



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 15, 2026 – 12:42 PM UTC

PDB ID : 3T0F / pdb_00003t0f
Title : IspH:HMBPP (substrate) structure of the E126D mutant
Authors : Span, I.; Graewert, T.; Bacher, A.; Eisenreich, W.; Groll, M.
Deposited on : 2011-07-20
Resolution : 1.90 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
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
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

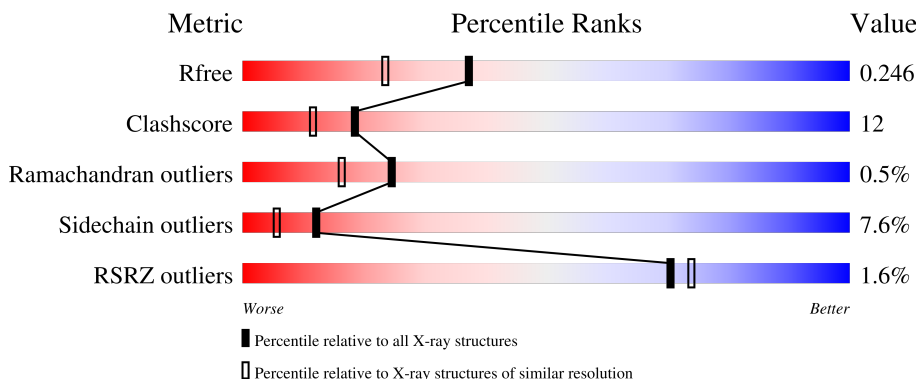
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	328	
1	B	328	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4995 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 4-hydroxy-3-methylbut-2-enyl diphosphate reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	314	Total	C	N	O	S	0	0	0
			2422	1509	438	465	10			
1	B	304	Total	C	N	O	S	0	0	0
			2345	1462	427	446	10			

There are 26 discrepancies between the modelled and reference sequences:

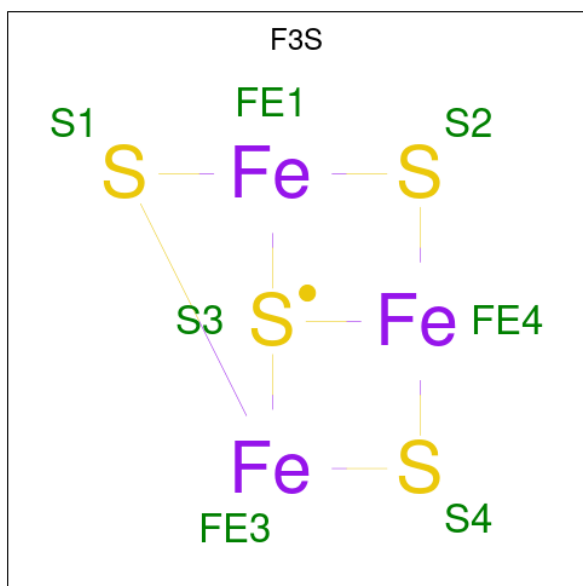
Chain	Residue	Modelled	Actual	Comment	Reference
A	-12	MET	-	expression tag	UNP P62623
A	-11	ARG	-	expression tag	UNP P62623
A	-10	GLY	-	expression tag	UNP P62623
A	-9	SER	-	expression tag	UNP P62623
A	-8	HIS	-	expression tag	UNP P62623
A	-7	HIS	-	expression tag	UNP P62623
A	-6	HIS	-	expression tag	UNP P62623
A	-5	HIS	-	expression tag	UNP P62623
A	-4	HIS	-	expression tag	UNP P62623
A	-3	HIS	-	expression tag	UNP P62623
A	-2	GLY	-	expression tag	UNP P62623
A	-1	SER	-	expression tag	UNP P62623
A	126	ASP	GLU	engineered mutation	UNP P62623
B	-12	MET	-	expression tag	UNP P62623
B	-11	ARG	-	expression tag	UNP P62623
B	-10	GLY	-	expression tag	UNP P62623
B	-9	SER	-	expression tag	UNP P62623
B	-8	HIS	-	expression tag	UNP P62623
B	-7	HIS	-	expression tag	UNP P62623
B	-6	HIS	-	expression tag	UNP P62623
B	-5	HIS	-	expression tag	UNP P62623
B	-4	HIS	-	expression tag	UNP P62623
B	-3	HIS	-	expression tag	UNP P62623
B	-2	GLY	-	expression tag	UNP P62623
B	-1	SER	-	expression tag	UNP P62623

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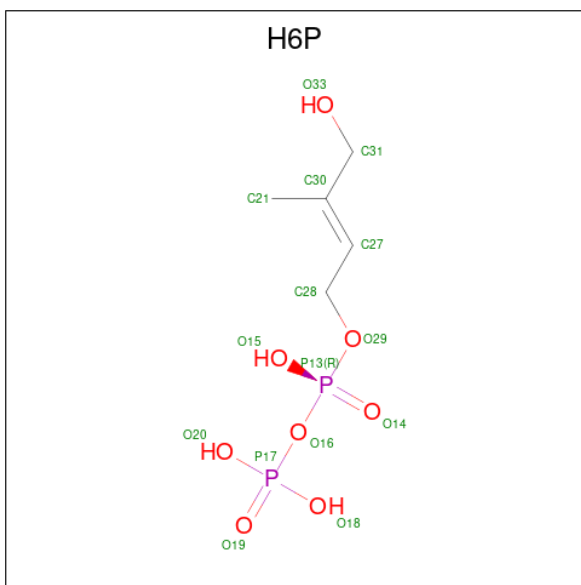
Chain	Residue	Modelled	Actual	Comment	Reference
B	126	ASP	GLU	engineered mutation	UNP P62623

- Molecule 2 is FE3-S4 CLUSTER (CCD ID: F3S) (formula: Fe_3S_4).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	Fe	S	0	0
			7	3	4		
2	B	1	Total	Fe	S	0	0
			7	3	4		

- Molecule 3 is (2E)-4-hydroxy-3-methylbut-2-en-1-yl trihydrogen diphosphate (CCD ID: H6P) (formula: $\text{C}_5\text{H}_{12}\text{O}_8\text{P}_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	P	0	0
			15	5	8	2		
3	B	1	Total	C	O	P	0	0
			15	5	8	2		

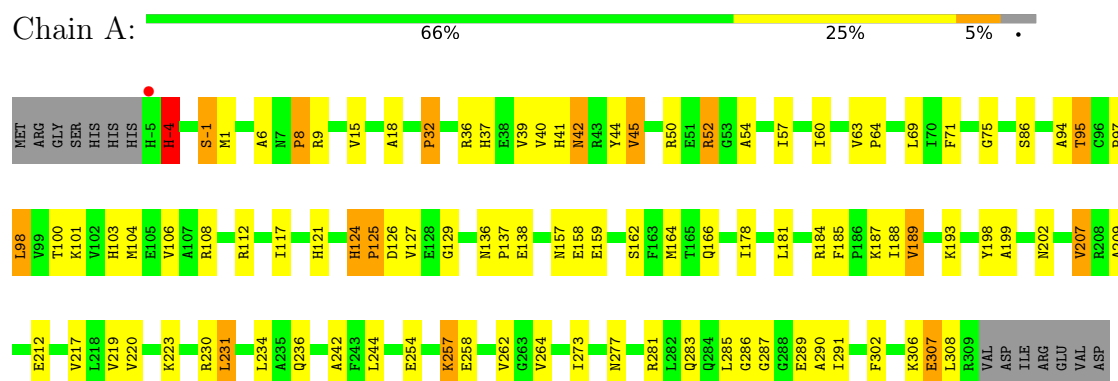
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	111	Total	O	0	0
			111	111		
4	B	73	Total	O	0	0
			73	73		

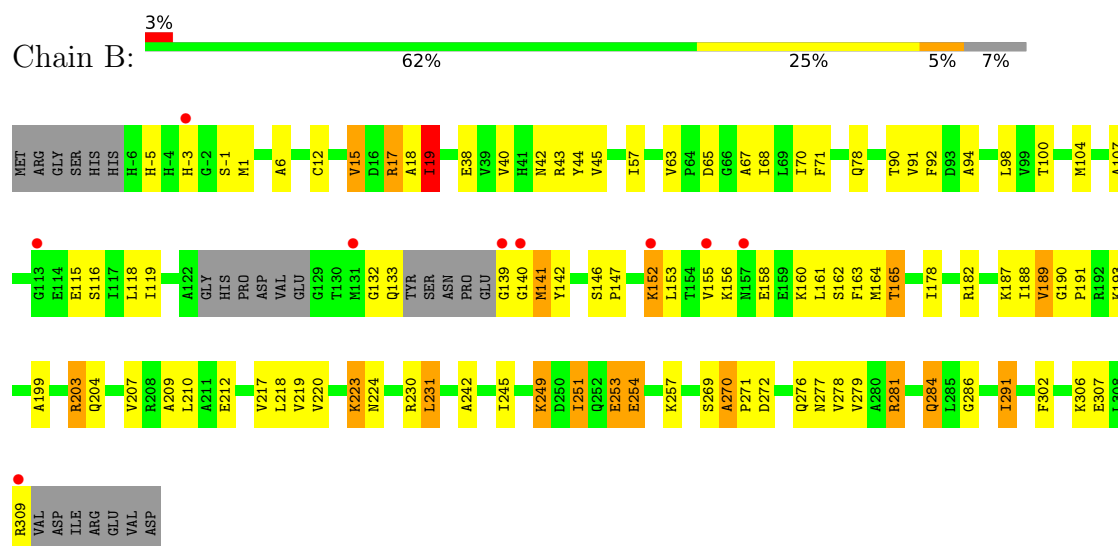
3 Residue-property plots [i](#)

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: 4-hydroxy-3-methylbut-2-enyl diphosphate reductase



- Molecule 1: 4-hydroxy-3-methylbut-2-enyl diphosphate reductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	64.85Å 80.18Å 114.51Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 1.90 10.00 – 1.90	Depositor EDS
% Data completeness (in resolution range)	100.0 (10.00-1.90) 98.4 (10.00-1.90)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.56 (at 1.90Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.201 , 0.242 0.213 , 0.246	Depositor DCC
R_{free} test set	2371 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	15.9	Xtriage
Anisotropy	0.879	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.46 , 41.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4995	wwPDB-VP
Average B, all atoms (Å ²)	10.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.88% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: F3S, H6P

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	1.58	20/2460 (0.8%)	1.47	17/3330 (0.5%)
1	B	1.57	16/2378 (0.7%)	1.45	21/3213 (0.7%)
All	All	1.57	36/4838 (0.7%)	1.46	38/6543 (0.6%)

All (36) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	220	VAL	CA-CB	8.34	1.64	1.54
1	B	278	VAL	CA-CB	7.69	1.63	1.54
1	B	15	VAL	C-O	7.62	1.32	1.24
1	A	39	VAL	N-CA	6.97	1.54	1.46
1	B	19	ILE	CA-C	-6.89	1.43	1.52
1	A	75	GLY	N-CA	6.88	1.52	1.45
1	A	103	HIS	N-CA	6.61	1.54	1.46
1	B	43	ARG	C-O	-6.59	1.15	1.24
1	B	12	CYS	CA-C	6.21	1.61	1.53
1	B	219	VAL	CA-CB	6.11	1.61	1.54
1	B	1	MET	C-O	-6.10	1.16	1.24
1	A	286	GLY	N-CA	6.04	1.53	1.45
1	A	41	HIS	C-O	5.95	1.31	1.24
1	B	1	MET	N-CA	5.87	1.53	1.46
1	B	91	VAL	CA-CB	5.86	1.61	1.54
1	A	18	ALA	CA-CB	-5.83	1.44	1.53
1	A	57	ILE	CA-CB	5.81	1.63	1.54
1	B	107	ALA	CA-CB	5.81	1.62	1.53
1	A	264	VAL	CA-CB	5.80	1.63	1.54
1	B	220	VAL	CA-CB	5.75	1.60	1.54
1	B	251	ILE	CA-C	-5.71	1.46	1.52
1	A	219	VAL	CA-C	-5.70	1.45	1.52
1	A	283	GLN	C-O	5.56	1.30	1.24
1	A	6	ALA	CA-CB	5.50	1.62	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	8	PRO	CG-CD	5.39	1.66	1.51
1	B	286	GLY	N-CA	5.39	1.52	1.45
1	A	37	HIS	N-CA	5.38	1.53	1.46
1	A	45	VAL	CB-CG1	5.31	1.70	1.52
1	A	54	ALA	N-CA	5.21	1.52	1.46
1	A	290	ALA	CA-CB	-5.18	1.45	1.53
1	A	42	ASN	CG-OD1	5.18	1.33	1.23
1	B	199	ALA	CA-CB	5.16	1.61	1.53
1	B	17	ARG	C-O	5.12	1.30	1.24
1	A	-1	SER	C-N	5.10	1.40	1.33
1	B	57	ILE	C-O	5.03	1.29	1.23
1	A	283	GLN	N-CA	5.02	1.52	1.46

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	-4	HIS	N-CA-C	11.75	131.56	114.16
1	A	-1	SER	N-CA-C	10.61	122.54	110.97
1	A	-1	SER	CA-C-N	8.58	134.16	121.72
1	A	-1	SER	C-N-CA	8.58	134.16	121.72
1	B	291	ILE	CA-C-N	-8.00	111.35	119.99
1	B	291	ILE	C-N-CA	-8.00	111.35	119.99
1	A	124	HIS	CA-C-N	6.66	128.16	119.84
1	A	124	HIS	C-N-CA	6.66	128.16	119.84
1	A	9	ARG	N-CA-C	-6.64	98.97	109.07
1	B	270	ALA	N-CA-C	6.58	117.94	109.72
1	B	203	ARG	N-CA-C	6.42	118.27	111.28
1	A	262	VAL	CA-C-N	-6.41	116.38	122.66
1	A	262	VAL	C-N-CA	-6.41	116.38	122.66
1	B	1	MET	N-CA-C	6.39	119.35	109.07
1	B	291	ILE	N-CA-C	6.38	114.92	107.77
1	A	231	LEU	N-CA-C	6.20	118.03	111.28
1	B	277	ASN	N-CA-C	5.89	117.78	111.36
1	B	63	VAL	CA-C-N	5.85	125.80	119.78
1	B	63	VAL	C-N-CA	5.85	125.80	119.78
1	B	38	GLU	N-CA-C	-5.78	101.72	110.28
1	A	277	ASN	N-CA-C	5.76	117.55	111.28
1	B	284	GLN	N-CA-CB	5.73	119.20	110.14
1	B	18	ALA	N-CA-C	5.65	117.44	111.28
1	A	291	ILE	CB-CA-C	-5.55	104.47	110.85
1	A	75	GLY	N-CA-C	5.54	119.66	112.68
1	A	188	ILE	N-CA-C	-5.53	101.92	109.55

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	224	ASN	CA-C-N	-5.50	115.69	122.84
1	B	224	ASN	C-N-CA	-5.50	115.69	122.84
1	B	94	ALA	N-CA-C	5.49	119.97	113.16
1	A	95	THR	CA-C-O	-5.43	115.77	120.88
1	B	284	GLN	CA-CB-CG	5.37	124.83	114.10
1	B	67	ALA	N-CA-C	5.25	117.30	110.53
1	A	98	LEU	CB-CG-CD1	5.24	126.41	110.70
1	A	285	LEU	N-CA-C	-5.15	106.59	112.92
1	B	71	PHE	N-CA-C	-5.09	102.75	110.28
1	B	204	GLN	CA-C-N	5.06	127.02	120.44
1	B	204	GLN	C-N-CA	5.06	127.02	120.44
1	B	279	VAL	N-CA-C	5.05	116.38	110.62

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	2422	0	2422	53	1
1	B	2345	0	2358	61	0
2	A	7	0	0	0	0
2	B	7	0	0	1	0
3	A	15	0	9	0	0
3	B	15	0	9	1	0
4	A	111	0	0	16	0
4	B	73	0	0	14	0
All	All	4995	0	4798	115	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 12.

All (115) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:254:GLU:HA	1:B:257:LYS:HD2	1.43	1.00
1:A:166:GLN:HG2	4:A:401:HOH:O	1.61	0.99
1:B:139:GLY:HA3	4:B:385:HOH:O	1.65	0.96
1:A:52:ARG:HD2	4:A:355:HOH:O	1.66	0.94
1:B:153:LEU:HB3	4:B:349:HOH:O	1.68	0.92
1:A:258:GLU:HG2	4:A:388:HOH:O	1.70	0.91
1:B:253:GLU:H	1:B:253:GLU:CD	1.79	0.90
1:B:223:LYS:HD2	4:B:325:HOH:O	1.72	0.88
1:B:254:GLU:HA	1:B:257:LYS:CD	2.04	0.87
1:A:125:PRO:HB3	4:A:417:HOH:O	1.81	0.81
1:A:-4:HIS:HB2	1:A:289:GLU:OE2	1.82	0.79
1:A:126:ASP:HB3	1:A:166:GLN:HE21	1.53	0.73
1:A:42:ASN:HD22	1:A:45:VAL:H	1.35	0.72
1:B:65:ASP:CB	4:B:360:HOH:O	2.38	0.71
1:B:68:ILE:HD12	1:B:90:THR:HB	1.73	0.68
1:B:162:SER:HA	1:B:189:VAL:O	1.92	0.68
1:B:119:ILE:HD12	1:B:165:THR:HG22	1.76	0.68
1:A:-4:HIS:CE1	4:A:424:HOH:O	2.45	0.68
1:B:245:ILE:HD13	1:B:251:ILE:HD11	1.77	0.67
1:A:212:GLU:CB	4:A:367:HOH:O	2.42	0.67
1:B:118:LEU:HB2	1:B:141:MET:CE	2.25	0.67
1:A:101:LYS:HD3	4:A:346:HOH:O	1.95	0.66
1:B:146:SER:HB2	1:B:147:PRO:HD2	1.76	0.66
1:A:178:ILE:HD13	1:A:193:LYS:O	1.96	0.65
1:B:42:ASN:HD22	1:B:45:VAL:H	1.43	0.65
1:A:212:GLU:HB3	4:A:367:HOH:O	1.95	0.64
1:A:126:ASP:HB3	1:A:166:GLN:NE2	2.12	0.63
1:B:253:GLU:CD	1:B:253:GLU:N	2.53	0.63
1:B:147:PRO:HD2	4:B:359:HOH:O	1.99	0.62
1:B:223:LYS:CD	4:B:325:HOH:O	2.34	0.62
1:B:163:PHE:CZ	1:B:191:PRO:HD2	2.34	0.62
1:A:136:ASN:HD22	1:A:137:PRO:HD2	1.65	0.61
1:B:65:ASP:CG	4:B:360:HOH:O	2.45	0.59
1:B:155:VAL:HG21	1:B:158:GLU:HB3	1.84	0.59
1:A:236:GLN:HG3	1:A:242:ALA:HB3	1.85	0.59
1:A:15:VAL:HG13	1:A:40:VAL:HG11	1.84	0.59
1:B:115:GLU:HA	1:B:140:GLY:O	2.03	0.58
1:A:108:ARG:HG3	1:A:112:ARG:NH2	2.19	0.58
1:A:254:GLU:HA	1:A:257:LYS:HD2	1.86	0.58
2:B:317:F3S:S2	3:B:998:H6P:H21A	2.43	0.57
1:B:291:ILE:HD11	4:B:368:HOH:O	2.04	0.57
1:A:-4:HIS:CB	1:A:289:GLU:OE2	2.52	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:245:ILE:CD1	1:B:251:ILE:HD11	2.34	0.56
1:B:152:LYS:O	1:B:152:LYS:HG3	2.04	0.56
1:B:209:ALA:O	1:B:212:GLU:HG2	2.06	0.56
1:A:117:ILE:HG21	1:A:181:LEU:HD21	1.88	0.55
1:A:32:PRO:HB3	4:A:364:HOH:O	2.08	0.54
1:A:212:GLU:HB2	4:A:367:HOH:O	2.07	0.54
1:B:281:ARG:HD3	4:B:347:HOH:O	2.07	0.54
1:B:164:MET:SD	1:B:164:MET:N	2.82	0.53
1:B:17:ARG:HD2	1:B:302:PHE:HB3	1.91	0.53
1:A:71:PHE:O	1:A:94:ALA:HB3	2.09	0.53
1:A:209:ALA:O	1:A:212:GLU:HG2	2.08	0.53
1:A:42:ASN:ND2	1:A:45:VAL:H	2.05	0.52
1:B:92:PHE:HE1	1:B:307:GLU:HG3	1.74	0.52
1:A:52:ARG:CD	4:A:355:HOH:O	2.38	0.52
1:B:218:LEU:HD13	1:B:251:ILE:HD11	1.91	0.51
1:B:-3:HIS:HB2	4:B:344:HOH:O	2.10	0.51
1:A:223:LYS:NZ	1:A:244:LEU:O	2.44	0.50
1:B:78:GLN:OE1	1:B:133:GLN:NE2	2.44	0.50
1:A:223:LYS:HD3	4:A:321:HOH:O	2.12	0.50
1:B:-5:HIS:HD2	4:B:344:HOH:O	1.94	0.50
1:B:203:ARG:HB3	1:B:231:LEU:HD11	1.94	0.49
1:B:218:LEU:HD13	1:B:251:ILE:CD1	2.42	0.49
1:A:36:ARG:HB2	1:A:69:LEU:HD11	1.95	0.49
1:B:42:ASN:ND2	1:B:44:TYR:HB3	2.28	0.49
1:A:8:PRO:HG2	1:A:202:ASN:HB3	1.95	0.48
1:B:115:GLU:OE1	1:B:142:TYR:OH	2.26	0.47
1:B:42:ASN:HD21	1:B:44:TYR:HB3	1.79	0.47
1:B:118:LEU:HB2	1:B:141:MET:HE1	1.97	0.47
1:A:306:LYS:HG3	1:A:307:GLU:N	2.28	0.47
1:B:182:ARG:HH21	1:B:182:ARG:HG2	1.80	0.46
1:B:100:THR:O	1:B:104:MET:HG2	2.16	0.46
1:A:217:VAL:O	1:A:242:ALA:HA	2.15	0.46
1:A:121:HIS:O	1:A:127:VAL:HG21	2.16	0.46
1:A:100:THR:O	1:A:104:MET:HG3	2.16	0.46
1:A:236:GLN:NE2	4:A:402:HOH:O	2.49	0.45
1:A:124:HIS:HA	1:A:125:PRO:HD2	1.66	0.45
1:B:15:VAL:HG13	1:B:40:VAL:HG11	1.98	0.45
1:B:163:PHE:CE2	1:B:190:GLY:HA3	2.52	0.45
1:A:157:ASN:OD1	1:A:159:GLU:HB2	2.16	0.45
1:B:178:ILE:HD13	1:B:193:LYS:O	2.17	0.45
1:B:15:VAL:O	1:B:19:ILE:HD12	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:17:ARG:HD2	1:B:302:PHE:CB	2.47	0.45
1:B:156:LYS:N	4:B:380:HOH:O	2.50	0.44
1:B:254:GLU:HA	1:B:257:LYS:CE	2.46	0.44
1:A:106:VAL:HA	1:A:164:MET:HE2	2.00	0.44
1:B:217:VAL:O	1:B:242:ALA:HA	2.16	0.44
1:A:1:MET:O	1:A:287:GLY:HA2	2.17	0.44
1:A:50:ARG:HH11	1:A:50:ARG:HD2	1.58	0.43
1:A:63:VAL:HA	1:A:64:PRO:HD3	1.75	0.43
1:A:184:ARG:HG2	1:A:185:PHE:CE1	2.54	0.43
1:A:207:VAL:HG12	1:A:234:LEU:HD13	2.01	0.42
1:A:129:GLY:HA3	4:A:347:HOH:O	2.19	0.42
1:A:97:PRO:HG2	1:A:302:PHE:CG	2.54	0.42
1:B:139:GLY:N	1:B:140:GLY:HA2	2.35	0.42
1:B:161:LEU:C	1:B:188:ILE:HG13	2.45	0.42
1:B:163:PHE:CE1	1:B:191:PRO:HD2	2.54	0.42
1:B:249:LYS:HB3	1:B:249:LYS:HE3	1.80	0.42
1:A:125:PRO:CB	4:A:417:HOH:O	2.55	0.42
1:A:42:ASN:HD21	1:A:44:TYR:HB3	1.85	0.42
1:A:308:LEU:HD23	1:A:308:LEU:HA	1.79	0.42
1:B:146:SER:HB2	1:B:147:PRO:CD	2.46	0.42
1:A:198:TYR:CG	1:A:199:ALA:N	2.88	0.41
1:A:162:SER:HA	1:A:189:VAL:O	2.20	0.41
1:B:6:ALA:HB2	1:B:210:LEU:HD22	2.02	0.41
1:B:272:ASP:O	1:B:276:GLN:HG3	2.20	0.41
1:B:153:LEU:HD23	4:B:349:HOH:O	2.20	0.41
1:A:52:ARG:NH2	4:A:355:HOH:O	2.54	0.41
1:B:115:GLU:HB3	1:B:142:TYR:HE1	1.84	0.41
1:A:136:ASN:HD22	1:A:137:PRO:CD	2.33	0.41
1:B:115:GLU:HB3	1:B:142:TYR:CE1	2.56	0.41
1:B:146:SER:HB2	4:B:359:HOH:O	2.21	0.41
1:B:270:ALA:HA	1:B:271:PRO:HD2	1.95	0.41
1:A:94:ALA:O	1:A:95:THR:C	2.64	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:86:SER:O	1:A:258:GLU:OE1[1_655]	2.11	0.09

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	312/328 (95%)	302 (97%)	9 (3%)	1 (0%)	36	29
1	B	298/328 (91%)	288 (97%)	8 (3%)	2 (1%)	18	10
All	All	610/656 (93%)	590 (97%)	17 (3%)	3 (0%)	24	16

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	132	GLY
1	B	152	LYS
1	A	125	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	261/274 (95%)	244 (94%)	17 (6%)	15	8
1	B	252/274 (92%)	230 (91%)	22 (9%)	9	4
All	All	513/548 (94%)	474 (92%)	39 (8%)	12	5

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-4	HIS
1	A	-1	SER
1	A	32	PRO

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Mol	Chain	Res	Type
1	A	52	ARG
1	A	60	ILE
1	A	98	LEU
1	A	138	GLU
1	A	158	GLU
1	A	187	LYS
1	A	189	VAL
1	A	207	VAL
1	A	230	ARG
1	A	231	LEU
1	A	257	LYS
1	A	273	ILE
1	A	281	ARG
1	A	307	GLU
1	B	-1	SER
1	B	19	ILE
1	B	70	ILE
1	B	98	LEU
1	B	116	SER
1	B	141	MET
1	B	160	LYS
1	B	165	THR
1	B	187	LYS
1	B	189	VAL
1	B	207	VAL
1	B	223	LYS
1	B	230	ARG
1	B	231	LEU
1	B	249	LYS
1	B	253	GLU
1	B	254	GLU
1	B	269	SER
1	B	281	ARG
1	B	284	GLN
1	B	306	LYS
1	B	309	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	-4	HIS
1	A	2	GLN

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Mol	Chain	Res	Type
1	A	42	ASN
1	A	133	GLN
1	A	136	ASN
1	A	213	GLN
1	A	229	ASN
1	A	236	GLN
1	A	252	GLN
1	A	277	ASN
1	B	-5	HIS
1	B	-3	HIS
1	B	42	ASN
1	B	78	GLN
1	B	133	GLN
1	B	277	ASN

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](#)

4 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	F3S	A	317	1	0,9,9	-	-	-		
3	H6P	A	998	-	13,14,14	1.96	3 (23%)	15,20,20	1.79	4 (26%)
3	H6P	B	998	-	13,14,14	1.16	1 (7%)	15,20,20	1.89	3 (20%)
2	F3S	B	317	1	0,9,9	-	-	-		

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
2	F3S	B	317	1	-	-	0/3/3/3
3	H6P	A	998	-	-	0/15/15/15	-
3	H6P	B	998	-	-	1/15/15/15	-
2	F3S	A	317	1	-	-	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	998	H6P	P13-O16	5.46	1.65	1.59
3	B	998	H6P	P13-O16	2.16	1.61	1.59
3	A	998	H6P	O29-C28	-2.14	1.40	1.44
3	A	998	H6P	P17-O18	-2.04	1.47	1.54

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	998	H6P	O20-P17-O18	4.53	124.78	107.80
3	A	998	H6P	O20-P17-O18	3.44	120.71	107.80
3	A	998	H6P	C28-C27-C30	3.00	131.12	126.20
3	A	998	H6P	O20-P17-O16	-2.99	94.62	104.64
3	A	998	H6P	O18-P17-O16	2.81	114.06	104.64
3	B	998	H6P	O33-C31-C30	-2.71	104.69	112.22
3	B	998	H6P	O20-P17-O16	-2.47	96.34	104.64

There are no chirality outliers.

All (1) torsion outliers are listed below:

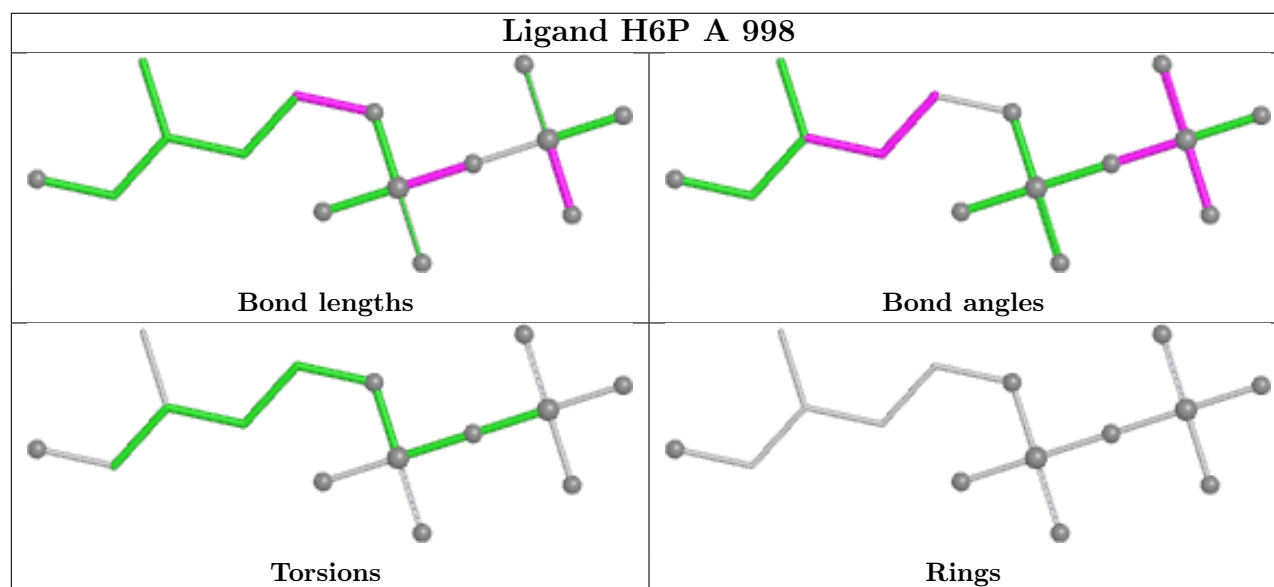
Mol	Chain	Res	Type	Atoms
3	B	998	H6P	C27-C28-O29-P13

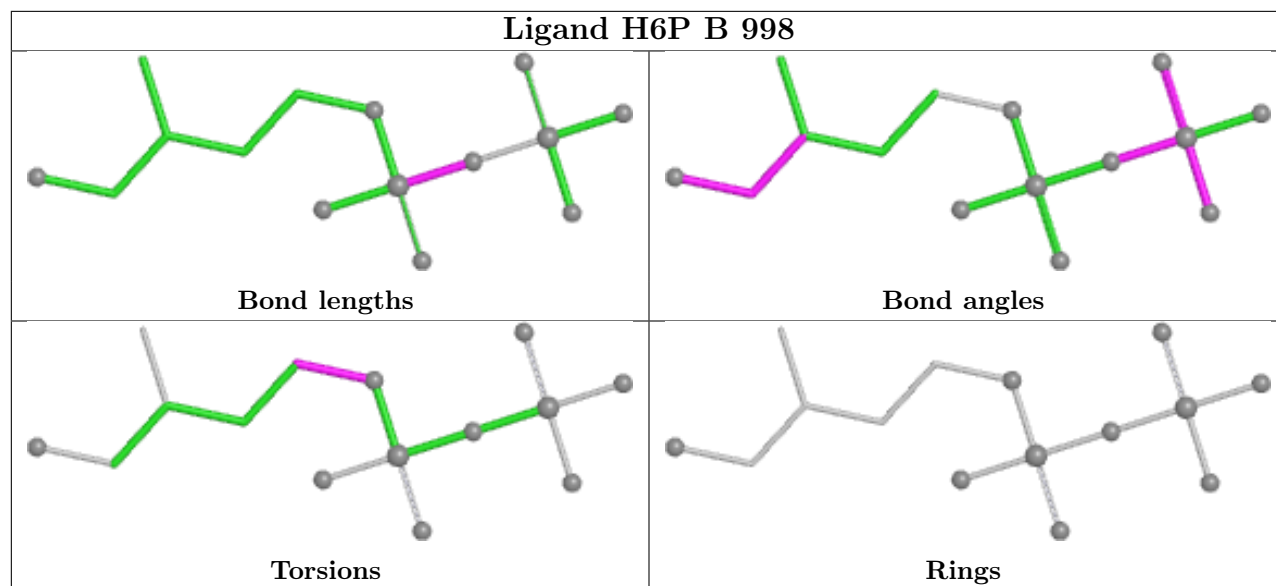
There are no ring outliers.

2 monomers are involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	998	H6P	1	0
2	B	317	F3S	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	314/328 (95%)	-0.41	1 (0%) 90 91	2, 7, 25, 49	0
1	B	304/328 (92%)	-0.24	9 (2%) 52 56	2, 8, 29, 35	0
All	All	618/656 (94%)	-0.33	10 (1%) 70 74	2, 8, 27, 49	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	-5	HIS	4.2
1	B	139	GLY	3.5
1	B	157	ASN	2.6
1	B	140	GLY	2.6
1	B	155	VAL	2.6
1	B	-3	HIS	2.5
1	B	131	MET	2.4
1	B	152	LYS	2.3
1	B	309	ARG	2.2
1	B	113	GLY	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

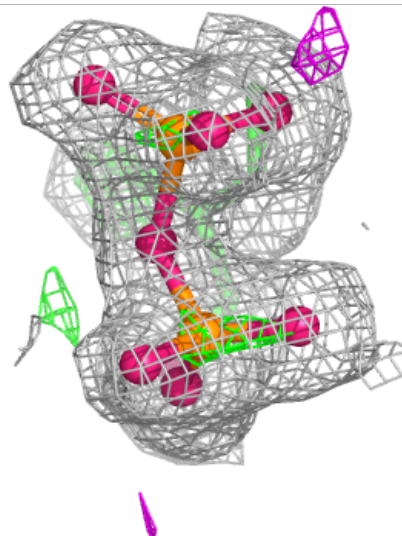
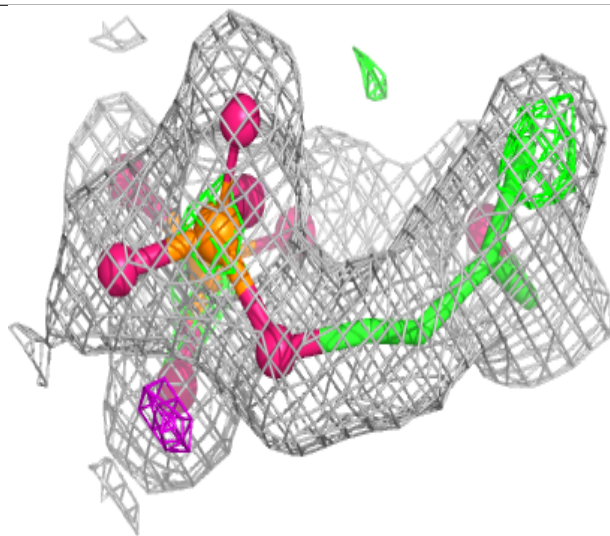
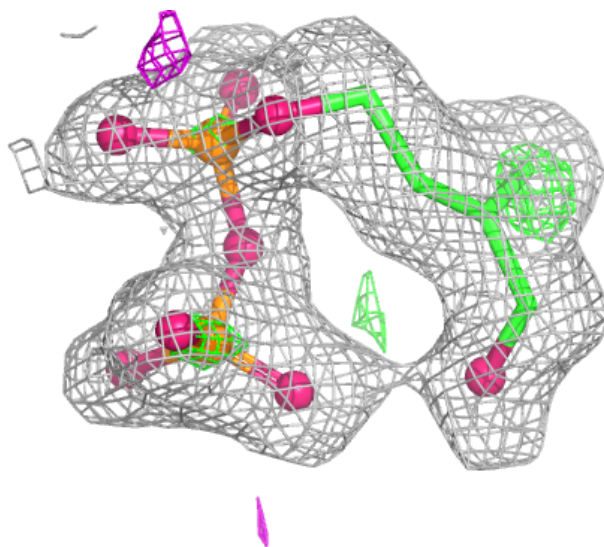
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	H6P	A	998	15/15	0.98	0.10	10,13,20,27	0
2	F3S	B	317	7/7	0.99	0.04	2,2,3,3	0
2	F3S	A	317	7/7	0.99	0.04	2,2,2,2	0
3	H6P	B	998	15/15	0.99	0.07	10,13,16,20	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

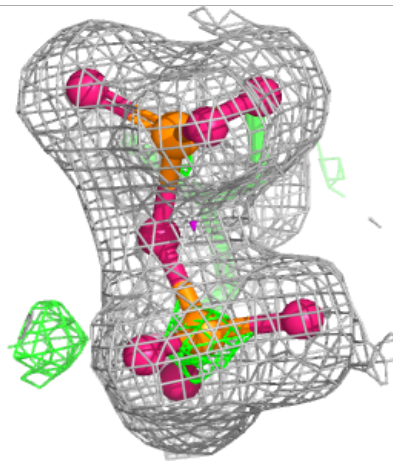
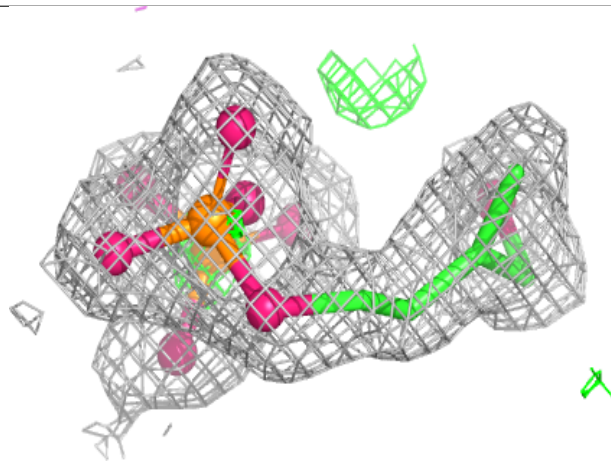
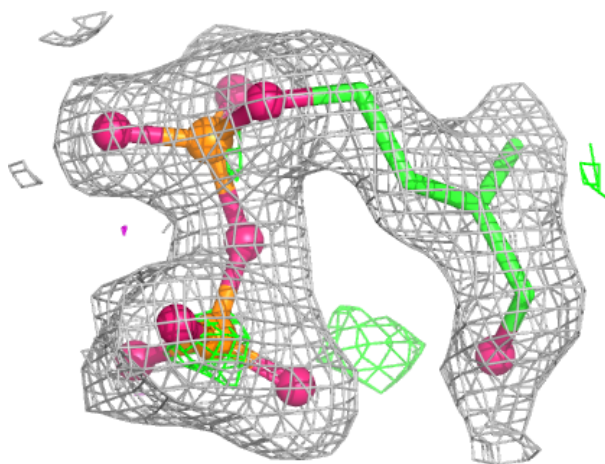
Electron density around H6P A 998:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around H6P B 998:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.