA	1:G:168:PRO:HD3	1.79	0.41
1:A:114:HIS:CE1	1:A:200:TYR:CD1	3.08	0.41
1:B:174:LEU:HD23	1:B:174:LEU:HA	1.72	0.41
1:F:171:LYS:HD3	1:F:171:LYS:HA	1.87	0.41
1:B:162:VAL:CG1	1:E:49:VAL:HA	2.51	0.41
1:G:127:LEU:HD21	1:G:173:PHE:HB3	2.01	0.41
1:H:51:SER:HB2	1:H:55:LYS:NZ	2.36	0.41
1:F:51:SER:O	1:F:55:LYS:HG3	2.21	0.41
1:H:51:SER:HB2	1:H:55:LYS:CE	2.51	0.41
1:H:119:TRP:CZ3	1:H:120:LEU:HD13	2.56	0.41
1:C:154:TRP:CZ2	1:C:158:GLN:HG3	2.56	0.41
1:H:169:SER:OG	1:H:172:VAL:HG23	2.21	0.41
1:D:215:LEU:HA	1:D:216:PRO:HD2	1.85	0.41
1:E:20:CYS:SG	1:E:27:VAL:HG13	2.60	0.41
1:B:167:LEU:HA	1:B:168:PRO:HD3	1.89	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:127:LEU:HD21	1:C:173:PHE:HB3	2.02	0.40
1:E:120:LEU:HD12	1:E:120:LEU:HA	1.81	0.40
1:G:60:LYS:HE2	1:G:112:ARG:HG2	2.02	0.40
1:H:51:SER:O	1:H:55:LYS:HG3	2.21	0.40
1:A:60:LYS:HE2	1:A:112:ARG:HG2	2.03	0.40
1:B:127:LEU:HA	1:B:127:LEU:HD12	1.92	0.40
1:E:124:LEU:HD23	1:E:124:LEU:HA	1.87	0.40
1:E:111:LEU:HB2	1:E:207:TYR:HE2	1.86	0.40
1:E:193:LEU:HD23	1:E:193:LEU:HA	1.81	0.40
1:G:10:LEU:O	1:G:14:LEU:HG	2.22	0.40
1:G:212:LEU:HD22	1:G:215:LEU:HD11	2.03	0.40
1:B:10:LEU:HB3	1:B:188:MET:CE	2.52	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	202/217 (93%)	198 (98%)	4 (2%)	0	100	100
1	B	203/217 (94%)	199 (98%)	4 (2%)	0	100	100
1	C	201/217 (93%)	198 (98%)	3 (2%)	0	100	100
1	D	196/217 (90%)	194 (99%)	2 (1%)	0	100	100
1	E	203/217 (94%)	200 (98%)	3 (2%)	0	100	100
1	F	196/217 (90%)	192 (98%)	3 (2%)	1 (0%)	24	34
1	G	199/217 (92%)	196 (98%)	3 (2%)	0	100	100
1	H	197/217 (91%)	194 (98%)	3 (2%)	0	100	100
All	All	1597/1736 (92%)	1571 (98%)	25 (2%)	1 (0%)	48	61

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	131	SER

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	178/194 (92%)	174 (98%)	4 (2%)	45	64
1	B	175/194 (90%)	168 (96%)	7 (4%)	28	42
1	C	180/194 (93%)	175 (97%)	5 (3%)	38	57
1	D	178/194 (92%)	174 (98%)	4 (2%)	45	64
1	E	184/194 (95%)	179 (97%)	5 (3%)	39	58
1	F	174/194 (90%)	170 (98%)	4 (2%)	44	62
1	G	179/194 (92%)	174 (97%)	5 (3%)	38	57
1	H	167/194 (86%)	163 (98%)	4 (2%)	43	61
All	All	1415/1552 (91%)	1377 (97%)	38 (3%)	39	58

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	77	THR
1	A	92	VAL
1	A	108	ARG
1	A	182	THR
1	B	49	VAL
1	B	77	THR
1	B	92	VAL
1	B	94	MET
1	B	108	ARG
1	B	173	PHE
1	B	182	THR
1	C	77	THR
1	C	92	VAL
1	C	108	ARG
1	C	135	ARG

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Mol	Chain	Res	Type
1	C	182	THR
1	D	77	THR
1	D	101	ARG
1	D	108	ARG
1	D	173	PHE
1	E	77	THR
1	E	92	VAL
1	E	108	ARG
1	E	173	PHE
1	E	182	THR
1	F	77	THR
1	F	92	VAL
1	F	108	ARG
1	F	182	THR
1	G	77	THR
1	G	92	VAL
1	G	108	ARG
1	G	135	ARG
1	G	182	THR
1	H	77	THR
1	H	92	VAL
1	H	108	ARG
1	H	182	THR

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

Mol	Chain	Res	Type
1	B	121	GLN
1	C	75	HIS
1	C	118	HIS
1	C	121	GLN
1	D	31	HIS
1	D	75	HIS
1	D	118	HIS
1	E	75	HIS
1	F	31	HIS
1	F	75	HIS
1	F	89	ASN
1	G	75	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

8 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	G	301	-	4,4,4	0.85	0	6,6,6	0.73	0
2	PO4	A	301	-	4,4,4	0.96	0	6,6,6	0.50	0
2	PO4	B	301	-	4,4,4	1.01	0	6,6,6	0.69	0
2	PO4	E	301	-	4,4,4	0.95	0	6,6,6	0.44	0
2	PO4	D	301	-	4,4,4	1.05	0	6,6,6	0.63	0
2	PO4	C	301	-	4,4,4	1.09	0	6,6,6	0.47	0
2	PO4	F	301	-	4,4,4	0.97	0	6,6,6	0.45	0
2	PO4	H	301	-	4,4,4	0.94	0	6,6,6	0.51	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	H	301	PO4	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	206/217 (94%)	-0.30	1 (0%) 87 89	29, 42, 64, 86	0
1	B	207/217 (95%)	-0.09	6 (2%) 53 55	29, 48, 79, 103	0
1	C	205/217 (94%)	-0.22	2 (0%) 79 80	32, 42, 71, 98	0
1	D	202/217 (93%)	0.26	12 (5%) 28 27	32, 54, 98, 123	0
1	E	207/217 (95%)	0.00	4 (1%) 66 66	31, 49, 76, 109	0
1	F	202/217 (93%)	0.54	16 (7%) 18 18	34, 68, 109, 132	0
1	G	203/217 (93%)	0.14	3 (1%) 72 73	33, 56, 81, 102	0
1	H	203/217 (93%)	0.38	14 (6%) 23 22	36, 58, 104, 122	0
All	All	1635/1736 (94%)	0.09	58 (3%) 47 48	29, 51, 96, 132	0

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	132	GLU	6.7
1	H	96	HIS	4.7
1	D	183	GLU	3.9
1	B	96	HIS	3.8
1	H	132	GLU	3.8
1	H	182	THR	3.7
1	F	135	ARG	3.6
1	F	96	HIS	3.5
1	F	155	ILE	3.4
1	E	8	PHE	3.4
1	H	170	ARG	3.4
1	H	136	THR	3.2
1	D	96	HIS	3.2
1	G	7	ASP	3.1
1	H	169	SER	3.1
1	B	136	THR	3.1

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Mol	Chain	Res	Type	RSRZ
1	D	136	THR	3.0
1	C	96	HIS	3.0
1	F	159	ALA	2.9
1	D	182	THR	2.9
1	H	183	GLU	2.7
1	D	169	SER	2.7
1	A	96	HIS	2.7
1	F	52	PHE	2.6
1	F	173	PHE	2.6
1	H	159	ALA	2.6
1	B	167	LEU	2.5
1	D	154	TRP	2.5
1	G	52	PHE	2.5
1	E	183	GLU	2.4
1	H	101	ARG	2.4
1	F	156	VAL	2.4
1	D	72	GLN	2.4
1	E	159	ALA	2.4
1	F	154	TRP	2.4
1	H	166	ALA	2.4
1	B	8	PHE	2.4
1	B	180	GLU	2.4
1	H	180	GLU	2.4
1	F	131	SER	2.3
1	F	166	ALA	2.3
1	D	131	SER	2.2
1	D	166	ALA	2.2
1	G	76	TYR	2.2
1	H	168	PRO	2.2
1	H	173	PHE	2.2
1	F	8	PHE	2.2
1	B	131	SER	2.1
1	F	103	PRO	2.1
1	D	172	VAL	2.1
1	F	67	LEU	2.1
1	H	154	TRP	2.1
1	D	8	PHE	2.0
1	E	134	ALA	2.0
1	D	155	ILE	2.0
1	F	76	TYR	2.0
1	C	162	VAL	2.0
1	F	95	ASP	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	F	301	5/5	0.81	0.12	80,90,99,105	0
2	PO4	H	301	5/5	0.87	0.09	82,83,87,88	0
2	PO4	G	301	5/5	0.88	0.14	72,73,78,79	0
2	PO4	E	301	5/5	0.90	0.09	64,66,69,87	0
2	PO4	A	301	5/5	0.94	0.07	62,64,69,71	0
2	PO4	D	301	5/5	0.94	0.06	67,69,79,81	0
2	PO4	C	301	5/5	0.95	0.08	50,55,66,67	0
2	PO4	B	301	5/5	0.96	0.07	47,51,60,62	0

6.5 Other polymers [i](#)

There are no such residues in this entry.