

Table VII Ligand abbreviations

Guidelines for the construction and use of ligand abbreviations are given in Section IR-4.4.4 and their use in the formulae of coordination complexes is described in Section IR-9.2.3.4. Abbreviations are listed in alphabetical order but those beginning with a numeral are listed by the first letter of the abbreviation (*e.g.* 2,3,2-tet appears under the letter 't').

Structural formulae of selected ligands are shown in Table VIII (numbered according to the present Table).

| <i>Number and abbreviation</i> | <i>Systematic name</i> | <i>Other name (from which abbreviation derived)</i> |
|--------------------------------|--|---|
| 1. 4-abu | 4-aminobutanoato | |
| 2. Ac | acetyl | |
| 3. acac | 2,4-dioxopentan-3-ido | acetylacetato |
| 4. acacen | 2,2'-[ethane-1,2-diylbis(azanylylidene)]bis(4-oxopentan-3-ido) | bis(acetylacetato)ethylenediamine |
| 5. ade | 1 <i>H</i> -purin-6-amine | adenine |
| 6. ado | 9-β-D-ribofuranosyl-9 <i>H</i> -purin-6-amine | adenosine |
| 7. adp | adenosine 5'-diphosphato(3-) | |
| 8. aet | 2-aminoethanethiolato | |
| 9. ala | 2-aminopropanoato | alaninato |
| 10. ama | 2-aminopropanedioato | aminomalonato |
| 11. amp | adenosine 5'-phosphato(2-) | adenosine monophosphato |

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| 12. [9]aneN ₃ (also tacn) | 1,4,7-triazonane | 1,4,7-triazacyclononane |
| 13. [12]aneN ₄ (also cyclen) | 1,4,7,10-tetraazacyclododecane | |
| 14. [14]aneN ₄ (also cyclam) | 1,4,8,11-tetraazacyclotetradecane | |
| 15. [18]aneP ₄ O ₂ | 1,10-dioxa-4,7,13,16-tetraphosphacyclooctadecane | |
| 16. [9]aneS ₃ | 1,4,7-trithionane | 1,4,7-trithiacyclononane |
| 17. [12]aneS ₄ | 1,4,7,10-tetrathiacyclododecane | |
| 18. arg | 2-amino-5-carbamimidamidopentanoato | argininato |
| 19. asn | 2,4-diamino-4-oxobutanoato | asparaginato |
| 20. asp | 2-aminobutanedioato | aspartato |
| 21. atmp | [nitrilotris(methylene)]tris(phosphonato)(6-) | aminotris(methylenephosphonato) |
| 22. atp | adenosine 5'-triphosphato(4-) | |
| 23. 2,3-bdta | 2,2',2'',2'''-(butane-2,3-diyl)dinitrido)tetraacetato | |
| 24. benzo-15-crown-5 | 2,3,5,6,8,9,11,12-octahydro-1,4,7,10,13-benzopentaoxacyclo-pentadecine | |
| 25. big | bis(carbamimidoyl)azanido | biguanid-3-ido |
| 26. biim | 2,2'-bi(1 <i>H</i> -imidazole)-1,1'-diido | 2,2'-biimidazolato |
| 27. binap | 1,1'-binaphthalene-2,2'-diylbis(diphenylphosphane) | |
| 28. bn | butane-2,3-diamine | |
| 29. bpy | 2,2'-bipyridine | |

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| 30. 4,4'-bpy | 4,4'-bipyridine | |
| 31. Bu | butyl | |
| 32. bzac | 1,3-dioxo-1-phenylbutan-2-ido | benzoylacetato |
| 33. bzim | benzimidazol-2-ido | |
| 34. Bz | benzyl | |
| 35. bztz | 1,3-benzothiazole | |
| 36. cat | benzene-1,2-diolato | catecholato |
| 37. cbdca | cyclobutane-1,1-dicarboxylato | |
| 38. cdta | 2,2',2'',2'''-(cyclohexane-1,2-diyl)dinitrilo)tetraacetato | |
| 39. C ₅ H ₄ Me | methylcyclopentadienyl | |
| 40. chxn (also, dach) | cyclohexane-1,2-diamine | |
| 41. cit | 2-hydroxypropane-1,2,3-tricarboxylato | citrato |
| 42. C ₅ Me ₅ ¹ | pentamethylcyclopentadienyl | |
| 43. cod | cycloocta-1,5-diene | |
| 44. cot | cycloocta-1,3,5,7-tetraene | |
| 45. Cp | cyclopentadienyl | |
| 46. cptn | cyclopentane-1,2-diamine | |
| 47. 18-crown-6 | 1,4,7,10,13,16-hexaoxacyclooctadecane | |

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| 48. crypt-211 | 4,7,13,18-tetraoxa-1,10-diazabicyclo[8.5.5]icosane | cryptand 211 |
| 49. crypt-222 | 4,7,13,16,21,24-hexaoxa-1,10-diazabicyclo[8.8.8]hexacosane | cryptand 222 |
| 50. Cy | cyclohexyl | |
| cyclam (see [14]aneN ₄ , No. 14) | | |
| cyclen (see [12]aneN ₄ , No. 13) | | |
| 51. cys | 2-amino-3-sulfanylpropanoato | cysteinato |
| 52. cyt | 4-aminopyrimidin-2(1 <i>H</i>)-one | cytosine |
| 53. dabco | 1,4-diazabicyclo[2.2.2]octane | |
| dach (see chxn, No. 40) | | |
| 54. dbm | 1,3-dioxo-1,3-diphenylpropan-2-ido | dibenzoylmethanato |
| 55. dea | 2,2'-azanediylbis(ethanolato) | diethanolaminato |
| 56. depe | ethane-1,2-diylbis(diethylphosphane) | 1,2-bis(diethylphosphino)ethane |
| 57. diars | benzene-1,2-diylbis(dimethylarsane) | |
| 58. dien | <i>N</i> -(2-aminoethyl)ethane-1,2-diamine | diethylenetriamine |
| 59. [14]1,3-dieneN ₄ | 1,4,8,11-tetraazacyclotetradeca-1,3-diene | |
| 60. diop | [2,2-dimethyl-1,3-dioxolane-4,5-diylbis(methylene)]bis(diphenylphosphane) | |
| 61. diox | 1,4-dioxane | |
| 62. dipamp | ethane-1,2-diylbis[(2-methoxyphenyl)phenylphosphane] | 'dimer of phenylanisylmethylphosphine' |

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| 63. dma | <i>N,N</i> -dimethylacetamide | dimethylacetamide |
| 64. dme | 1,2-dimethoxyethane | |
| 65. dmf | <i>N,N</i> -dimethylformamide | |
| 66. dmg | butane-2,3-diylidenebis(azanolate) | dimethylglyoximato |
| 67. dmpe | ethane-1,2-diylbis(dimethylphosphane) | 1,2-bis(dimethylphosphino)ethane |
| 68. dppm | methylenebis(dimethylphosphane) | bis(dimethylphosphino)methane |
| 69. dmso | (methanesulfinyl)methane | dimethyl sulfoxide |
| 70. dpm | 2,2,6,6-tetramethyl-3,5-dioxoheptan-4-ido | dipivaloylmethanato |
| 71. dppe | ethane-1,2-diylbis(diphenylphosphane) | 1,2-bis(diphenylphosphino)ethane |
| 72. dppf | 1,1'-bis(diphenylphosphanyl)ferrocene | |
| 73. dppm | methylenebis(diphenylphosphane) | bis(diphenylphosphino)methane |
| 74. dppp | propane-1,3-diylbis(diphenylphosphane) | 1,3-bis(diphenylphosphino)propane |
| 75. dtmpa | [(phosphonatomethyl)azanediylbis(ethane-2,1-diyl)dinitrilobis(methylene)]tetrakis(phosphonato)(10-) | diethylenetriaminepentakis(methylenephosphonato) |
| 76. dtpa | 2,2',2'',2'''-[(carboxylatomethyl)azanediylbis(ethane-2,1-diyl)dinitriolo]tetraacetato | diethylenetriaminepentaacetato |
| 77. ea | 2-aminoethanolato | ethanolaminato |
| 78. edda | 2,2'-[ethane-1,2-diylbis(azanediyl)]diacetato | ethylenediaminediacetato |
| 79. edta | 2,2',2'',2'''-(ethane-1,2-diyl)dinitriolo]tetraacetato | ethylenediaminetetraacetato |
| 80. edtmpa | [ethane-1,2-diyldinitrioltetrakis(methylene)]tetrakis(phosphonato)(8-) | ethylenediaminetetrakis(methylenephosphonato) |

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| 81. egta | 2,2',2'',2'''-[ethane-1,2-diylbis(oxyethane-2,1-diyl)trilo)]tetraacetato | ethylene glycol-bis(2-aminoethyl)-N,N,N',N'-tetraacetic acid |
| 82. en | ethane-1,2-diamine | |
| 83. Et | ethyl | |
| 84. Et ₂ dtc | N,N-diethylcarbamodithioato | N,N-diethyldithiocarbamato |
| 85. fod | 6,6,7,7,8,8,8-heptafluoro-2,2-dimethyl-3,5-dioxooctan-4-ido | |
| 86. fta | 1,1,l-trifluoro-2,4-dioxopentan-3-ido | trifluoroacetylacetonato |
| 87. gln | 2,5-diamino-5-oxopentanoato | glutaminato |
| 88. glu | 2-aminopentanedioato | glutamato |
| 89. gly | aminoethanoato, or 2-aminoacetato | glycinato |
| 90. gua | 2-amino-9H-purin-6(1H)-one | guanine |
| 91. guo | 2-amino-9-β-D-ribofuranosyl-9H-purin-6(1H)-one | guanosine |
| 92. hdtmpa | [(hexane-1,6-diyl)dinitrilo)tetrakis(methylene)]tetrakis(phosphonato)(8-) | hexamethylenediaminetetrakis(methylenephosphonato) |
| 93. hedp | 1-hydroxyethane-1,1-bis(phosphonato)(4-) | 1-hydroxyethane-1,1-diphosphonato |
| 94. hfa | 1,1,1,5,5-hexafluoropentane-2,4-dioxopentan-3-ido | hexafluoroacetylacetonato |
| 95. his | 2-amino-3-(imidazol-4-yl)propanoato | histidinato |
| 96. hmpa | N,N,N',N'',N''-hexamethylphosphoric triamide | |
| 97. hmta | 1,3,5,7-tetraazatricyclo[3.3.1.1 ^{3,7}]decane | hexamethylenetetraamine |
| 98. ida | 2,2'-azanediyl diacetato | iminodiacetato |

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| 99. ile | 2-amino-3-methylpentanoato | isoleucinato |
| 100. im | imidazol-1-ido | |
| 101. isn | pyridine-4-carboxamide | isonicotinamide |
| 102. leu | 2-amino-4-methylpentanoato | leucinato |
| 103. lut | 2,6-dimethylpyridine | lutidine |
| 104. lys | 2,6-diaminohexanoato | lysinato |
| 105. mal | 2-hydroxybutanedioato | malato |
| 106. male | (Z)-butenedioato | maleato |
| 107. malo | propanedioato | malonato |
| 108. Me | methyl | |
| 109. 2-Mepy | 2-methylpyridine | |
| 110. met | 2-amino-4-(methylsulfanyl)butanoato | methioninato |
| 111. mnt | 1,2-dicyanoethene-1,2-dithiolato | maleonitriledithiolato |
| 112. napy | 1,8-naphthyridine | |
| 113. nbd | bicyclo[2.2.1]hepta-2,5-diene | norbornadiene |
| 114. nia | pyridine-3-carboxamide | nicotinamide |
| 115. nmp | <i>N</i> -methylpyrrolidine | |
| 116. nta | 2,2',2"-nitrilotriacetato | |

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| 117. oep | 2,3,7,8,12,13,17,18-octaethylporphyrin-21,23-ido | |
| 118. ox | ethanedioato | oxalato |
| 119. pc | phthalocyanine-29,31-diido | |
| 120. 1,2-pdta | 2,2',2'',2'''-(propane-1,2-diyldinitrilo)tetraacetato | 1,2-propylenediaminetetraacetato |
| 121. 1,3-pdta | 2,2',2'',2'''-(propane-1,3-diyldinitrilo)tetraacetato | 1,3-propylenediaminetetraacetato |
| 122. Ph | phenyl | |
| 123. phe | 2-amino-3-phenylpropanoato | phenylalaninato |
| 124. phen | 1,10-phenanthroline | |
| 125. pip | piperidine | |
| 126. pn | propane-1,2-diamine | |
| 127. pmdien | 2,2'-(methylazanediyl)bis(<i>N,N</i> -dimethylethanamine) | <i>N,N,N',N'',N''-pentamethyldiethylenetriamine</i> |
| 128. ppIX | 2,18-bis(carboxyethyl)-3,7,12,17-tetramethyl-8,13-divinylporphyrin-21,23-diido | protoporphyrinato IX |
| 129. pro | pyrrolidine-2-carboxylato | prolinato |
| 130. ptn | pentane-2,4-diamine | |
| 131. py | pyridine | |
| 132. pyz | pyrazine | |
| 133. pz | 1 <i>H</i> -pyrazol-1-ido | |
| 134. qdt | quinoxaline-2,3-dithiolato | |

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| 135. quin | quinolin-8-olato | |
| 136. sal | 2-hydroxybenzoato | salicylato |
| 137. salan | 2-[(phenylimino)methyl]phenolato | salicylideneanilinato |
| 138. saldien | 2,2'-[azanediylbis(ethane-2,1-diylazanylylidene)methanylylidene)]diphenolato | bis(salicylidene)diethylenetriaminato |
| 139. salen | 2,2'-[ethane-1,2-diylbis(azanylylidene)methanylylidene)]diphenolato | bis(salicylidene)ethylenediaminato |
| 140. salgly | <i>N</i> -(2-oxidobenzylidene)glycinato | salicylideneglycinato |
| 141. salpn | 2,2'-[propane-1,2-diylbis(azanylylidene)methanylylidene)]diphenolato | bis(salicylidene)propylenediaminato |
| 142. saltn | 2,2'-[propane-1,3-diylbis(azanylylidene)methanylylidene)]diphenolato | bis(salicylidene)trimethylenediaminato |
| 143. sdta | 2,2',2'',2'''-(1,2-diphenylethane-1,2-diyl)dinitrilo)tetraacetato | stilbenediaminetetraacetato |
| 144. sep ² | 1,3,6,8,10,13,16,19-octaazabicyclo[6.6.6]icosane | |
| 145. ser | 2-amino-3-hydroxypropanoato | serinato |
| 146. stien ³ | 1,2-diphenylethane-1,2-diamine | |
| tacn (see [9]aneN ₃ , No. 12) | | |
| 147. tap | propane-1,2,3-triamine | 1,2,3-triaminopropane |
| 148. tart | 2,3-dihydroxybutanedioato | tartrato |
| 149. tcne | ethenetetracarbonitrile | tetracyanoethylene |
| 150. tcnq | 2,2'-(cyclohexa-2,5-diene-1,4-diylidene)bis(propane-1,3-dinitrile) | tetracyanoquinodimethane |
| 151. tdt | 4-methylbenzene-1,2-dithiolato | |

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| 152. tea | 2,2',2"-nitrilotris(ethanolato) | triethanolaminato |
| 153. terpy | 2,2':6',2"-terpyridine | terpyridine |
| 154. 2,3,2-tet | <i>N,N'</i> -bis(2-aminoethyl)propane-1,3-diamine | 1,4,8,11-tetraazaundecane |
| 155. 3,3,3-tet | <i>N,N'</i> -bis(3-aminopropyl)propane-1,3-diamine | 1,5,9,13-tetraazatridecane |
| 156. tetren | <i>N,N'</i> -(azanediylidethane-2,1-diyl)bis(ethane-1,2-diamine) | tetraethylenepentamine |
| 157. tfa | trifluoroacetato | |
| 158. thf | oxolane | tetrahydrofuran |
| 159. thiox | 1,4-oxathiane | thioxane |
| 160. thr | 2-amino-3-hydroxybutanoato | threoninato |
| 161. tht | thiolane | tetrahydrothiophene |
| 162. thy | 5-methylpyrimidine-2,4(<i>1H,3H</i>)-dione | thymine |
| 163. tmen | <i>N,N,N',N'</i> -tetramethylethane-1,2-diamine | |
| 164. tmp | 5,10,15,20-tetrakis(2,4,6-trimethylphenyl)porphyrin-21,23-diido | 5,10,15,20-tetramesitylporphyrin-21,23-diido |
| 165. tn | propane-1,3-diamine | trimethylenediamine |
| 166. Tol (<i>o</i> -, <i>m</i> - or <i>p</i> -) | 2-, 3- or 4-methylphenyl | tolyl (<i>o</i> -, <i>m</i> - or <i>p</i> -) |
| 167. Tp | hydridotris(pyrazolido- <i>N</i>)borato(1-), or tris(1 <i>H</i> -pyrazol-1-yl)boranuido | hydrotris(pyrazolyl)borato |
| 168. Tp ⁴ | hydridotris(3,5-dimethylpyrazolido- <i>N</i>)borato(1-) | hydrotris(3,5-dimethylpyrazolyl)borato |
| 169. tpp | 5,10,15,20-tetraphenylporphyrin-21,23-diido | |

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| 170. tren | <i>N,N</i> -bis(2-aminoethyl)ethane-1,2-diamine | tris(2-aminoethyl)amine |
| 171. trien | <i>N,N'</i> -bis(2-aminoethyl)ethane-1,2-diamine | triethylenetetramine |
| 172. triphos ⁵ | [(phenylphosphanediyI)bis(ethane-2,1-diyI)]bis(diphenylphosphane) | |
| 173. tris | 2-amino-2-(hydroxymethyl)propane-1,3-diol | aminotris(hydroxymethyl)methane |
| 174. trp | 2-amino-3-(1 <i>H</i> -indol-3-yl)propanoato | tryptophanato |
| 175. tsalen | 2,2'-[ethane-1,2-diylbis(azanylylidene)methanylylidene]dibenzene-thiolato | bis(thiosalicylidene)ethylenediaminato |
| 176. ttfa | 4,4,4-trifluoro-1,3-dioxo-1-(2-thienyl)butan-2-ido | thenoyltrifluoroacetonato |
| 177. ttha | [2,2',2'',2'''-(ethane-1,2-diylbis{[(carboxylatomethyl)azanediyl]ethane-2,1-diyI}nitrilo)]tetraacetato | triethylenetetraminehexaacetato |
| 178. ttp | 5,10,15,20-tetrakis(4-methylphenyl)porphyrin-21,23-diido | 5,10,15,20-tetra- <i>p</i> -tolylporphyrin-21,23-diido |
| 179. tu | thiourea | |
| 180. tyr | 2-amino-3-(4-hydroxyphenyl)propanoato | tyrosinato |
| 181. tz | 1,3-thiazole | thiazole |
| 182. ura | pyrimidine-2,4(1 <i>H</i> ,3 <i>H</i>)-dione | uracil |
| 183. val | 2-amino-3-methylbutanoato | valinato |

Footnotes

- The use of the abbreviation Cp^{*} for pentamethylcyclopentadienyl is discouraged. It can lead to confusion because the asterisk, *, is also used to represent an excited state, an optically active substance, a radioactive substance, etc.

2. The abbreviation derives from the non-systematic name sepulchrate which incorrectly implies that the ligand is anionic.
3. The abbreviation derives from the non-systematic name stilbenediamine which incorrectly implies the presence of a C=C double bond in the ligand.
4. The use of Tp' is preferred to Tp* for the reasons given in Note 1. A general procedure for abbreviating substituted hydridotris(pyrazolido-*N*)borate ligands has been proposed {see S. Trofimenko, *Chem. Rev.*, **93**, 943 (1993)}. For example, Tp' becomes Tp^{Me₂}, the superscript denoting the methyl groups at the 3- and 5-positions of the pyrazolyl rings.
5. The abbreviation triphos should not be used for the ligand PhP(CH₂PPh₂)₃.